



## Illinois Environmental Protection Agency Laboratory

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G106** Lab Sample ID: **21F1058-01**

Matrix: Water Collected By: MW Date/Time Collected: 06/25/21 15:10

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

Method: 8260 Prepared: 06/27/21 10:50

Units: ug/L Analyzed: 06/27/21 18:11

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
1,1,1,2-Tetrachloroethane	< 2.0	Y	2.0
1,1,1-Trichloroethane	< 2.0	Y	2.0
1,1,2,2-Tetrachloroethane	< 2.0	Y	2.0
1,1,2-Trichloroethane	< 2.0	Y	2.0
1,1-Dichloroethane	< 2.0	Y	2.0
1,1-Dichloroethene	< 2.0	Y	2.0
1,1-Dichloropropene	< 2.0	Y	2.0
1,2,3-Trichloropropane	< 2.0	Y	2.0
1,2-Dibromoethane	< 2.0	Y	2.0
1,2-Dichloroethane	< 2.0	Y	2.0
1,2-Dichloropropane	< 2.0	Y	2.0
1,3-Dichloropropane	< 2.0	Y	2.0
2,2-Dichloropropane	< 2.0	Y	2.0
<b>2-Butanone (MEK)</b>	<b>14</b>	Y	10
2-Hexanone (MBK)	< 5.0	Y	5.0
4-Methyl-2-pentanone (MIBK)	< 10	Y	10
<b>Acetone</b>	<b>200</b>	L, Y	10
Benzene	< 2.0	Y	2.0
Bromobenzene	< 2.0	Y	2.0
Bromochloromethane	< 2.0	Y	2.0
Bromodichloromethane	< 2.0	Y	2.0
Bromoform	< 5.0	Y	5.0
Bromomethane	< 5.0	Y	5.0

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

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### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 19:09

<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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<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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<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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<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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Units: ug/L Analyzed: 06/29/21 19:09

<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: 06/29/21 16:08

Units: ug/L Analyzed: 07/01/21 10:33

<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: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 10:50

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Aluminum	10400		100
Antimony	< 10.0		10.0
Arsenic	< 10.0		10.0
Barium	164		10.0
Beryllium	< 1.00		1.00
Boron	242	B1	25.0
Cadmium	9.72		3.00

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

Method: SW-846 6010 Prepared: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 12:43

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Calcium	290000		3000
Chromium	24.9		5.00
Cobalt	27.0		10.0
Copper	116		10.0
Hardness	1240000		1980
Iron	29900		150
Lead	17.4		5.00
Magnesium	126000		300
Manganese	1040		15.0
Nickel	33.9		5.00
Potassium	5520		1400
Selenium	< 25.0	B1	25.0
Silver	5.05		3.00
Sodium	198000		300
Strontium	350		5.00
Thallium	< 10.0		10.0
Vanadium	31.8		5.00
Zinc	128		25.0

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### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G107** Lab Sample ID: **21F1058-02**

Matrix: Water Collected By: SP Date/Time Collected: 06/25/21 16:30

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

Method: 8260 Prepared: 06/27/21 10:50

Units: ug/L Analyzed: 06/27/21 18:32

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

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Carbon disulfide	< 2.0	Y	2.0
Carbon tetrachloride	< 2.0	Y	2.0
Chlorobenzene	< 2.0	Y	2.0
Chloroethane	< 2.0	Y	2.0
Chloroform	< 2.0	Y	2.0
Chloromethane	< 2.0	Y	2.0
cis-1,2-Dichloroethene	< 2.0	Y	2.0
cis-1,3-Dichloropropene	< 2.0	Y	2.0
Dibromochloromethane	< 5.0	Y	5.0
Dibromomethane	< 2.0	Y	2.0
Ethylbenzene	< 2.0	Y	2.0
Isopropylbenzene	< 2.0	Y	2.0
Methyl tert-butyl ether	< 2.0	Y	2.0
Methylene chloride	< 5.0	Y	5.0
Styrene	< 2.0	Y	2.0
Tetrachloroethene	< 2.0	Y	2.0
Toluene	< 2.0	Y	2.0
trans-1,2-Dichloroethene	< 2.0	Y	2.0
trans-1,3-Dichloropropene	< 5.0	Y	5.0
Trichloroethene	< 2.0	Y	2.0
Trichlorofluoromethane	< 2.0	Y	2.0
Vinyl chloride	< 2.0	Y	2.0
Xylenes, total	< 2.0	Y	2.0

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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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G107** Lab Sample ID: **21F1058-02**

Matrix: Water Collected By: SP Date/Time Collected: 06/25/21 16:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 19:43

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G107** Lab Sample ID: **21F1058-02**

Matrix: Water Collected By: SP Date/Time Collected: 06/25/21 16:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 19:43

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G107** Lab Sample ID: **21F1058-02**

Matrix: Water Collected By: SP Date/Time Collected: 06/25/21 16:30

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 19:43

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G107** Lab Sample ID: **21F1058-02**

Matrix: Water Collected By: SP Date/Time Collected: 06/25/21 16:30

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

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 19:43

<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: 06/29/21 16:08

Units: ug/L Analyzed: 07/01/21 10:35

<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: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 10:54

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
<b>Aluminum</b>	<b>4870</b>		100
Antimony	< 10.0		10.0
Arsenic	< 10.0		10.0
<b>Barium</b>	<b>70.1</b>		10.0
Beryllium	< 1.00		1.00
<b>Boron</b>	<b>54.7</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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G107** Lab Sample ID: **21F1058-02**

Matrix: Water Collected By: SP Date/Time Collected: 06/25/21 16:30

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

Method: SW-846 6010 Prepared: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 10:54

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
<b>Calcium</b>	<b>154000</b>		300
<b>Chromium</b>	<b>13.0</b>		5.00
Cobalt	< 10.0		10.0
<b>Copper</b>	<b>15.9</b>		10.0
<b>Hardness</b>	<b>631000</b>		1980
<b>Iron</b>	<b>8810</b>		150
Lead	< 5.00		5.00
<b>Magnesium</b>	<b>59800</b>		300
<b>Manganese</b>	<b>312</b>		15.0
<b>Nickel</b>	<b>10.9</b>		5.00
<b>Potassium</b>	<b>1960</b>		1400
Selenium	< 25.0	B1	25.0
<b>Silver</b>	<b>3.26</b>		3.00
<b>Sodium</b>	<b>11400</b>		300
<b>Strontium</b>	<b>90.0</b>		5.00
Thallium	< 10.0		10.0
<b>Vanadium</b>	<b>10.9</b>		5.00
<b>Zinc</b>	<b>28.4</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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G108** Lab Sample ID: **21F1058-03**

Matrix: Water Collected By: LC Date/Time Collected: 06/25/21 17:50

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

Method: 8260 Prepared: 06/27/21 10:50

Units: ug/L Analyzed: 06/27/21 18:53

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
1,1,1,2-Tetrachloroethane	< 2.0	Y	2.0
1,1,1-Trichloroethane	< 2.0	Y	2.0
1,1,2,2-Tetrachloroethane	< 2.0	Y	2.0
1,1,2-Trichloroethane	< 2.0	Y	2.0
1,1-Dichloroethane	< 2.0	Y	2.0
1,1-Dichloroethene	< 2.0	Y	2.0
1,1-Dichloropropene	< 2.0	Y	2.0
1,2,3-Trichloropropane	< 2.0	Y	2.0
1,2-Dibromoethane	< 2.0	Y	2.0
1,2-Dichloroethane	< 2.0	Y	2.0
1,2-Dichloropropane	< 2.0	Y	2.0
1,3-Dichloropropane	< 2.0	Y	2.0
2,2-Dichloropropane	< 2.0	Y	2.0
2-Butanone (MEK)	< 10	Y	10
2-Hexanone (MBK)	< 5.0	Y	5.0
4-Methyl-2-pentanone (MIBK)	< 10	Y	10
Acetone	< 10	Y	10
Benzene	< 2.0	Y	2.0
Bromobenzene	< 2.0	Y	2.0
Bromochloromethane	< 2.0	Y	2.0
Bromodichloromethane	< 2.0	Y	2.0
Bromoform	< 5.0	Y	5.0
Bromomethane	< 5.0	Y	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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G108** Lab Sample ID: **21F1058-03**

Matrix: Water Collected By: LC Date/Time Collected: 06/25/21 17:50

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

Method: 8260 Prepared: 06/27/21 10:50

Units: ug/L Analyzed: 06/27/21 18:53

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Carbon disulfide	< 2.0	Y	2.0
Carbon tetrachloride	< 2.0	Y	2.0
Chlorobenzene	< 2.0	Y	2.0
Chloroethane	< 2.0	Y	2.0
Chloroform	< 2.0	Y	2.0
Chloromethane	< 2.0	Y	2.0
cis-1,2-Dichloroethene	< 2.0	Y	2.0
cis-1,3-Dichloropropene	< 2.0	Y	2.0
Dibromochloromethane	< 5.0	Y	5.0
Dibromomethane	< 2.0	Y	2.0
Ethylbenzene	< 2.0	Y	2.0
Isopropylbenzene	< 2.0	Y	2.0
Methyl tert-butyl ether	< 2.0	Y	2.0
Methylene chloride	< 5.0	Y	5.0
Styrene	< 2.0	Y	2.0
<b>Tetrachloroethene</b>	<b>6.9</b>	Y	2.0
Toluene	< 2.0	Y	2.0
trans-1,2-Dichloroethene	< 2.0	Y	2.0
trans-1,3-Dichloropropene	< 5.0	Y	5.0
Trichloroethene	< 2.0	Y	2.0
Trichlorofluoromethane	< 2.0	Y	2.0
Vinyl chloride	< 2.0	Y	2.0
Xylenes, total	< 2.0	Y	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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G108** Lab Sample ID: **21F1058-03**

Matrix: Water Collected By: LC Date/Time Collected: 06/25/21 17:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:18

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G108** Lab Sample ID: **21F1058-03**

Matrix: Water Collected By: LC Date/Time Collected: 06/25/21 17:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:18

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G108** Lab Sample ID: **21F1058-03**

Matrix: Water Collected By: LC Date/Time Collected: 06/25/21 17:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:18

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G108** Lab Sample ID: **21F1058-03**

Matrix: Water Collected By: LC Date/Time Collected: 06/25/21 17:50

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:18

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G108** Lab Sample ID: **21F1058-03**

Matrix: Water Collected By: LC Date/Time Collected: 06/25/21 17:50

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

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:18

<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: 06/29/21 16:08

Units: ug/L Analyzed: 07/01/21 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: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 11:04

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
<b>Aluminum</b>	<b>41900</b>		100
Antimony	< 10.0		10.0
Arsenic	< 10.0		10.0
<b>Barium</b>	<b>348</b>		10.0
Beryllium	< 1.00		1.00
Boron	< 25.0	B1	25.0
<b>Cadmium</b>	<b>23.1</b>		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G108** Lab Sample ID: **21F1058-03**

Matrix: Water Collected By: LC Date/Time Collected: 06/25/21 17:50

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

Method: SW-846 6010 Prepared: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 12:46

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Calcium	1010000		3000
Chromium	344		5.00
Cobalt	62.8		10.0
Copper	161		10.0
Hardness	4310000		1980
Iron	84900		150
Lead	38.3		5.00
Magnesium	432000		3000
Manganese	2850		15.0
Nickel	184		5.00
Potassium	9690		1400
Selenium	< 25.0	B1	25.0
Silver	13.2		3.00
Sodium	54000		300
Strontium	474		5.00
Thallium	< 10.0		10.0
Vanadium	95.4		5.00
Zinc	271		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G109** Lab Sample ID: **21F1058-04**

Matrix: Water Collected By: SP Date/Time Collected: 06/26/21 9:35

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

Method: 8260 Prepared: 06/27/21 10:50

Units: ug/L Analyzed: 06/27/21 19:14

<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

825 N. Rutledge Springfield, Illinois 62702 217.782.9780

### LABORATORY RESULTS

Name: **CHEMTOOL**

Project/Facility Number: 2010355004 Date Received : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G109** Lab Sample ID: **21F1058-04**

Matrix: Water Collected By: SP Date/Time Collected: 06/26/21 9:35

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

Method: 8260 Prepared: 06/27/21 10:50

Units: ug/L Analyzed: 06/27/21 19:14

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G109** Lab Sample ID: **21F1058-04**

Matrix: Water Collected By: SP Date/Time Collected: 06/26/21 9:35

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:52

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
1,2,4,5-Tetrachlorobenzene	< 1.5		1.5
1,2,4-Trichlorobenzene	< 1.5		1.5
1,2-Dichlorobenzene	< 1.5		1.5
1,2-Dinitrobenzene	< 1.5		1.5
1,3-Dichlorobenzene	< 1.5		1.5
1,3-Dinitrobenzene	< 5.0		5.0
1,4-Dichlorobenzene	< 1.5		1.5
1,4-Dinitrobenzene	< 5.0		5.0
1-Chloronaphthalene	< 1.5		1.5
1-Naphthylamine	< 5.0		5.0
2,2-Oxybis(1-chloropropane)	< 1.5		1.5
2,3,4,6-Tetrachlorophenol	< 1.5		1.5
2,4,5-Trichlorophenol	< 1.5		1.5
2,4,6-Trichlorophenol	< 1.5		1.5
2,4-Dichlorophenol	< 1.5		1.5
2,4-Dimethylphenol	< 1.5		1.5
2,4-Dinitrophenol	< 5.0		5.0
2,4-Dinitrotoluene	< 5.0		5.0
2,6-Dichlorophenol	< 1.5		1.5
2,6-Dinitrotoluene	< 1.5		1.5
2-Chloronaphthalene	< 1.5		1.5
2-Chlorophenol	< 1.5		1.5
2-Methylnaphthalene	< 1.5		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G109** Lab Sample ID: **21F1058-04**

Matrix: Water Collected By: SP Date/Time Collected: 06/26/21 9:35

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:52

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
2-Methylphenol	< 1.5		1.5
2-Naphthylamine	< 5.0		5.0
2-Nitroaniline	< 1.5		1.5
2-Nitrophenol	< 5.0		5.0
2-Picoline	< 1.5		1.5
3,3-Dichlorobenzidine	< 1.5		1.5
3-Nitroaniline	< 1.5		1.5
4,6-Dinitro-2-methylphenol	< 5.0		5.0
4-Bromophenyl phenyl ether	< 1.5		1.5
4-Chloro-3-methylphenol	< 1.5		1.5
4-Chloroaniline	< 1.5		1.5
4-Chlorophenyl phenyl ether	< 1.5		1.5
4-Methylphenol	< 1.5		1.5
4-Nitroaniline	< 1.5		1.5
4-Nitrobiphenyl	< 5.0		5.0
4-Nitrophenol	< 5.0		5.0
5-Nitroacenaphthene	< 5.0		5.0
7,12-Dimethylbenzo(a)anthracene	< 5.0		5.0
Acenaphthene	< 1.5		1.5
Acenaphthylene	< 1.5		1.5
Acetophenone	< 1.5		1.5
Anthracene	< 1.5		1.5
Azobenzene	< 1.5		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G109** Lab Sample ID: **21F1058-04**

Matrix: Water Collected By: SP Date/Time Collected: 06/26/21 9:35

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:52

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Benzo(a)anthracene	< 1.5		1.5
Benzo(a)pyrene	< 1.5		1.5
Benzo(b)fluoranthene	< 1.5		1.5
Benzo(ghi)perylene	< 5.0		5.0
Benzo(k)fluoranthene	< 1.5		1.5
Bis(2-chloroethoxy)methane	< 1.5		1.5
Bis(2-chloroethyl)ether	< 1.5		1.5
Bis(2-ethylhexyl)phthalate	< 5.0		5.0
Butyl benzyl phthalate	< 5.0		5.0
Carbazole	< 1.5		1.5
Chrysene	< 1.5		1.5
Dibenzo(a,h)anthracene	< 5.0		5.0
Dibenzofuran	< 1.5		1.5
Diethylphthalate	< 1.5		1.5
Dimethylphthalate	< 1.5		1.5
Di-n-butylphthalate	< 1.5		1.5
Di-n-octylphthalate	< 5.0		5.0
Diphenylamine	< 1.5		1.5
Ethyl methanesulfonate	< 1.5		1.5
Fluoranthene	< 1.5		1.5
Fluorene	< 1.5		1.5
Hexachlorobenzene	< 1.5		1.5
Hexachlorobutadiene	< 1.5		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G109** Lab Sample ID: **21F1058-04**

Matrix: Water Collected By: SP Date/Time Collected: 06/26/21 9:35

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:52

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Hexachlorocyclopentadiene	< 1.5		1.5
Hexachloroethane	< 1.5		1.5
Hexachloropropene	< 1.5		1.5
Indeno(1,2,3-cd)pyrene	< 5.0		5.0
Isodrin	< 1.5		1.5
Isophorone	< 1.5		1.5
Isosafrole	< 1.5		1.5
Mestranol	< 5.0		5.0
Methyl methanesulfonate	< 1.5		1.5
Naphthalene	< 1.5		1.5
Nitrobenzene	< 1.5		1.5
N-Nitrosodi-n-butylamine	< 1.5		1.5
N-Nitrosodi-n-propylamine	< 1.5		1.5
N-Nitrosopiperidine	< 1.5		1.5
p-Dimethylaminoazobenzene	< 1.5		1.5
Pentachlorobenzene	< 1.5		1.5
Pentachloronitrobenzene	< 1.5		1.5
Pentachlorophenol	< 5.0		5.0
Phenacetin	< 1.5		1.5
Phenanthrene	< 1.5		1.5
Phenol	< 1.5		1.5
Pronamide	< 1.5		1.5
Pyrene	< 1.5		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G109** Lab Sample ID: **21F1058-04**

Matrix: Water Collected By: SP Date/Time Collected: 06/26/21 9:35

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

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 20:52

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

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

Method: 245.1 Prepared: 06/29/21 16:08

Units: ug/L Analyzed: 07/01/21 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: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 11:08

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
<b>Aluminum</b>	<b>573</b>		100
Antimony	< 10.0		10.0
Arsenic	< 10.0		10.0
<b>Barium</b>	<b>50.3</b>		10.0
Beryllium	< 1.00		1.00
<b>Boron</b>	<b>36.3</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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G109** Lab Sample ID: **21F1058-04**

Matrix: Water Collected By: SP Date/Time Collected: 06/26/21 9:35

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

Method: SW-846 6010 Prepared: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 11:08

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
<b>Calcium</b>	<b>107000</b>		300
<b>Chromium</b>	<b>13.5</b>		5.00
Cobalt	< 10.0		10.0
<b>Copper</b>	<b>36.0</b>		10.0
<b>Hardness</b>	<b>474000</b>		1980
<b>Iron</b>	<b>8640</b>		150
Lead	< 5.00		5.00
<b>Magnesium</b>	<b>50200</b>		300
<b>Manganese</b>	<b>51.3</b>		15.0
<b>Nickel</b>	<b>7.93</b>		5.00
Potassium	< 1400		1400
Selenium	< 25.0	B1	25.0
Silver	< 3.00		3.00
<b>Sodium</b>	<b>8050</b>		300
<b>Strontium</b>	<b>73.5</b>		5.00
Thallium	< 10.0		10.0
<b>Vanadium</b>	<b>6.65</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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G110** Lab Sample ID: **21F1058-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/26/21 10:45

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

Method: 8260 Prepared: 06/27/21 08:00

Units: ug/L Analyzed: 06/27/21 19:26

<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	O1	10
Acetone	< 10	O2	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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G110** Lab Sample ID: **21F1058-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/26/21 10:45

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

Method: 8260 Prepared: 06/27/21 08:00

Units: ug/L Analyzed: 06/27/21 19:26

<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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G110** Lab Sample ID: **21F1058-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/26/21 10:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 21:27

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
1,2,4,5-Tetrachlorobenzene	< 1.5		1.5
1,2,4-Trichlorobenzene	< 1.5		1.5
1,2-Dichlorobenzene	< 1.5		1.5
1,2-Dinitrobenzene	< 1.5		1.5
1,3-Dichlorobenzene	< 1.5		1.5
1,3-Dinitrobenzene	< 5.0		5.0
1,4-Dichlorobenzene	< 1.5		1.5
1,4-Dinitrobenzene	< 5.0		5.0
1-Chloronaphthalene	< 1.5		1.5
1-Naphthylamine	< 5.0		5.0
2,2-Oxybis(1-chloropropane)	< 1.5		1.5
2,3,4,6-Tetrachlorophenol	< 1.5		1.5
2,4,5-Trichlorophenol	< 1.5		1.5
2,4,6-Trichlorophenol	< 1.5		1.5
2,4-Dichlorophenol	< 1.5		1.5
2,4-Dimethylphenol	< 1.5		1.5
2,4-Dinitrophenol	< 5.0		5.0
2,4-Dinitrotoluene	< 5.0		5.0
2,6-Dichlorophenol	< 1.5		1.5
2,6-Dinitrotoluene	< 1.5		1.5
2-Chloronaphthalene	< 1.5		1.5
2-Chlorophenol	< 1.5		1.5
2-Methylnaphthalene	< 1.5		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G110** Lab Sample ID: **21F1058-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/26/21 10:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 21:27

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
2-Methylphenol	< 1.5		1.5
2-Naphthylamine	< 5.0		5.0
2-Nitroaniline	< 1.5		1.5
2-Nitrophenol	< 5.0		5.0
2-Picoline	< 1.5		1.5
3,3-Dichlorobenzidine	< 1.5		1.5
3-Nitroaniline	< 1.5		1.5
4,6-Dinitro-2-methylphenol	< 5.0		5.0
4-Bromophenyl phenyl ether	< 1.5		1.5
4-Chloro-3-methylphenol	< 1.5		1.5
4-Chloroaniline	< 1.5		1.5
4-Chlorophenyl phenyl ether	< 1.5		1.5
4-Methylphenol	< 1.5		1.5
4-Nitroaniline	< 1.5		1.5
4-Nitrobiphenyl	< 5.0		5.0
4-Nitrophenol	< 5.0		5.0
5-Nitroacenaphthene	< 5.0		5.0
7,12-Dimethylbenzo(a)anthracene	< 5.0		5.0
Acenaphthene	< 1.5		1.5
Acenaphthylene	< 1.5		1.5
Acetophenone	< 1.5		1.5
Anthracene	< 1.5		1.5
Azobenzene	< 1.5		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G110** Lab Sample ID: **21F1058-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/26/21 10:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 21:27

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Benzo(a)anthracene	< 1.5		1.5
Benzo(a)pyrene	< 1.5		1.5
Benzo(b)fluoranthene	< 1.5		1.5
Benzo(ghi)perylene	< 5.0		5.0
Benzo(k)fluoranthene	< 1.5		1.5
Bis(2-chloroethoxy)methane	< 1.5		1.5
Bis(2-chloroethyl)ether	< 1.5		1.5
Bis(2-ethylhexyl)phthalate	< 5.0		5.0
Butyl benzyl phthalate	< 5.0		5.0
Carbazole	< 1.5		1.5
Chrysene	< 1.5		1.5
Dibenzo(a,h)anthracene	< 5.0		5.0
Dibenzofuran	< 1.5		1.5
Diethylphthalate	< 1.5		1.5
Dimethylphthalate	< 1.5		1.5
Di-n-butylphthalate	< 1.5		1.5
Di-n-octylphthalate	< 5.0		5.0
Diphenylamine	< 1.5		1.5
Ethyl methanesulfonate	< 1.5		1.5
Fluoranthene	< 1.5		1.5
Fluorene	< 1.5		1.5
Hexachlorobenzene	< 1.5		1.5
Hexachlorobutadiene	< 1.5		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G110** Lab Sample ID: **21F1058-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/26/21 10:45

### Semivolatiles by GC/MS

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 21:27

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Hexachlorocyclopentadiene	< 1.5		1.5
Hexachloroethane	< 1.5		1.5
Hexachloropropene	< 1.5		1.5
Indeno(1,2,3-cd)pyrene	< 5.0		5.0
Isodrin	< 1.5		1.5
Isophorone	< 1.5		1.5
Isosafrole	< 1.5		1.5
Mestranol	< 5.0		5.0
Methyl methanesulfonate	< 1.5		1.5
Naphthalene	< 1.5		1.5
Nitrobenzene	< 1.5		1.5
N-Nitrosodi-n-butylamine	< 1.5		1.5
N-Nitrosodi-n-propylamine	< 1.5		1.5
N-Nitrosopiperidine	< 1.5		1.5
p-Dimethylaminoazobenzene	< 1.5		1.5
Pentachlorobenzene	< 1.5		1.5
Pentachloronitrobenzene	< 1.5		1.5
Pentachlorophenol	< 5.0		5.0
Phenacetin	< 1.5		1.5
Phenanthrene	< 1.5		1.5
Phenol	< 1.5		1.5
Pronamide	< 1.5		1.5
Pyrene	< 1.5		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G110** Lab Sample ID: **21F1058-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/26/21 10:45

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

Method: 8270 Prepared: 06/28/21 12:48

Units: ug/L Analyzed: 06/29/21 21:27

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

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

Method: 245.1 Prepared: 06/29/21 16:08

Units: ug/L Analyzed: 07/01/21 10:47

<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: 06/30/21 15:59

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

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
<b>Aluminum</b>	<b>5590</b>		100
<b>Antimony</b>	<b>16.3</b>		10.0
Arsenic	< 10.0		10.0
<b>Barium</b>	<b>108</b>		10.0
Beryllium	< 1.00		1.00
Boron	< 25.0	B1	25.0
<b>Cadmium</b>	<b>8.09</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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

Client Sample ID: **G110** Lab Sample ID: **21F1058-05**

Matrix: Water Collected By: MW Date/Time Collected: 06/26/21 10:45

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

Method: SW-846 6010 Prepared: 06/30/21 15:59

Units: ug/L Analyzed: 07/02/21 12:50

<u>Analyte</u>	<u>Result</u>	<u>Qualifier</u>	<u>Reporting Limit</u>
Calcium	205000		3000
Chromium	1580		5.00
Cobalt	38.8		10.0
Copper	56.9		10.0
Hardness	889000		1980
Iron	30700		150
Lead	15.9		5.00
Magnesium	91900		300
Manganese	1660		15.0
Nickel	86.3		5.00
Potassium	2270		1400
Selenium	< 25.0	B1	25.0
Silver	5.89		3.00
Sodium	9310		300
Strontium	103		5.00
Thallium	< 10.0		10.0
Vanadium	44.5		5.00
Zinc	46.9		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 : 06/27/21

Funding Code: CS29 B50 Temperature C: 7.00

#### **Notes and Definitions**

- Y The laboratory analysis was performed on an unpreserved or improperly preserved sample.
- O2 Quality control sample failed low - possible low bias or false non-detect result.
- O1 Quality control sample failed high - possible high bias or false positive result.
- L Actual value not known, but known to be greater than value shown. Value shown is the highest acceptable level for quantitation. (For bacteria, result calculated as if the smallest filtration volume had a count of 200).
- 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

Methods 8260 & 8270: Samples received at the laboratory outside of the acceptable temperature requirements were Y qualified.

Method 8260: Dilution for acetone could not be analyzed for sample 21F1058-01. Concentration known to be greater than value reported.

Method 8270: Tentatively Identified Compounds (TICs) were detected in the semi-volatile analysis of sample 21F1058-01. Please contact the laboratory if additional information about the TICs is needed.

Method 8270: Insufficient sample volume received to perform a matrix spike and matrix spike duplicate for the batch containing samples 21F1058-01, -02, -03, -04 and -05. NELAC and method requirements were not met.

Report Authorized by:

Tom Weiss  
Laboratory Manager

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