



## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W9** Lab Sample ID: **21G0031-01**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 13:30

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 14:50

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| <b>Acetone</b>              | <b>92</b>     |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W9** Lab Sample ID: **21G0031-01**

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### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 14:50

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W9** Lab Sample ID: **21G0031-01**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 13:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:11

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W9** Lab Sample ID: **21G0031-01**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 13:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:11

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W9** Lab Sample ID: **21G0031-01**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 13:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:11

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W9** Lab Sample ID: **21G0031-01**

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### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:11

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W9** Lab Sample ID: **21G0031-01**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 13:30

### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:11

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 10:37

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | 0.19          |                  | 0.06                   |

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:18

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Aluminum       | 5080          |                  | 100                    |
| Antimony       | < 10.0        |                  | 10.0                   |
| Arsenic        | < 10.0        |                  | 10.0                   |
| Barium         | 129           |                  | 10.0                   |
| Beryllium      | < 1.00        |                  | 1.00                   |
| Boron          | 293           | B1               | 25.0                   |
| Cadmium        | 4.93          |                  | 3.00                   |

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Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W9** Lab Sample ID: **21G0031-01**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 13:30

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:25

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Calcium        | 220000        |                  | 3000                   |
| Chromium       | 11.9          |                  | 5.00                   |
| Cobalt         | 20.0          |                  | 10.0                   |
| Copper         | 62.0          |                  | 10.0                   |
| Hardness       | 964000        |                  | 1980                   |
| Iron           | 15000         |                  | 150                    |
| Lead           | 8.20          |                  | 5.00                   |
| Magnesium      | 101000        |                  | 300                    |
| Manganese      | 415           |                  | 15.0                   |
| Nickel         | 16.6          |                  | 5.00                   |
| Potassium      | 4560          |                  | 1400                   |
| Selenium       | < 25.0        | B1               | 25.0                   |
| Silver         | 3.99          |                  | 3.00                   |
| Sodium         | 169000        |                  | 300                    |
| Strontium      | 328           |                  | 5.00                   |
| Thallium       | < 10.0        |                  | 10.0                   |
| Vanadium       | 14.4          |                  | 5.00                   |
| Zinc           | 56.9          |                  | 25.0                   |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W13** Lab Sample ID: **21G0031-02**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 14:30

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 15:10

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W13** Lab Sample ID: **21G0031-02**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 14:30

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 15:10

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W13** Lab Sample ID: **21G0031-02**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 14:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:45

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W13** Lab Sample ID: **21G0031-02**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 14:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:45

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W13** Lab Sample ID: **21G0031-02**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 14:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:45

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W13** Lab Sample ID: **21G0031-02**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 14:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:45

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W13** Lab Sample ID: **21G0031-02**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 14:30

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 18:45

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 10:39

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:21

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>2450</b>   |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>52.9</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>60.3</b>   | B1               | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W13** Lab Sample ID: **21G0031-02**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 14:30

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:21

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>140000</b> |                  | 300                    |
| <b>Chromium</b>  | <b>7.19</b>   |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| <b>Copper</b>    | <b>10.1</b>   |                  | 10.0                   |
| <b>Hardness</b>  | <b>578000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>4610</b>   |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>55200</b>  |                  | 300                    |
| <b>Manganese</b> | <b>158</b>    |                  | 15.0                   |
| <b>Nickel</b>    | <b>5.33</b>   |                  | 5.00                   |
| Potassium        | < 1400        |                  | 1400                   |
| Selenium         | < 25.0        | B1               | 25.0                   |
| <b>Silver</b>    | <b>3.22</b>   |                  | 3.00                   |
| <b>Sodium</b>    | <b>10600</b>  |                  | 300                    |
| <b>Strontium</b> | <b>79.5</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| <b>Vanadium</b>  | <b>5.77</b>   |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W54** Lab Sample ID: **21G0031-03**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:15

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 15:31

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W54** Lab Sample ID: **21G0031-03**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:15

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 15:31

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| <b>Tetrachloroethene</b>  | <b>7.6</b>    |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W54** Lab Sample ID: **21G0031-03**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:15

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:20

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W54** Lab Sample ID: **21G0031-03**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:15

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:20

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W54** Lab Sample ID: **21G0031-03**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:15

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:20

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W54** Lab Sample ID: **21G0031-03**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:15

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:20

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W54** Lab Sample ID: **21G0031-03**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:15

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:20

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 10:42

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:23

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>18000</b>  |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>166</b>    |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| Boron           | < 25.0        | B1               | 25.0                   |
| <b>Cadmium</b>  | <b>8.58</b>   |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W54** Lab Sample ID: **21G0031-03**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:15

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:27

| <u>Analyte</u>   | <u>Result</u>  | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|----------------|------------------|------------------------|
| <b>Calcium</b>   | <b>399000</b>  |                  | 3000                   |
| <b>Chromium</b>  | <b>129</b>     |                  | 5.00                   |
| <b>Cobalt</b>    | <b>20.5</b>    |                  | 10.0                   |
| <b>Copper</b>    | <b>60.0</b>    |                  | 10.0                   |
| <b>Hardness</b>  | <b>1730000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>34000</b>   |                  | 150                    |
| <b>Lead</b>      | <b>16.3</b>    |                  | 5.00                   |
| <b>Magnesium</b> | <b>179000</b>  |                  | 300                    |
| <b>Manganese</b> | <b>875</b>     |                  | 15.0                   |
| <b>Nickel</b>    | <b>72.4</b>    |                  | 5.00                   |
| <b>Potassium</b> | <b>5640</b>    |                  | 1400                   |
| Selenium         | < 25.0         | B1               | 25.0                   |
| <b>Silver</b>    | <b>6.10</b>    |                  | 3.00                   |
| <b>Sodium</b>    | <b>46100</b>   |                  | 300                    |
| <b>Strontium</b> | <b>219</b>     |                  | 5.00                   |
| Thallium         | < 10.0         |                  | 10.0                   |
| <b>Vanadium</b>  | <b>38.2</b>    |                  | 5.00                   |
| <b>Zinc</b>      | <b>107</b>     |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W3R** Lab Sample ID: **21G0031-04**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:55

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 15:52

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W3R** Lab Sample ID: **21G0031-04**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:55

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 15:52

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W3R** Lab Sample ID: **21G0031-04**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:54

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W3R** Lab Sample ID: **21G0031-04**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:54

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W3R** Lab Sample ID: **21G0031-04**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:54

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W3R** Lab Sample ID: **21G0031-04**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:54

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W3R** Lab Sample ID: **21G0031-04**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:55

### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 19:54

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 10:44

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:26

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>429</b>    |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>45.2</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>38.7</b>   |                  | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W3R** Lab Sample ID: **21G0031-04**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 15:55

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:26

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>110000</b> |                  | 300                    |
| <b>Chromium</b>  | <b>8.96</b>   |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| <b>Copper</b>    | <b>18.7</b>   |                  | 10.0                   |
| <b>Hardness</b>  | <b>493000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>5260</b>   |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>52900</b>  |                  | 300                    |
| <b>Manganese</b> | <b>39.0</b>   |                  | 15.0                   |
| Nickel           | < 5.00        |                  | 5.00                   |
| Potassium        | < 1400        |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>8750</b>   |                  | 300                    |
| <b>Strontium</b> | <b>73.3</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

### Volatile Organic Compounds by GC/MS

Method: 524.3 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 11:27

| <u>Analyte</u>           | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|--------------------------|---------------|------------------|------------------------|
| 1,1,1-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1,2-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1-Dichloroethene       | < 0.50        |                  | 0.50                   |
| 1,2-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| 1,2-Dichloroethane       | < 0.50        |                  | 0.50                   |
| 1,2-Dichloropropane      | < 0.50        |                  | 0.50                   |
| 1,4-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| Benzene                  | < 0.50        |                  | 0.50                   |
| Carbon tetrachloride     | < 0.50        |                  | 0.50                   |
| Chlorobenzene            | < 0.50        |                  | 0.50                   |
| cis-1,2-Dichloroethene   | < 0.50        |                  | 0.50                   |
| Ethylbenzene             | < 0.50        |                  | 0.50                   |
| Methyl tert-butyl ether  | < 0.50        |                  | 0.50                   |
| Methylene chloride       | < 0.50        |                  | 0.50                   |
| Styrene                  | < 0.50        |                  | 0.50                   |
| Tetrachloroethene        | < 0.50        |                  | 0.50                   |
| Toluene                  | < 0.50        |                  | 0.50                   |
| trans-1,2-Dichloroethene | < 0.50        |                  | 0.50                   |
| Trichloroethene          | < 0.50        |                  | 0.50                   |
| Vinyl chloride           | < 0.50        | J5               | 0.50                   |
| Xylenes, total           | < 0.50        |                  | 0.50                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 16:12

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 16:12

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 20:28

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 20:28

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 20:28

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 20:28

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 20:28

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 10:46

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:29

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Aluminum       | < 100         |                  | 100                    |
| Antimony       | < 10.0        |                  | 10.0                   |
| Arsenic        | < 10.0        |                  | 10.0                   |
| <b>Barium</b>  | <b>21.1</b>   |                  | 10.0                   |
| Beryllium      | < 1.00        |                  | 1.00                   |
| <b>Boron</b>   | <b>30.7</b>   |                  | 25.0                   |
| Cadmium        | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W19B** Lab Sample ID: **21G0031-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:30

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:29

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>80700</b>  |                  | 300                    |
| Chromium         | < 5.00        |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>369000</b> |                  | 1980                   |
| Iron             | < 150         |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>40700</b>  |                  | 300                    |
| <b>Manganese</b> | <b>527</b>    |                  | 15.0                   |
| Nickel           | < 5.00        |                  | 5.00                   |
| <b>Potassium</b> | <b>2330</b>   |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>30500</b>  |                  | 300                    |
| <b>Strontium</b> | <b>97.8</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

### Volatile Organic Compounds by GC/MS

Method: 524.3 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 12:04

| <u>Analyte</u>           | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|--------------------------|---------------|------------------|------------------------|
| 1,1,1-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1,2-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1-Dichloroethene       | < 0.50        |                  | 0.50                   |
| 1,2-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| 1,2-Dichloroethane       | < 0.50        |                  | 0.50                   |
| 1,2-Dichloropropane      | < 0.50        |                  | 0.50                   |
| 1,4-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| Benzene                  | < 0.50        |                  | 0.50                   |
| Carbon tetrachloride     | < 0.50        |                  | 0.50                   |
| Chlorobenzene            | < 0.50        |                  | 0.50                   |
| cis-1,2-Dichloroethene   | < 0.50        |                  | 0.50                   |
| Ethylbenzene             | < 0.50        |                  | 0.50                   |
| Methyl tert-butyl ether  | < 0.50        |                  | 0.50                   |
| Methylene chloride       | < 0.50        |                  | 0.50                   |
| Styrene                  | < 0.50        |                  | 0.50                   |
| Tetrachloroethene        | < 0.50        |                  | 0.50                   |
| Toluene                  | < 0.50        |                  | 0.50                   |
| trans-1,2-Dichloroethene | < 0.50        |                  | 0.50                   |
| Trichloroethene          | < 0.50        |                  | 0.50                   |
| Vinyl chloride           | < 0.50        | J5               | 0.50                   |
| Xylenes, total           | < 0.50        |                  | 0.50                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 16:33

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 16:33

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:03

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:03

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:03

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:03

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:03

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 10:49

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:37

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Aluminum       | < 100         |                  | 100                    |
| Antimony       | < 10.0        |                  | 10.0                   |
| Arsenic        | < 10.0        |                  | 10.0                   |
| <b>Barium</b>  | <b>21.1</b>   |                  | 10.0                   |
| Beryllium      | < 1.00        |                  | 1.00                   |
| <b>Boron</b>   | <b>35.0</b>   |                  | 25.0                   |
| Cadmium        | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **219B** Lab Sample ID: **21G0031-06**

Matrix: Water Collected By: MW Date/Time Collected: 06/30/21 18:45

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:37

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>81000</b>  |                  | 300                    |
| Chromium         | < 5.00        |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>372000</b> |                  | 1980                   |
| Iron             | < 150         |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>41200</b>  |                  | 300                    |
| <b>Manganese</b> | <b>520</b>    |                  | 15.0                   |
| Nickel           | < 5.00        |                  | 5.00                   |
| <b>Potassium</b> | <b>2360</b>   |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>30800</b>  |                  | 300                    |
| <b>Strontium</b> | <b>97.6</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W12R** Lab Sample ID: **21G0031-07**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 13:45

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 16:54

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W12R** Lab Sample ID: **21G0031-07**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 13:45

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 16:54

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W12R** Lab Sample ID: **21G0031-07**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 13:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:37

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W12R** Lab Sample ID: **21G0031-07**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 13:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:37

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W12R** Lab Sample ID: **21G0031-07**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 13:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:37

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W12R** Lab Sample ID: **21G0031-07**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 13:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:37

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W12R** Lab Sample ID: **21G0031-07**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 13:45

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 21:37

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 10:55

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:40

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>166</b>    |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>20.3</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>131</b>    |                  | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W12R** Lab Sample ID: **21G0031-07**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 13:45

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:40

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>79300</b>  |                  | 300                    |
| Chromium         | < 5.00        |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>332000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>309</b>    |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>32500</b>  |                  | 300                    |
| Manganese        | < 15.0        |                  | 15.0                   |
| Nickel           | < 5.00        |                  | 5.00                   |
| Potassium        | < 1400        |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>2210</b>   |                  | 300                    |
| <b>Strontium</b> | <b>62.7</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W22** Lab Sample ID: **21G0031-08**

Matrix: Water Collected By: AG Date/Time Collected: 06/30/21 14:10

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 17:14

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W22** Lab Sample ID: **21G0031-08**

Matrix: Water Collected By: AG Date/Time Collected: 06/30/21 14:10

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 17:14

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W22** Lab Sample ID: **21G0031-08**

Matrix: Water Collected By: AG Date/Time Collected: 06/30/21 14:10

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:12

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W22** Lab Sample ID: **21G0031-08**

Matrix: Water Collected By: AG Date/Time Collected: 06/30/21 14:10

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:12

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W22** Lab Sample ID: **21G0031-08**

Matrix: Water Collected By: AG Date/Time Collected: 06/30/21 14:10

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:12

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W22** Lab Sample ID: **21G0031-08**

Matrix: Water Collected By: AG Date/Time Collected: 06/30/21 14:10

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:12

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W22** Lab Sample ID: **21G0031-08**

Matrix: Water Collected By: AG Date/Time Collected: 06/30/21 14:10

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:12

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 10:58

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:43

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>221</b>    |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>46.2</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>31.1</b>   | B1               | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W22** Lab Sample ID: **21G0031-08**

Matrix: Water Collected By: AG Date/Time Collected: 06/30/21 14:10

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:43

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>105000</b> |                  | 300                    |
| Chromium         | < 5.00        |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>487000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>559</b>    |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>54400</b>  |                  | 300                    |
| <b>Manganese</b> | <b>19.3</b>   |                  | 15.0                   |
| Nickel           | < 5.00        |                  | 5.00                   |
| Potassium        | < 1400        |                  | 1400                   |
| Selenium         | < 25.0        | B1               | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>13400</b>  |                  | 300                    |
| <b>Strontium</b> | <b>58.0</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W45** Lab Sample ID: **21G0031-09**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 14:50

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 17:35

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W45** Lab Sample ID: **21G0031-09**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 14:50

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 17:35

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W45** Lab Sample ID: **21G0031-09**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 14:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:46

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W45** Lab Sample ID: **21G0031-09**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 14:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:46

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W45** Lab Sample ID: **21G0031-09**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 14:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:46

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W45** Lab Sample ID: **21G0031-09**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 14:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:46

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W45** Lab Sample ID: **21G0031-09**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 14:50

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 22:46

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:00

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:46

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Aluminum       | 750           |                  | 100                    |
| Antimony       | < 10.0        |                  | 10.0                   |
| Arsenic        | < 10.0        |                  | 10.0                   |
| Barium         | 25.7          |                  | 10.0                   |
| Beryllium      | < 1.00        |                  | 1.00                   |
| Boron          | < 25.0        |                  | 25.0                   |
| Cadmium        | < 3.00        |                  | 3.00                   |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W45** Lab Sample ID: **21G0031-09**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 14:50

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:46

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>74100</b>  |                  | 300                    |
| Chromium         | < 5.00        |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>336000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>1080</b>   |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>36700</b>  |                  | 300                    |
| <b>Manganese</b> | <b>34.5</b>   |                  | 15.0                   |
| Nickel           | < 5.00        |                  | 5.00                   |
| <b>Potassium</b> | <b>1580</b>   |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>1660</b>   |                  | 300                    |
| <b>Strontium</b> | <b>38.3</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W20R** Lab Sample ID: **21G0031-10**

Matrix: Water Collected By: TAR Date/Time Collected: 06/30/21 13:05

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 17:55

| <u>Analyte</u>               | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane    | < 2.0         |                  | 2.0                    |
| <b>1,1,1-Trichloroethane</b> | <b>2.3</b>    |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane    | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane        | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane           | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene           | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene          | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane       | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane            | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane          | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane          | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane          | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)             | < 10          |                  | 10                     |
| 2-Hexanone (MBK)             | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK)  | < 10          |                  | 10                     |
| Acetone                      | < 10          |                  | 10                     |
| Benzene                      | < 2.0         |                  | 2.0                    |
| Bromobenzene                 | < 2.0         |                  | 2.0                    |
| Bromochloromethane           | < 2.0         |                  | 2.0                    |
| Bromodichloromethane         | < 2.0         |                  | 2.0                    |
| Bromoform                    | < 5.0         |                  | 5.0                    |
| Bromomethane                 | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W20R** Lab Sample ID: **21G0031-10**

Matrix: Water Collected By: TAR Date/Time Collected: 06/30/21 13:05

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 17:55

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W20R** Lab Sample ID: **21G0031-10**

Matrix: Water Collected By: TAR Date/Time Collected: 06/30/21 13:05

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:21

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W20R** Lab Sample ID: **21G0031-10**

Matrix: Water Collected By: TAR Date/Time Collected: 06/30/21 13:05

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:21

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W20R** Lab Sample ID: **21G0031-10**

Matrix: Water Collected By: TAR Date/Time Collected: 06/30/21 13:05

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:21

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W20R** Lab Sample ID: **21G0031-10**

Matrix: Water Collected By: TAR Date/Time Collected: 06/30/21 13:05

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:21

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W20R** Lab Sample ID: **21G0031-10**

Matrix: Water Collected By: TAR Date/Time Collected: 06/30/21 13:05

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:21

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:02

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:48

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>1710</b>   |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>49.5</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>32.4</b>   | B1               | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W20R** Lab Sample ID: **21G0031-10**

Matrix: Water Collected By: TAR Date/Time Collected: 06/30/21 13:05

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:48

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>131000</b> |                  | 300                    |
| <b>Chromium</b>  | <b>8.77</b>   |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| <b>Copper</b>    | <b>11.1</b>   |                  | 10.0                   |
| <b>Hardness</b>  | <b>550000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>4570</b>   |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>54300</b>  |                  | 300                    |
| <b>Manganese</b> | <b>490</b>    |                  | 15.0                   |
| <b>Nickel</b>    | <b>12.6</b>   |                  | 5.00                   |
| <b>Potassium</b> | <b>1400</b>   |                  | 1400                   |
| Selenium         | < 25.0        | B1               | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>8070</b>   |                  | 300                    |
| <b>Strontium</b> | <b>87.5</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| <b>Vanadium</b>  | <b>6.92</b>   |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W46** Lab Sample ID: **21G0031-11**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 16:40

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 18:16

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W46** Lab Sample ID: **21G0031-11**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 16:40

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 18:16

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W46** Lab Sample ID: **21G0031-11**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 16:40

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:56

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W46** Lab Sample ID: **21G0031-11**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 16:40

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:56

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W46** Lab Sample ID: **21G0031-11**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 16:40

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:56

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W46** Lab Sample ID: **21G0031-11**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 16:40

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:56

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W46** Lab Sample ID: **21G0031-11**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 16:40

### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/12/21 23:56

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:07

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:51

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>442</b>    |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>31.9</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| Boron           | < 25.0        |                  | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W46** Lab Sample ID: **21G0031-11**

Matrix: Water Collected By: JO Date/Time Collected: 06/30/21 16:40

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:51

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>95100</b>  |                  | 300                    |
| <b>Chromium</b>  | <b>99.5</b>   |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>423000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>2610</b>   |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>45100</b>  |                  | 300                    |
| <b>Manganese</b> | <b>160</b>    |                  | 15.0                   |
| <b>Nickel</b>    | <b>10.1</b>   |                  | 5.00                   |
| Potassium        | < 1400        |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>10200</b>  |                  | 300                    |
| <b>Strontium</b> | <b>58.9</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W15** Lab Sample ID: **21G0031-12**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:03

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 18:37

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W15** Lab Sample ID: **21G0031-12**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:03

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 18:37

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W15** Lab Sample ID: **21G0031-12**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:03

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:05

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W15** Lab Sample ID: **21G0031-12**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:03

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:05

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W15** Lab Sample ID: **21G0031-12**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:03

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:05

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| <b>Di-n-butylphthalate</b> | <b>2.0</b>    | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W15** Lab Sample ID: **21G0031-12**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:03

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:05

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W15** Lab Sample ID: **21G0031-12**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:03

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:05

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:09

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:56

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>448</b>    |                  | 100                    |
| <b>Antimony</b> | <b>84.9</b>   |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>160</b>    |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| Boron           | < 25.0        | B1               | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W15** Lab Sample ID: **21G0031-12**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:03

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 11:56

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Calcium        | 190000        |                  | 300                    |
| Chromium       | 11600         |                  | 50.0                   |
| Cobalt         | 21.3          |                  | 10.0                   |
| Copper         | 47.6          |                  | 10.0                   |
| Hardness       | 757000        |                  | 1980                   |
| Iron           | 11000         |                  | 150                    |
| Lead           | < 5.00        |                  | 5.00                   |
| Magnesium      | 68400         |                  | 300                    |
| Manganese      | 466           |                  | 15.0                   |
| Nickel         | 974           |                  | 5.00                   |
| Potassium      | 5760          |                  | 1400                   |
| Selenium       | < 25.0        | B1               | 25.0                   |
| Silver         | 3.82          |                  | 3.00                   |
| Sodium         | 451000        |                  | 3000                   |
| Strontium      | 343           |                  | 5.00                   |
| Thallium       | < 10.0        |                  | 10.0                   |
| Vanadium       | 45.0          |                  | 5.00                   |
| Zinc           | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **WIR** Lab Sample ID: **21G0031-13**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:44

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 18:57

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **WIR** Lab Sample ID: **21G0031-13**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:44

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 18:57

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **WIR** Lab Sample ID: **21G0031-13**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:44

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/13/21 00:30

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | Y                | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **WIR** Lab Sample ID: **21G0031-13**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:44

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/13/21 00:30

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **WIR** Lab Sample ID: **21G0031-13**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:44

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/13/21 00:30

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **WIR** Lab Sample ID: **21G0031-13**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:44

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/13/21 00:30

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | Y                | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **WIR** Lab Sample ID: **21G0031-13**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:44

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/06/21 09:30

Units: ug/L Analyzed: 07/13/21 00:30

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:12

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:00

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>7520</b>   |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>78.3</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| Boron           | < 25.0        | B1               | 25.0                   |
| <b>Cadmium</b>  | <b>3.36</b>   |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **WIR** Lab Sample ID: **21G0031-13**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 13:44

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:00

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>170000</b> |                  | 300                    |
| <b>Chromium</b>  | <b>61.2</b>   |                  | 5.00                   |
| <b>Cobalt</b>    | < 10.0        |                  | 10.0                   |
| <b>Copper</b>    | <b>22.1</b>   |                  | 10.0                   |
| <b>Hardness</b>  | <b>724000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>13100</b>  |                  | 150                    |
| <b>Lead</b>      | <b>5.29</b>   |                  | 5.00                   |
| <b>Magnesium</b> | <b>72700</b>  |                  | 300                    |
| <b>Manganese</b> | <b>534</b>    |                  | 15.0                   |
| <b>Nickel</b>    | <b>19.0</b>   |                  | 5.00                   |
| <b>Potassium</b> | <b>3270</b>   |                  | 1400                   |
| <b>Selenium</b>  | < 25.0        | B1               | 25.0                   |
| <b>Silver</b>    | <b>3.19</b>   |                  | 3.00                   |
| <b>Sodium</b>    | <b>45900</b>  |                  | 300                    |
| <b>Strontium</b> | <b>130</b>    |                  | 5.00                   |
| <b>Thallium</b>  | < 10.0        |                  | 10.0                   |
| <b>Vanadium</b>  | <b>17.9</b>   |                  | 5.00                   |
| <b>Zinc</b>      | <b>31.3</b>   |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W17** Lab Sample ID: **21G0031-14**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 14:40

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 21:12

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W17** Lab Sample ID: **21G0031-14**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 14:40

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 21:12

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W17** Lab Sample ID: **21G0031-14**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 14:40

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:40

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W17** Lab Sample ID: **21G0031-14**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 14:40

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:40

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W17** Lab Sample ID: **21G0031-14**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 14:40

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:40

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W17** Lab Sample ID: **21G0031-14**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 14:40

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:40

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W17** Lab Sample ID: **21G0031-14**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 14:40

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 13:40

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:14

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| <b>Mercury</b> | <b>0.10</b>   |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:03

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>8500</b>   |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>78.9</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| Boron           | < 25.0        |                  | 25.0                   |
| <b>Cadmium</b>  | <b>4.24</b>   |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W17** Lab Sample ID: **21G0031-14**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 14:40

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:39

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>211000</b> |                  | 3000                   |
| <b>Chromium</b>  | <b>94.4</b>   |                  | 5.00                   |
| <b>Cobalt</b>    | < 10.0        |                  | 10.0                   |
| <b>Copper</b>    | <b>28.8</b>   |                  | 10.0                   |
| <b>Hardness</b>  | <b>877000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>15900</b>  |                  | 150                    |
| <b>Lead</b>      | <b>6.28</b>   |                  | 5.00                   |
| <b>Magnesium</b> | <b>85200</b>  |                  | 300                    |
| <b>Manganese</b> | <b>639</b>    |                  | 15.0                   |
| <b>Nickel</b>    | <b>27.1</b>   |                  | 5.00                   |
| <b>Potassium</b> | <b>3690</b>   |                  | 1400                   |
| <b>Selenium</b>  | < 25.0        |                  | 25.0                   |
| <b>Silver</b>    | <b>3.44</b>   |                  | 3.00                   |
| <b>Sodium</b>    | <b>10800</b>  |                  | 300                    |
| <b>Strontium</b> | <b>184</b>    |                  | 5.00                   |
| <b>Thallium</b>  | < 10.0        |                  | 10.0                   |
| <b>Vanadium</b>  | <b>24.9</b>   |                  | 5.00                   |
| <b>Zinc</b>      | <b>39.9</b>   |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Volatile Organic Compounds by GC/MS

Method: 524.3 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 12:41

| <u>Analyte</u>           | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|--------------------------|---------------|------------------|------------------------|
| 1,1,1-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1,2-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1-Dichloroethene       | < 0.50        |                  | 0.50                   |
| 1,2-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| 1,2-Dichloroethane       | < 0.50        |                  | 0.50                   |
| 1,2-Dichloropropane      | < 0.50        |                  | 0.50                   |
| 1,4-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| Benzene                  | < 0.50        |                  | 0.50                   |
| Carbon tetrachloride     | < 0.50        |                  | 0.50                   |
| Chlorobenzene            | < 0.50        |                  | 0.50                   |
| cis-1,2-Dichloroethene   | < 0.50        |                  | 0.50                   |
| Ethylbenzene             | < 0.50        |                  | 0.50                   |
| Methyl tert-butyl ether  | < 0.50        |                  | 0.50                   |
| Methylene chloride       | < 0.50        |                  | 0.50                   |
| Styrene                  | < 0.50        |                  | 0.50                   |
| Tetrachloroethene        | < 0.50        |                  | 0.50                   |
| Toluene                  | < 0.50        |                  | 0.50                   |
| trans-1,2-Dichloroethene | < 0.50        |                  | 0.50                   |
| Trichloroethene          | < 0.50        |                  | 0.50                   |
| Vinyl chloride           | < 0.50        | J5               | 0.50                   |
| Xylenes, total           | < 0.50        |                  | 0.50                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 21:34

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 21:34

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:14

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:14

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:14

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:14

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:14

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:16

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:11

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>452</b>    |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>19.2</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>114</b>    |                  | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **G103D** Lab Sample ID: **21G0031-15**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:11

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>95700</b>  |                  | 300                    |
| <b>Chromium</b>  | <b>9.36</b>   |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>425000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>907</b>    |                  | 150                    |
| <b>Lead</b>      | <b>8.62</b>   |                  | 5.00                   |
| <b>Magnesium</b> | <b>45200</b>  |                  | 300                    |
| <b>Manganese</b> | <b>23.4</b>   |                  | 15.0                   |
| <b>Nickel</b>    | <b>8.95</b>   |                  | 5.00                   |
| <b>Potassium</b> | <b>1950</b>   |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>19800</b>  |                  | 300                    |
| <b>Strontium</b> | <b>72.4</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Volatile Organic Compounds by GC/MS

Method: 524.3 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 13:18

| <u>Analyte</u>           | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|--------------------------|---------------|------------------|------------------------|
| 1,1,1-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1,2-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1-Dichloroethene       | < 0.50        |                  | 0.50                   |
| 1,2-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| 1,2-Dichloroethane       | < 0.50        |                  | 0.50                   |
| 1,2-Dichloropropane      | < 0.50        |                  | 0.50                   |
| 1,4-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| Benzene                  | < 0.50        |                  | 0.50                   |
| Carbon tetrachloride     | < 0.50        |                  | 0.50                   |
| Chlorobenzene            | < 0.50        |                  | 0.50                   |
| cis-1,2-Dichloroethene   | < 0.50        |                  | 0.50                   |
| Ethylbenzene             | < 0.50        |                  | 0.50                   |
| Methyl tert-butyl ether  | < 0.50        |                  | 0.50                   |
| Methylene chloride       | < 0.50        |                  | 0.50                   |
| Styrene                  | < 0.50        |                  | 0.50                   |
| Tetrachloroethene        | < 0.50        |                  | 0.50                   |
| Toluene                  | < 0.50        |                  | 0.50                   |
| trans-1,2-Dichloroethene | < 0.50        |                  | 0.50                   |
| Trichloroethene          | < 0.50        |                  | 0.50                   |
| Vinyl chloride           | < 0.50        | J5               | 0.50                   |
| Xylenes, total           | < 0.50        |                  | 0.50                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 21:55

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 21:55

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:49

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:49

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:49

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:49

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 14:49

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:23

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:14

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>332</b>    |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>18.2</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>112</b>    |                  | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **X103D** Lab Sample ID: **21G0031-16**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 17:55

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:14

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>93500</b>  |                  | 300                    |
| <b>Chromium</b>  | <b>6.64</b>   |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>413000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>663</b>    |                  | 150                    |
| <b>Lead</b>      | <b>6.23</b>   |                  | 5.00                   |
| <b>Magnesium</b> | <b>43600</b>  |                  | 300                    |
| <b>Manganese</b> | <b>18.4</b>   |                  | 15.0                   |
| <b>Nickel</b>    | <b>7.04</b>   |                  | 5.00                   |
| <b>Potassium</b> | <b>1600</b>   |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>19400</b>  |                  | 300                    |
| <b>Strontium</b> | <b>70.6</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

### Volatile Organic Compounds by GC/MS

Method: 524.3 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 13:55

| <u>Analyte</u>           | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|--------------------------|---------------|------------------|------------------------|
| 1,1,1-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1,2-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1-Dichloroethene       | < 0.50        |                  | 0.50                   |
| 1,2-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| 1,2-Dichloroethane       | < 0.50        |                  | 0.50                   |
| 1,2-Dichloropropane      | < 0.50        |                  | 0.50                   |
| 1,4-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| Benzene                  | < 0.50        |                  | 0.50                   |
| Carbon tetrachloride     | < 0.50        |                  | 0.50                   |
| Chlorobenzene            | < 0.50        |                  | 0.50                   |
| cis-1,2-Dichloroethene   | < 0.50        |                  | 0.50                   |
| Ethylbenzene             | < 0.50        |                  | 0.50                   |
| Methyl tert-butyl ether  | < 0.50        |                  | 0.50                   |
| Methylene chloride       | < 0.50        | J3               | 0.50                   |
| Styrene                  | < 0.50        | J3               | 0.50                   |
| Tetrachloroethene        | < 0.50        |                  | 0.50                   |
| Toluene                  | < 0.50        |                  | 0.50                   |
| trans-1,2-Dichloroethene | < 0.50        |                  | 0.50                   |
| Trichloroethene          | < 0.50        |                  | 0.50                   |
| Vinyl chloride           | < 0.50        | J3, J5           | 0.50                   |
| Xylenes, total           | < 0.50        |                  | 0.50                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 09:04

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 09:04

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         | J3               | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         | J3               | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:23

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | J3, O2Y          | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:23

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:23

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:23

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:23

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:25

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:17

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>6580</b>   |                  | 100                    |
| <b>Antimony</b> | <b>14.5</b>   |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>85.5</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| Boron           | < 25.0        | B1               | 25.0                   |
| <b>Cadmium</b>  | <b>6.19</b>   |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W29** Lab Sample ID: **21G0031-17**

Matrix: Water Collected By: KHW Date/Time Collected: 06/30/21 19:00

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:17

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Calcium        | 123000        |                  | 300                    |
| Chromium       | 2070          |                  | 5.00                   |
| Cobalt         | 56.3          |                  | 10.0                   |
| Copper         | 48.9          |                  | 10.0                   |
| Hardness       | 573000        |                  | 1980                   |
| Iron           | 27800         |                  | 150                    |
| Lead           | 10.7          |                  | 5.00                   |
| Magnesium      | 64500         |                  | 300                    |
| Manganese      | 924           |                  | 15.0                   |
| Nickel         | 244           |                  | 5.00                   |
| Potassium      | 3080          |                  | 1400                   |
| Selenium       | < 25.0        | B1               | 25.0                   |
| Silver         | < 3.00        |                  | 3.00                   |
| Sodium         | 56700         |                  | 300                    |
| Strontium      | 80.1          |                  | 5.00                   |
| Thallium       | < 10.0        |                  | 10.0                   |
| Vanadium       | 36.3          |                  | 5.00                   |
| Zinc           | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W41** Lab Sample ID: **21G0031-18**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 13:30

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 09:57

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W41** Lab Sample ID: **21G0031-18**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 13:30

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 09:57

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W41** Lab Sample ID: **21G0031-18**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 13:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:58

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W41** Lab Sample ID: **21G0031-18**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 13:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:58

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W41** Lab Sample ID: **21G0031-18**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 13:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:58

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W41** Lab Sample ID: **21G0031-18**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 13:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:58

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W41** Lab Sample ID: **21G0031-18**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 13:30

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 15:58

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:28

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:19

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>410</b>    |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>59.1</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>78.7</b>   | B1               | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W41** Lab Sample ID: **21G0031-18**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 13:30

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:19

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>108000</b> |                  | 300                    |
| <b>Chromium</b>  | <b>550</b>    |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| <b>Copper</b>    | <b>34.4</b>   |                  | 10.0                   |
| <b>Hardness</b>  | <b>439000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>4090</b>   |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>40900</b>  |                  | 300                    |
| <b>Manganese</b> | <b>42.9</b>   |                  | 15.0                   |
| <b>Nickel</b>    | <b>37.3</b>   |                  | 5.00                   |
| <b>Potassium</b> | <b>1400</b>   |                  | 1400                   |
| Selenium         | < 25.0        | B1               | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>60900</b>  |                  | 300                    |
| <b>Strontium</b> | <b>121</b>    |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W53** Lab Sample ID: **21G0031-19**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:20

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 10:18

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W53** Lab Sample ID: **21G0031-19**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:20

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 10:18

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W53** Lab Sample ID: **21G0031-19**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:20

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 16:32

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W53** Lab Sample ID: **21G0031-19**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:20

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 16:32

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W53** Lab Sample ID: **21G0031-19**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:20

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 16:32

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W53** Lab Sample ID: **21G0031-19**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:20

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 16:32

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W53** Lab Sample ID: **21G0031-19**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:20

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 16:32

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:30

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:22

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>2050</b>   |                  | 100                    |
| Antimony        | < 10.0        |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>63.7</b>   |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| <b>Boron</b>    | <b>35.4</b>   |                  | 25.0                   |
| Cadmium         | < 3.00        |                  | 3.00                   |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W53** Lab Sample ID: **21G0031-19**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:20

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:22

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>64700</b>  |                  | 300                    |
| <b>Chromium</b>  | <b>96.8</b>   |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| <b>Copper</b>    | <b>12.7</b>   |                  | 10.0                   |
| <b>Hardness</b>  | <b>273000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>4890</b>   |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>27000</b>  |                  | 300                    |
| <b>Manganese</b> | <b>133</b>    |                  | 15.0                   |
| <b>Nickel</b>    | <b>49.3</b>   |                  | 5.00                   |
| <b>Potassium</b> | <b>2420</b>   |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>529000</b> |                  | 3000                   |
| <b>Strontium</b> | <b>131</b>    |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| <b>Vanadium</b>  | <b>6.04</b>   |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W52** Lab Sample ID: **21G0031-20**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:45

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 10:39

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W52** Lab Sample ID: **21G0031-20**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:45

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 10:39

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| <b>Tetrachloroethene</b>  | <b>5.4</b>    |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W52** Lab Sample ID: **21G0031-20**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:07

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W52** Lab Sample ID: **21G0031-20**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:07

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W52** Lab Sample ID: **21G0031-20**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:07

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W52** Lab Sample ID: **21G0031-20**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:07

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W52** Lab Sample ID: **21G0031-20**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:45

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:07

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:32

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | 0.12          |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:44

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Aluminum       | 47400         |                  | 1000                   |
| Antimony       | < 10.0        |                  | 10.0                   |
| Arsenic        | 17.7          |                  | 10.0                   |
| Barium         | 379           |                  | 10.0                   |
| Beryllium      | < 1.00        |                  | 1.00                   |
| Boron          | < 25.0        | B1               | 25.0                   |
| Cadmium        | 21.8          |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W52** Lab Sample ID: **21G0031-20**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 14:45

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:44

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Calcium        | 554000        |                  | 3000                   |
| Chromium       | 2390          |                  | 5.00                   |
| Cobalt         | 58.3          |                  | 10.0                   |
| Copper         | 198           |                  | 10.0                   |
| Hardness       | 2430000       |                  | 1980                   |
| Iron           | 102000        |                  | 1500                   |
| Lead           | 37.2          |                  | 5.00                   |
| Magnesium      | 255000        |                  | 3000                   |
| Manganese      | 2040          |                  | 15.0                   |
| Nickel         | 739           |                  | 5.00                   |
| Potassium      | 9350          |                  | 1400                   |
| Selenium       | < 25.0        | B1               | 25.0                   |
| Silver         | 8.74          |                  | 3.00                   |
| Sodium         | 74900         |                  | 300                    |
| Strontium      | 377           |                  | 5.00                   |
| Thallium       | < 10.0        |                  | 10.0                   |
| Vanadium       | 112           |                  | 5.00                   |
| Zinc           | 265           |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

### Volatile Organic Compounds by GC/MS

Method: 524.3 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 14:32

| <u>Analyte</u>           | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|--------------------------|---------------|------------------|------------------------|
| 1,1,1-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1,2-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1-Dichloroethene       | < 0.50        |                  | 0.50                   |
| 1,2-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| 1,2-Dichloroethane       | < 0.50        |                  | 0.50                   |
| 1,2-Dichloropropane      | < 0.50        |                  | 0.50                   |
| 1,4-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| Benzene                  | < 0.50        |                  | 0.50                   |
| Carbon tetrachloride     | < 0.50        |                  | 0.50                   |
| Chlorobenzene            | < 0.50        |                  | 0.50                   |
| cis-1,2-Dichloroethene   | < 0.50        |                  | 0.50                   |
| Ethylbenzene             | < 0.50        |                  | 0.50                   |
| Methyl tert-butyl ether  | < 0.50        |                  | 0.50                   |
| Methylene chloride       | < 0.50        |                  | 0.50                   |
| Styrene                  | < 0.50        |                  | 0.50                   |
| Tetrachloroethene        | < 0.50        |                  | 0.50                   |
| Toluene                  | < 0.50        |                  | 0.50                   |
| trans-1,2-Dichloroethene | < 0.50        |                  | 0.50                   |
| Trichloroethene          | < 0.50        |                  | 0.50                   |
| Vinyl chloride           | < 0.50        | J5               | 0.50                   |
| Xylenes, total           | < 0.50        |                  | 0.50                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 10:59

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 10:59

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:41

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | J1 Y             | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | J1 Y             | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | J1 Y             | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | J1 Y             | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | J1 Y             | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | J1 Y             | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | J1 Y             | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | J1 Y             | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | J1 Y             | 1.5                    |
| 1-Naphthylamine             | < 5.0         | J1 Y             | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | J1 Y             | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | J1 Y             | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | J1 Y             | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | J1 Y             | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | J1 Y             | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | J1 Y             | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | J1, O2Y          | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | J1 Y             | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | J1 Y             | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | J1 Y             | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | J1 Y             | 1.5                    |
| 2-Chlorophenol              | < 1.5         | J1 Y             | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | J1 Y             | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:41

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | J1 Y             | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | J1 Y             | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | J1 Y             | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | J1 Y             | 5.0                    |
| 2-Picoline                      | < 1.5         | J1 Y             | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | J1 Y             | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | J1 Y             | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | J1 Y             | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | J1 Y             | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | J1 Y             | 1.5                    |
| 4-Methylphenol                  | < 1.5         | J1 Y             | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | J1 Y             | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | J1, O2Y          | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | J1 Y             | 1.5                    |
| Acenaphthylene                  | < 1.5         | J1 Y             | 1.5                    |
| Acetophenone                    | < 1.5         | J1 Y             | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:41

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | J1 Y             | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | J1 Y             | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | J1 Y             | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | J1 Y             | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | J1 Y             | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | J1 Y             | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | J1 Y             | 1.5                    |
| Diethylphthalate           | < 1.5         | J1 Y             | 1.5                    |
| Dimethylphthalate          | < 1.5         | J1 Y             | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | J1 Y             | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | J1 Y             | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | J1 Y             | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:41

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | J1 Y             | 1.5                    |
| Hexachloroethane          | < 1.5         | J1 Y             | 1.5                    |
| Hexachloropropene         | < 1.5         | J1 Y             | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | J1 Y             | 1.5                    |
| Isosafrole                | < 1.5         | J1 Y             | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | J1 Y             | 1.5                    |
| Naphthalene               | < 1.5         | J1 Y             | 1.5                    |
| Nitrobenzene              | < 1.5         | J1 Y             | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | J1 Y             | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | J1 Y             | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | J1 Y             | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | J1 Y             | 1.5                    |
| Pentachlorobenzene        | < 1.5         | J1 Y             | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | J1 Y             | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | J1 Y             | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 17:41

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | J1 Y             | 1.5                    |
| Safrole        | < 1.5         | J1 Y             | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:44

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:29

| <u>Analyte</u>  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------|---------------|------------------|------------------------|
| <b>Aluminum</b> | <b>1210</b>   |                  | 100                    |
| <b>Antimony</b> | <b>69.5</b>   |                  | 10.0                   |
| Arsenic         | < 10.0        |                  | 10.0                   |
| <b>Barium</b>   | <b>215</b>    |                  | 10.0                   |
| Beryllium       | < 1.00        |                  | 1.00                   |
| Boron           | < 25.0        | B1               | 25.0                   |
| <b>Cadmium</b>  | <b>19.2</b>   |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W37** Lab Sample ID: **21G0031-21**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 16:35

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 12:29

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>122000</b> |                  | 300                    |
| <b>Chromium</b>  | <b>9890</b>   |                  | 50.0                   |
| <b>Cobalt</b>    | <b>143</b>    |                  | 10.0                   |
| <b>Copper</b>    | <b>219</b>    |                  | 10.0                   |
| <b>Hardness</b>  | <b>504000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>68400</b>  |                  | 1500                   |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>48100</b>  |                  | 300                    |
| <b>Manganese</b> | <b>1190</b>   |                  | 15.0                   |
| <b>Nickel</b>    | <b>1150</b>   |                  | 5.00                   |
| <b>Potassium</b> | <b>2200</b>   |                  | 1400                   |
| Selenium         | < 25.0        | B1               | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>121000</b> |                  | 300                    |
| <b>Strontium</b> | <b>99.8</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| <b>Vanadium</b>  | <b>36.8</b>   |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

### Volatile Organic Compounds by GC/MS

Method: 524.3 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 15:08

| <u>Analyte</u>           | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|--------------------------|---------------|------------------|------------------------|
| 1,1,1-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1,2-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1-Dichloroethene       | < 0.50        |                  | 0.50                   |
| 1,2-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| 1,2-Dichloroethane       | < 0.50        |                  | 0.50                   |
| 1,2-Dichloropropane      | < 0.50        |                  | 0.50                   |
| 1,4-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| Benzene                  | < 0.50        |                  | 0.50                   |
| Carbon tetrachloride     | < 0.50        |                  | 0.50                   |
| Chlorobenzene            | < 0.50        |                  | 0.50                   |
| cis-1,2-Dichloroethene   | < 0.50        |                  | 0.50                   |
| Ethylbenzene             | < 0.50        |                  | 0.50                   |
| Methyl tert-butyl ether  | < 0.50        |                  | 0.50                   |
| Methylene chloride       | < 0.50        |                  | 0.50                   |
| Styrene                  | < 0.50        |                  | 0.50                   |
| Tetrachloroethene        | < 0.50        |                  | 0.50                   |
| Toluene                  | < 0.50        |                  | 0.50                   |
| trans-1,2-Dichloroethene | < 0.50        |                  | 0.50                   |
| Trichloroethene          | < 0.50        |                  | 0.50                   |
| Vinyl chloride           | < 0.50        | J5               | 0.50                   |
| Xylenes, total           | < 0.50        |                  | 0.50                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 11:20

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 11:20

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:16

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:16

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:16

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:16

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:16

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:51

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:07

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Aluminum       | < 100         |                  | 100                    |
| Antimony       | < 10.0        |                  | 10.0                   |
| Arsenic        | < 10.0        |                  | 10.0                   |
| Barium         | < 10.0        |                  | 10.0                   |
| Beryllium      | < 1.00        |                  | 1.00                   |
| Boron          | < 25.0        | B1               | 25.0                   |
| Cadmium        | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **FB** Lab Sample ID: **21G0031-22**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 18:15

### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:07

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Calcium        | < 300         |                  | 300                    |
| Chromium       | < 5.00        |                  | 5.00                   |
| Cobalt         | < 10.0        |                  | 10.0                   |
| Copper         | < 10.0        |                  | 10.0                   |
| Hardness       | < 1980        |                  | 1980                   |
| Iron           | < 150         |                  | 150                    |
| Lead           | < 5.00        |                  | 5.00                   |
| Magnesium      | < 300         |                  | 300                    |
| Manganese      | < 15.0        |                  | 15.0                   |
| Nickel         | < 5.00        |                  | 5.00                   |
| Potassium      | < 1400        |                  | 1400                   |
| Selenium       | < 25.0        | B1               | 25.0                   |
| Silver         | < 3.00        |                  | 3.00                   |
| Sodium         | < 300         |                  | 300                    |
| Strontium      | < 5.00        |                  | 5.00                   |
| Thallium       | < 10.0        |                  | 10.0                   |
| Vanadium       | < 5.00        |                  | 5.00                   |
| Zinc           | < 25.0        |                  | 25.0                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **TB1** Lab Sample ID: **21G0031-23**

Matrix: Water Collected By: Date/Time Collected: 06/30/21 19:50

### Volatile Organic Compounds by GC/MS

Method: 524.3 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/09/21 15:45

| <u>Analyte</u>           | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|--------------------------|---------------|------------------|------------------------|
| 1,1,1-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1,2-Trichloroethane    | < 0.50        |                  | 0.50                   |
| 1,1-Dichloroethene       | < 0.50        |                  | 0.50                   |
| 1,2-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| 1,2-Dichloroethane       | < 0.50        |                  | 0.50                   |
| 1,2-Dichloropropane      | < 0.50        |                  | 0.50                   |
| 1,4-Dichlorobenzene      | < 0.50        |                  | 0.50                   |
| Benzene                  | < 0.50        |                  | 0.50                   |
| Carbon tetrachloride     | < 0.50        |                  | 0.50                   |
| Chlorobenzene            | < 0.50        |                  | 0.50                   |
| cis-1,2-Dichloroethene   | < 0.50        |                  | 0.50                   |
| Ethylbenzene             | < 0.50        |                  | 0.50                   |
| Methyl tert-butyl ether  | < 0.50        |                  | 0.50                   |
| Methylene chloride       | < 0.50        |                  | 0.50                   |
| Styrene                  | < 0.50        |                  | 0.50                   |
| Tetrachloroethene        | < 0.50        |                  | 0.50                   |
| Toluene                  | < 0.50        |                  | 0.50                   |
| trans-1,2-Dichloroethene | < 0.50        |                  | 0.50                   |
| Trichloroethene          | < 0.50        |                  | 0.50                   |
| Vinyl chloride           | < 0.50        | J5               | 0.50                   |
| Xylenes, total           | < 0.50        |                  | 0.50                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **TB1** Lab Sample ID: **21G0031-23**

Matrix: Water Collected By: Date/Time Collected: 06/30/21 19:50

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 12:23

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **TB1** Lab Sample ID: **21G0031-23**

Matrix: Water Collected By: Date/Time Collected: 06/30/21 19:50

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 12:23

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

|                          |                 |                 |          |
|--------------------------|-----------------|-----------------|----------|
| Name:                    | <b>CHEMTOOL</b> | Date Received : | 07/01/21 |
| Project/Facility Number: | 2010355004      | Temperature C:  | 3.00     |
| Funding Code:            | CS29 B50        |                 |          |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W6** Lab Sample ID: **21G0031-24**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 19:50

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 12:44

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,1,1,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,1-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1,2,2-Tetrachloroethane   | < 2.0         |                  | 2.0                    |
| 1,1,2-Trichloroethane       | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloroethene          | < 2.0         |                  | 2.0                    |
| 1,1-Dichloropropene         | < 2.0         |                  | 2.0                    |
| 1,2,3-Trichloropropane      | < 2.0         |                  | 2.0                    |
| 1,2-Dibromoethane           | < 2.0         |                  | 2.0                    |
| 1,2-Dichloroethane          | < 2.0         |                  | 2.0                    |
| 1,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 1,3-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2,2-Dichloropropane         | < 2.0         |                  | 2.0                    |
| 2-Butanone (MEK)            | < 10          |                  | 10                     |
| 2-Hexanone (MBK)            | < 5.0         |                  | 5.0                    |
| 4-Methyl-2-pentanone (MIBK) | < 10          |                  | 10                     |
| Acetone                     | < 10          |                  | 10                     |
| Benzene                     | < 2.0         |                  | 2.0                    |
| Bromobenzene                | < 2.0         |                  | 2.0                    |
| Bromochloromethane          | < 2.0         |                  | 2.0                    |
| Bromodichloromethane        | < 2.0         |                  | 2.0                    |
| Bromoform                   | < 5.0         |                  | 5.0                    |
| Bromomethane                | < 5.0         |                  | 5.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W6** Lab Sample ID: **21G0031-24**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 19:50

### **Volatiles Organic Compounds by Purge and Trap GC/MS**

Method: 8260 Prepared: 07/09/21 08:00

Units: ug/L Analyzed: 07/12/21 12:44

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Carbon disulfide          | < 2.0         |                  | 2.0                    |
| Carbon tetrachloride      | < 2.0         |                  | 2.0                    |
| Chlorobenzene             | < 2.0         |                  | 2.0                    |
| Chloroethane              | < 2.0         |                  | 2.0                    |
| Chloroform                | < 2.0         |                  | 2.0                    |
| Chloromethane             | < 2.0         |                  | 2.0                    |
| cis-1,2-Dichloroethene    | < 2.0         |                  | 2.0                    |
| cis-1,3-Dichloropropene   | < 2.0         |                  | 2.0                    |
| Dibromochloromethane      | < 5.0         |                  | 5.0                    |
| Dibromomethane            | < 2.0         |                  | 2.0                    |
| Ethylbenzene              | < 2.0         |                  | 2.0                    |
| Isopropylbenzene          | < 2.0         |                  | 2.0                    |
| Methyl tert-butyl ether   | < 2.0         |                  | 2.0                    |
| Methylene chloride        | < 5.0         |                  | 5.0                    |
| Styrene                   | < 2.0         |                  | 2.0                    |
| Tetrachloroethene         | < 2.0         |                  | 2.0                    |
| Toluene                   | < 2.0         |                  | 2.0                    |
| trans-1,2-Dichloroethene  | < 2.0         |                  | 2.0                    |
| trans-1,3-Dichloropropene | < 5.0         |                  | 5.0                    |
| Trichloroethene           | < 2.0         |                  | 2.0                    |
| Trichlorofluoromethane    | < 2.0         |                  | 2.0                    |
| Vinyl chloride            | < 2.0         |                  | 2.0                    |
| Xylenes, total            | < 2.0         |                  | 2.0                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W6** Lab Sample ID: **21G0031-24**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 19:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:50

| <u>Analyte</u>              | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|-----------------------------|---------------|------------------|------------------------|
| 1,2,4,5-Tetrachlorobenzene  | < 1.5         | Y                | 1.5                    |
| 1,2,4-Trichlorobenzene      | < 1.5         | Y                | 1.5                    |
| 1,2-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,2-Dinitrobenzene          | < 1.5         | Y                | 1.5                    |
| 1,3-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,3-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1,4-Dichlorobenzene         | < 1.5         | Y                | 1.5                    |
| 1,4-Dinitrobenzene          | < 5.0         | Y                | 5.0                    |
| 1-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 1-Naphthylamine             | < 5.0         | Y                | 5.0                    |
| 2,2-Oxybis(1-chloropropane) | < 1.5         | Y                | 1.5                    |
| 2,3,4,6-Tetrachlorophenol   | < 1.5         | Y                | 1.5                    |
| 2,4,5-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4,6-Trichlorophenol       | < 1.5         | Y                | 1.5                    |
| 2,4-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dimethylphenol          | < 1.5         | Y                | 1.5                    |
| 2,4-Dinitrophenol           | < 5.0         | O2 Y             | 5.0                    |
| 2,4-Dinitrotoluene          | < 5.0         | Y                | 5.0                    |
| 2,6-Dichlorophenol          | < 1.5         | Y                | 1.5                    |
| 2,6-Dinitrotoluene          | < 1.5         | Y                | 1.5                    |
| 2-Chloronaphthalene         | < 1.5         | Y                | 1.5                    |
| 2-Chlorophenol              | < 1.5         | Y                | 1.5                    |
| 2-Methylnaphthalene         | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W6** Lab Sample ID: **21G0031-24**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 19:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:50

| <u>Analyte</u>                  | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------------|---------------|------------------|------------------------|
| 2-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 2-Naphthylamine                 | < 5.0         | Y                | 5.0                    |
| 2-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 2-Nitrophenol                   | < 5.0         | Y                | 5.0                    |
| 2-Picoline                      | < 1.5         | Y                | 1.5                    |
| 3,3-Dichlorobenzidine           | < 1.5         | Y                | 1.5                    |
| 3-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4,6-Dinitro-2-methylphenol      | < 5.0         | Y                | 5.0                    |
| 4-Bromophenyl phenyl ether      | < 1.5         | Y                | 1.5                    |
| 4-Chloro-3-methylphenol         | < 1.5         | Y                | 1.5                    |
| 4-Chloroaniline                 | < 1.5         | Y                | 1.5                    |
| 4-Chlorophenyl phenyl ether     | < 1.5         | Y                | 1.5                    |
| 4-Methylphenol                  | < 1.5         | Y                | 1.5                    |
| 4-Nitroaniline                  | < 1.5         | Y                | 1.5                    |
| 4-Nitrobiphenyl                 | < 5.0         | Y                | 5.0                    |
| 4-Nitrophenol                   | < 5.0         | O2 Y             | 5.0                    |
| 5-Nitroacenaphthene             | < 5.0         | Y                | 5.0                    |
| 7,12-Dimethylbenzo(a)anthracene | < 5.0         | Y                | 5.0                    |
| Acenaphthene                    | < 1.5         | Y                | 1.5                    |
| Acenaphthylene                  | < 1.5         | Y                | 1.5                    |
| Acetophenone                    | < 1.5         | Y                | 1.5                    |
| Anthracene                      | < 1.5         | Y                | 1.5                    |
| Azobenzene                      | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W6** Lab Sample ID: **21G0031-24**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 19:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:50

| <u>Analyte</u>             | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------------------|---------------|------------------|------------------------|
| Benzo(a)anthracene         | < 1.5         | Y                | 1.5                    |
| Benzo(a)pyrene             | < 1.5         | Y                | 1.5                    |
| Benzo(b)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Benzo(ghi)perylene         | < 5.0         | Y                | 5.0                    |
| Benzo(k)fluoranthene       | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethoxy)methane | < 1.5         | Y                | 1.5                    |
| Bis(2-chloroethyl)ether    | < 1.5         | Y                | 1.5                    |
| Bis(2-ethylhexyl)phthalate | < 5.0         | Y                | 5.0                    |
| Butyl benzyl phthalate     | < 5.0         | Y                | 5.0                    |
| Carbazole                  | < 1.5         | Y                | 1.5                    |
| Chrysene                   | < 1.5         | Y                | 1.5                    |
| Dibenzo(a,h)anthracene     | < 5.0         | Y                | 5.0                    |
| Dibenzofuran               | < 1.5         | Y                | 1.5                    |
| Diethylphthalate           | < 1.5         | Y                | 1.5                    |
| Dimethylphthalate          | < 1.5         | Y                | 1.5                    |
| Di-n-butylphthalate        | < 1.5         | Y                | 1.5                    |
| Di-n-octylphthalate        | < 5.0         | Y                | 5.0                    |
| Diphenylamine              | < 1.5         | Y                | 1.5                    |
| Ethyl methanesulfonate     | < 1.5         | Y                | 1.5                    |
| Fluoranthene               | < 1.5         | Y                | 1.5                    |
| Fluorene                   | < 1.5         | Y                | 1.5                    |
| Hexachlorobenzene          | < 1.5         | Y                | 1.5                    |
| Hexachlorobutadiene        | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W6** Lab Sample ID: **21G0031-24**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 19:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:50

| <u>Analyte</u>            | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|---------------------------|---------------|------------------|------------------------|
| Hexachlorocyclopentadiene | < 1.5         | Y                | 1.5                    |
| Hexachloroethane          | < 1.5         | Y                | 1.5                    |
| Hexachloropropene         | < 1.5         | Y                | 1.5                    |
| Indeno(1,2,3-cd)pyrene    | < 5.0         | Y                | 5.0                    |
| Isodrin                   | < 1.5         | Y                | 1.5                    |
| Isophorone                | < 1.5         | Y                | 1.5                    |
| Isosafrole                | < 1.5         | Y                | 1.5                    |
| Mestranol                 | < 5.0         | Y                | 5.0                    |
| Methyl methanesulfonate   | < 1.5         | Y                | 1.5                    |
| Naphthalene               | < 1.5         | Y                | 1.5                    |
| Nitrobenzene              | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-butylamine  | < 1.5         | Y                | 1.5                    |
| N-Nitrosodi-n-propylamine | < 1.5         | Y                | 1.5                    |
| N-Nitrosopiperidine       | < 1.5         | Y                | 1.5                    |
| p-Dimethylaminoazobenzene | < 1.5         | Y                | 1.5                    |
| Pentachlorobenzene        | < 1.5         | Y                | 1.5                    |
| Pentachloronitrobenzene   | < 1.5         | Y                | 1.5                    |
| Pentachlorophenol         | < 5.0         | O2 Y             | 5.0                    |
| Phenacetin                | < 1.5         | Y                | 1.5                    |
| Phenanthrene              | < 1.5         | Y                | 1.5                    |
| Phenol                    | < 1.5         | Y                | 1.5                    |
| Pronamide                 | < 1.5         | Y                | 1.5                    |
| Pyrene                    | < 1.5         | Y                | 1.5                    |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W6** Lab Sample ID: **21G0031-24**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 19:50

#### **Semivolatiles by GC/MS**

Method: 8270 Prepared: 07/07/21 13:40

Units: ug/L Analyzed: 07/13/21 18:50

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Pyridine       | < 1.5         | Y                | 1.5                    |
| Safrole        | < 1.5         | Y                | 1.5                    |

#### **Mercury by EPA Method 245.1**

Method: 245.1 Prepared: 07/06/21 08:15

Units: ug/L Analyzed: 07/07/21 11:53

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Mercury        | < 0.06        |                  | 0.06                   |

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:10

| <u>Analyte</u> | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|----------------|---------------|------------------|------------------------|
| Aluminum       | < 100         |                  | 100                    |
| Antimony       | < 10.0        |                  | 10.0                   |
| Arsenic        | < 10.0        |                  | 10.0                   |
| <b>Barium</b>  | <b>25.3</b>   |                  | 10.0                   |
| Beryllium      | < 1.00        |                  | 1.00                   |
| Boron          | < 25.0        |                  | 25.0                   |
| Cadmium        | < 3.00        |                  | 3.00                   |

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## Illinois Environmental Protection Agency Laboratory

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

Client Sample ID: **W6** Lab Sample ID: **21G0031-24**

Matrix: Water Collected By: MR Date/Time Collected: 06/30/21 19:50

#### **Metals by EPA 6000/7000 Series Methods**

Method: SW-846 6010 Prepared: 07/02/21 08:50

Units: ug/L Analyzed: 07/08/21 13:10

| <u>Analyte</u>   | <u>Result</u> | <u>Qualifier</u> | <u>Reporting Limit</u> |
|------------------|---------------|------------------|------------------------|
| <b>Calcium</b>   | <b>131000</b> |                  | 300                    |
| Chromium         | < 5.00        |                  | 5.00                   |
| Cobalt           | < 10.0        |                  | 10.0                   |
| Copper           | < 10.0        |                  | 10.0                   |
| <b>Hardness</b>  | <b>557000</b> |                  | 1980                   |
| <b>Iron</b>      | <b>156</b>    |                  | 150                    |
| Lead             | < 5.00        |                  | 5.00                   |
| <b>Magnesium</b> | <b>55700</b>  |                  | 300                    |
| Manganese        | < 15.0        |                  | 15.0                   |
| Nickel           | < 5.00        |                  | 5.00                   |
| Potassium        | < 1400        |                  | 1400                   |
| Selenium         | < 25.0        |                  | 25.0                   |
| Silver           | < 3.00        |                  | 3.00                   |
| <b>Sodium</b>    | <b>2650</b>   | V                | 300                    |
| <b>Strontium</b> | <b>99.1</b>   |                  | 5.00                   |
| Thallium         | < 10.0        |                  | 10.0                   |
| Vanadium         | < 5.00        |                  | 5.00                   |
| Zinc             | < 25.0        |                  | 25.0                   |

*The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety. Test results meet all requirements of NELAC (accredited by Florida DOH #E37645). If you have any questions about this report, please contact Tom Weiss, Laboratory Manager, at 217.782.9780.*

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## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 07/01/21

Funding Code: CS29 B50 Temperature C: 3.00

#### **Notes and Definitions**

- Y The laboratory analysis was performed on an unpreserved or improperly preserved sample.
- V Indicates the analyte was detected in both the sample and the associated method blank and was outside method blank acceptance criteria.
- O2 Quality control sample failed low - possible low bias or false non-detect result.
- J5 Blank spike failed high, result was less than the reporting limit - impact on data may be minimal.
- J3 The reported value failed to meet the established quality control criteria for either precision or accuracy possibly due to matrix effects.
- J1 Surrogate compound recovery limits have not been met.
- B1 The sample matrix caused possible effects on measurement. The result may be biased low.
- ND Analyte NOT DETECTED at or above the reporting limit
- \* Non-NELAP accredited

Method 8270: Samples are Yqualified due to a tripped breaker shutting down the refrigerator where 8270 extracts were stored prior to analysis.

Method 8270: Tentatively Identified Compounds (TICs) were detected in the semi-volatile analysis of samples 21G0031-13. Please contact the laboratory if additional information about the TICs is needed.

Method 8270: There was insufficient amount of sample received to perform a matrix spike duplicate. A matrix spike was analyzed, however, NELAC and method requirements were not met due to the absence of a duplicate matrix spike.

Report Authorized by:

Tom Weiss  
Laboratory Manager

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