

Patoka Lake Regional Water District

WATER QUALITY DATA 2021

Inorganic Contaminants(2021)

	MCL	D.L.	RESULT
	mg/L	mg/L	mg/L
Antimony	0.006	0.001	BDL
Arsenic	0.01	0.001	BDL
Barium	2	0.002	0.025
Beryllium	0.004	0.0003	BDL
Cadmium	0.005	0.001	BDL
Chromium	1	0.0009	BDL
Cyanide, (Free)	0.2	0.02	BDL
Fluoride	4	0.1	0.6
Mercury	0.002	0.0001	BDL
Nickel	0.1	0.001	BDL
Nitrate	10	0.1	BDL
Selenium	0.05	0.002	BDL
Sodium	No MCL	0.1	2.9
Thallium	0.002	0.0003	BDL

Definitions

"MCL"	means maximum contaminant level
"BDL"	means below detectable limit
"pCi/L"	means picocuries per liter
"D.L."	means detectable limit
2012 "mg/L"	means part per million or milligrams per liter
"NTU"	means nephelometric turbidity unit
"µg/L"	means part per billion or micrograms per liter
"U.C."	means unregulated contaminants

Radioactive Contaminants(2020)

	MCL	RESULT
		pCi/L
Radium-228 2020	.17+ .41	pCi/L
Gross Alpha 2020	15 1.7+ .9	pCi/L

Synthetic Organic Contaminants(2021)

	MCL	D.L.	RESULT
	ug/L	ug/L	ug/L
Alachlor(Lasso)	2021 2	0.1	BDL
Atrazine	2021 3	0.1	BDL
Benzo(a)pyrene	2021 0.2	0.02	BDL
Carbofuran	2021 40	0.9	BDL
Chlordane(alpha & gamma)	2021 2	0.1	BDL
2,4-D	2021 70	0.1	0.2
Dalapon	2021 200	1	BDL
DBCP	2021 0.2	0.01	BDL
Dinoseb	2021 7	0.1	BDL
2,3,7,8-TCDD(Dioxin)	2021 30 pg/L	5.0 pg/L	BDL
Diquat	2021 20	0.4	BDL
Di(2-ethylhexyl)adipate	2021 400	0.6	BDL
Di(2-ethylhexyl)phthalate	2021 6	0.6	BDL
Endothall	2021 100	9	BDL
Endrin	2021 2	0.01	BDL
Ethylene Dibromide(EDB)	2021 50 ng/L	10 ng/L	BDL
Glyphosate (Round-Up)	2019 700	6	BDL
Heptachlor	2021 0.4	0.04	BDL
Heptachlor Epoxide	2021 0.2	0.02	BDL
Hexachlorobenzene	2021 1	0.1	BDL
Hexachlorocyclopentadiene	2021 50	0.1	BDL
gamma-BHG Lindane	2021 0.2	0.02	BDL
Methoxychlor	2021 40	0.1	BDL
Oxamyl(Vydate)	2021 200	1	BDL
Pentachlorophenol	2021 1	0.04	BDL
Picloram(Tordon)	2021 500	0.1	BDL
PCBs	2019 0.5	0.5	BDL
Simazine	2021 4	0.07	BDL
2,4,5-TP(Silvex)	2021 50	0.1	BDL
Toxaphene	2021 3	1	BDL

Volatile Organic Contaminants(2021)

	MCL	D.L.	RESULT
	ug/L	ug/L	ug/L
Benzene	5	0.5	BDL
Carbon Tetrachloride	5	0.5	BDL
Chlorobenzene	100	0.5	BDL
1,2-Dichlorobenzene	600	0.5	BDL
1,4-Dichlorobenzene	75	0.5	BDL
1,2-Dichloroethane	5	0.5	BDL
1,1-Dichloroethylene	7	0.5	BDL
cis-1,2 Dichloroethylene	70	0.5	BDL
trans-1,2-Dichloroethylene	100	0.5	BDL
Dichloromethane	5	0.5	BDL
1,2-Dichloropropane	5	0.5	BDL
Ethylbenzene	700	0.5	BDL
Styrene	100	0.5	BDL
Tetrachloroethylene	5	0.5	BDL
Toluene	1000	0.5	BDL
1,2,4-Trichlorobenzene	70	0.5	BDL
1,1,1-Trichloroethane	200	0.5	BDL
1,1,2-Trichloroethane	5	0.5	BDL
Trichloroethylene	5	0.5	BDL
Vinyl Chloride	2	0.2	BDL
Total Xylenes	10000	0.5	BDL
Methy-T-butyl ether	NO MCL	0.5	BDL
TOTAL TRIHalomethanes(4)	80	0.5	41.7
Bromodichloromethane		0.5	4.9
Bromoform		0.5	BDL
Chlorodibromomethane		0.5	BDL
Chloroform		0.5	36.7
	MCL	RESULT	
	µg/L	µg/L	
Haloacetic Acids 5 (4)	60	34.9 Average	
	2021	Range	25
Total Trihalomethanes(4)	80	41.7 Average	45
	2021	Range	20.4
	MCL	RESULT	
Lead 90th percentile	2020 15ug/L		3.7ug/L
Copper 90th percentile	2020 1.3mg/L		0.17mg/L

Total Organic Carbon (TOC)

	MCL	Range	27.9% - 40.5%
	25%	Average	34%
Percent Removal TOC	Running	Average<25%	

Highest Turbidity Measurement 2021

.25 on 6/20/2021 & 8/24/2021

110 South Hill Street
 South Bend, IN 46617
 Tel: (574) 233-4777
 Fax: (574) 233-8207
 1 800 332 4345

Laboratory Report

Client: Patoka Lake Regional Water & Sewer District
 Attn: Jerry Allstott
 2647 North State Road 545
 Dubois, IN 47527

Report: 525439
 Priority: Standard Written
 Status: Final
 PWS ID: IN5219012

Sample Information					
EEA ID #	Client ID	Method	Collected Date / Time	Collected By:	Received Date / Time
4966010	EP	200.8	07/27/21 11:35	Client	07/28/21 08:30
4966010	EP	245.1	07/27/21 11:35	Client	07/28/21 08:30
4966010	EP	200.7	07/27/21 11:35	Client	07/28/21 08:30
4966011	EP	353.2	07/27/21 11:36	Client	07/28/21 08:30
4966011	EP	4500-F- C	07/27/21 11:36	Client	07/28/21 08:30
4966012	EP	335.4	07/27/21 11:37	Client	07/28/21 08:30
4966013	EP	524.2	07/27/21 11:30	Client	07/28/21 08:30

Report Summary

Note: This data was submitted electronically to IDEM for compliance.

Detailed quantitative results are presented on the following pages. The results presented relate only to the samples provided for analysis.

We appreciate the opportunity to provide you with this analysis. If you have any questions concerning this report, please do not hesitate to call Traci Chlebowski at (574) 233-4777.

Note: This report may not be reproduced, except in full, without written approval from EEA.

 ASM

 Authorized Signature Title
 Client Name: Patoka Lake Regional Water & Sewer District
 Report #: 525439

08/05/2021

 Date

Sampling Point: EP

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
57-12-5	Cyanide, Total	335.4	0.2 *	0.02	< 0.02	mg/L	08/02/21 12:39	08/02/21 13:47	4966012
14797-55-8	Nitrate	353.2	10 *	0.1	< 0.1	mg/L	---	07/28/21 15:45	4966011
16984-48-8	Fluoride	4500-F- C	4 *	0.1	0.6	mg/L	---	08/02/21 09:53	4966011

Metals									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
7440-23-5	Sodium \$	200.7	---	0.1	2.9	mg/L	---	07/29/21 13:20	4966010
7440-36-0	Antimony	200.8	6 *	1.0	< 1.0	ug/L	---	08/05/21 14:09	4966010
7440-38-2	Arsenic	200.8	10 *	1.0	< 1.0	ug/L	---	08/05/21 14:09	4966010
7440-39-3	Barium	200.8	2000 *	2.0	25	ug/L	---	08/05/21 14:09	4966010
7440-41-7	Beryllium	200.8	4 *	0.3	< 0.3	ug/L	---	08/05/21 14:09	4966010
7440-43-9	Cadmium	200.8	5 *	1.0	< 1.0	ug/L	---	08/05/21 14:09	4966010
7440-47-3	Chromium	200.8	100 *	0.9	< 0.9	ug/L	---	08/05/21 14:09	4966010
7440-02-0	Nickel	200.8	100 *	1.0	< 1.0	ug/L	---	08/05/21 14:09	4966010
7782-49-2	Selenium	200.8	50 *	2.0	< 2.0	ug/L	---	08/05/21 14:09	4966010
7440-28-0	Thallium	200.8	2 *	0.3	< 0.3	ug/L	---	08/05/21 14:09	4966010
7439-97-6	Mercury	245.1	2 *	0.1	< 0.1	ug/L	07/29/21 16:20	07/30/21 13:41	4966010

Volatile Organic Chemicals									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
71-43-2	Benzene	524.2	5 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
56-23-5	Carbon tetrachloride	524.2	5 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
108-90-7	Chlorobenzene	524.2	100 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
95-50-1	1,2-Dichlorobenzene	524.2	600 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
106-46-7	1,4-Dichlorobenzene	524.2	75 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
107-06-2	1,2-Dichloroethane	524.2	5 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
75-35-4	1,1-Dichloroethylene	524.2	7 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
156-59-2	cis-1,2-Dichloroethylene	524.2	70 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
156-60-5	trans-1,2-Dichloroethylene	524.2	100 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
75-09-2	Dichloromethane	524.2	5 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
78-87-5	1,2-Dichloropropane	524.2	5 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
100-41-4	Ethylbenzene	524.2	700 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
100-42-5	Styrene	524.2	100 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
127-18-4	Tetrachloroethylene	524.2	5 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
108-88-3	Toluene	524.2	1000 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
120-82-1	1,2,4-Trichlorobenzene	524.2	70 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
71-55-6	1,1,1-Trichloroethane	524.2	200 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
79-00-5	1,1,2-Trichloroethane	524.2	5 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
79-01-6	Trichloroethylene	524.2	5 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
75-01-4	Vinyl chloride	524.2	2 *	0.2	< 0.2	ug/L	---	07/30/21 11:35	4966013
95-47-6	1,2-Xylene	524.2	---	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
179601-23-1	1,3 + 1,4-Xylene	524.2	---	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013
1330-20-7	Xylenes, Total	524.2	10000 *	0.5	< 0.5	ug/L	---	07/30/21 11:35	4966013

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: EP001

PWS ID: IN5219012

Semi-volatile Organic Chemicals									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed	EEA ID #
15972-60-8	Alachlor	525.2	2 *	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
309-00-2	Aldrin \$	525.2	---	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
1912-24-9	Atrazine	525.2	3 *	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
50-32-8	Benzo(a)pyrene	525.2	0.2 *	0.02	< 0.02	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
58-89-9	gamma-BHC (Lindane)	525.2	0.2 *	0.02	< 0.02	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
23184-66-9	Butachlor \$	525.2	---	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
60-57-1	Dieldrin \$	525.2	---	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
103-23-1	Di(2-ethylhexyl)adipate	525.2	400 *	0.6	< 0.6	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
117-81-7	Di(2-ethylhexyl)phthalate	525.2	6 *	0.6	< 0.6	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
72-20-8	Endrin	525.2	2 *	0.01	< 0.01	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
76-44-8	Heptachlor	525.2	0.4 *	0.04	< 0.04	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
1024-57-3	Heptachlor epoxide	525.2	0.2 *	0.02	< 0.02	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
118-74-1	Hexachlorobenzene	525.2	1 *	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
77-47-4	Hexachlorocyclopentadiene	525.2	50 *	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
72-43-5	Methoxychlor	525.2	40 *	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
51218-45-2	Metolachlor \$	525.2	---	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
21087-64-9	Metribuzin \$	525.2	---	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
1918-16-7	Propachlor \$	525.2	---	0.1	< 0.1	ug/L	02/16/21 09:00	02/17/21 00:29	4832601
122-34-9	Simazine	525.2	4 *	0.07	< 0.07	ug/L	02/16/21 09:00	02/17/21 00:29	4832601

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: ETDS

PWS ID: IN5219012

Semi-volatile Organic Chemicals									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed	EEA ID #
15972-60-8	Alachlor	525.2	2 *	0.1	< 0.1	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
1912-24-9	Atrazine	525.2	3 *	0.1	< 0.1	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
50-32-8	Benzo(a)pyrene	525.2	0.2 *	0.02	< 0.02	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
58-89-9	gamma-BHC (Lindane)	525.2	0.2 *	0.02	< 0.02	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
103-23-1	Di(2-ethylhexyl)adipate	525.2	400 *	0.6	< 0.6	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
117-81-7	Di(2-ethylhexyl)phthalate	525.2	6 *	0.6	< 0.6	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
72-20-8	Endrin	525.2	2 *	0.01	< 0.01	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
76-44-8	Heptachlor	525.2	0.4 *	0.04	< 0.04	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
1024-57-3	Heptachlor epoxide	525.2	0.2 *	0.02	< 0.02	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
118-74-1	Hexachlorobenzene	525.2	1 *	0.1	< 0.1	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
77-47-4	Hexachlorocyclopentadiene	525.2	50 *	0.1	< 0.1	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
72-43-5	Methoxychlor	525.2	40 *	0.1	< 0.1	ug/L	05/19/21 08:02	05/20/21 07:13	4898972
122-34-9	Simazine	525.2	4 *	0.07	< 0.07	ug/L	05/19/21 08:02	05/20/21 07:13	4898972

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: EP

PWS ID: IN5219012

Semi-volatile Organic Chemicals									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed	EEA ID #
15972-60-8	Alachlor	525.2	2 *	0.1	< 0.1	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
1912-24-9	Atrazine	525.2	3 *	0.1	< 0.1	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
50-32-8	Benzo(a)pyrene	525.2	0.2 *	0.02	< 0.02	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
58-89-9	gamma-BHC (Lindane)	525.2	0.2 *	0.02	< 0.02	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
103-23-1	Di(2-ethylhexyl)adipate	525.2	400 *	0.6	< 0.6	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
117-81-7	Di(2-ethylhexyl)phthalate	525.2	6 *	0.6	< 0.6	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
72-20-8	Endrin	525.2	2 *	0.01	< 0.01	ug/L	08/12/21 07:49	08/20/21 14:51	4980921
76-44-8	Heptachlor	525.2	0.4 *	0.04	< 0.04	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
1024-57-3	Heptachlor epoxide	525.2	0.2 *	0.02	< 0.02	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
118-74-1	Hexachlorobenzene	525.2	1 *	0.1	< 0.1	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
77-47-4	Hexachlorocyclopentadiene	525.2	50 *	0.1	< 0.1	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
72-43-5	Methoxychlor	525.2	40 *	0.1	< 0.1	ug/L	08/12/21 07:49	08/14/21 11:39	4980921
122-34-9	Simazine	525.2	4 *	0.07	< 0.07	ug/L	08/12/21 07:49	08/14/21 11:39	4980921

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-7353-1

Client Sample ID: EP

Lab Sample ID: 810-7353-1

Date Collected: 11/10/21 11:00

Matrix: Drinking Water

Date Received: 11/11/21 08:15

PWSID Number: IN5219012

Method: 525.2 - Semivolatile Organic Compounds (GC/MS)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Heptachlor epoxide	<0.019		0.019	ug/L		11/12/21 08:43	11/17/21 14:29	1
Di(2-ethylhexyl)adipate	<0.57		0.57	ug/L		11/12/21 08:43	11/17/21 14:29	1
Di (2-ethylhexyl)phthalate	<0.57		0.57	ug/L		11/12/21 08:43	11/17/21 14:29	1
Hexachlorobenzene	<0.095		0.095	ug/L		11/12/21 08:43	11/17/21 14:29	1
Simazine	<0.067		0.067	ug/L		11/12/21 08:43	11/17/21 14:29	1
Alachlor	<0.095		0.095	ug/L		11/12/21 08:43	11/17/21 14:29	1
Atrazine	<0.095		0.095	ug/L		11/12/21 08:43	11/17/21 14:29	1
Benzo[a]pyrene	<0.019		0.019	ug/L		11/12/21 08:43	11/17/21 14:29	1
gamma-BHC (Lindane)	<0.019		0.019	ug/L		11/12/21 08:43	11/17/21 14:29	1
Endrin	<0.0095		0.0095	ug/L		11/12/21 08:43	11/17/21 14:29	1
Methoxychlor	<0.095		0.095	ug/L		11/12/21 08:43	11/17/21 14:29	1
Heptachlor	<0.038		0.038	ug/L		11/12/21 08:43	11/17/21 14:29	1
Hexachlorocyclopentadiene	<0.095		0.095	ug/L		11/12/21 08:43	11/17/21 14:29	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
2-Nitro-m-xylene	100		70 - 130	11/12/21 08:43	11/17/21 14:29	1
Perylene-d12	102		70 - 130	11/12/21 08:43	11/17/21 14:29	1
Triphenylphosphate	121		70 - 130	11/12/21 08:43	11/17/21 14:29	1

Method: 548.1 - Endothall (GC/MS)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Endothall	<9.0		9.0	ug/L		11/15/21 07:09	11/16/21 10:39	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
2,4-Dichlorophenoxyacetic acid (Surr)	97		70 - 130	11/15/21 07:09	11/16/21 10:39	1

Method: 504.1 - EDB, DBCP and 1,2,3-TCP (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
1,2-Dibromoethane	<0.010		0.010	ug/L		11/15/21 11:00	11/15/21 22:59	1
1,2-Dibromo-3-Chloropropane	<0.010		0.010	ug/L		11/15/21 11:00	11/15/21 22:59	1

Method: 505 - Organochlorine Pesticides/PCBs (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Chlordane (technical)	<0.10		0.10	ug/L		11/15/21 12:26	11/15/21 22:10	1
Toxaphene	<1.0		1.0	ug/L		11/15/21 12:26	11/15/21 22:10	1

Method: 515.3 - Herbicides (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
2,4,5-TP (Silvex)	<0.10		0.10	ug/L		11/15/21 08:06	11/18/21 04:18	1
Dalapon	<1.0		1.0	ug/L		11/15/21 08:06	11/18/21 04:18	1
Dinoseb	<0.10		0.10	ug/L		11/15/21 08:06	11/18/21 04:18	1
Pentachlorophenol	<0.040		0.040	ug/L		11/15/21 08:06	11/18/21 04:18	1
Picloram	<0.10		0.10	ug/L		11/15/21 08:06	11/18/21 04:18	1
2,4-D	<0.10		0.10	ug/L		11/15/21 08:06	11/18/21 04:18	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
2,4-Dichlorophenylacetic acid	120		70 - 130	11/15/21 08:06	11/18/21 04:18	1

Method: 531.2 - Carbamate Pesticides (HPLC) - Dissolved

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Carbofuran	<0.90		0.90	ug/L			11/16/21 17:26	1

Eurofins Eaton Analytical - South Bend

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-7353-1

Client Sample ID: EP

Date Collected: 11/10/21 11:00

Date Received: 11/11/21 08:15

Lab Sample ID: 810-7353-1

Matrix: Drinking Water

PWSID Number: IN5219012

Method: 531.2 - Carbamate Pesticides (HPLC) - Dissolved (Continued)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Oxamyl	<1.0		1.0	ug/L			11/16/21 17:26	1

Method: 549.2 - Diquat and Paraquat (HPLC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Diquat	<0.40		0.40	ug/L		11/15/21 06:09	11/22/21 15:41	1

Sampling Point: IN5938

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.71	mg/L	---	01/22/21 02:53	4813345

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP001

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	1.87	mg/L	---	01/22/21 03:13	4813346

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP002

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	1.88	mg/L	---	01/22/21 03:32	4813347

\$ The state of origin does not offer certification for this parameter.

Sampling Point: ETDS

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	1.94	mg/L	---	01/22/21 04:12	4813348

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: IN5938 Raw

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.86	mg/L	---	03/01/21 17:53	4842238

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP001

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	1.99	mg/L	---	03/01/21 18:13	4842239

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP002

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.00	mg/L	---	03/01/21 18:33	4842240

\$ The state of origin does not offer certification for this parameter.

Sampling Point: ETDS

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.04	mg/L	---	03/01/21 18:52	4842241

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: IN 5938

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	3.23	mg/L	---	03/30/21 17:12	4861678

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP 001

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.33	mg/L	---	03/30/21 17:31	4861679

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP 002

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.05	mg/L	---	03/30/21 17:51	4861680

\$ The state of origin does not offer certification for this parameter.

Sampling Point: ETDS

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.17	mg/L	---	03/30/21 18:11	4861681

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Client Name: Patoka Lake Regional Water & Sewer District

Report #: 514672

Sampling Point: IN 5938

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	3.41	mg/L	---	04/15/21 12:24	4871957

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP001

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.39	mg/L	---	04/15/21 12:44	4871958

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP002

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.45	mg/L	---	04/15/21 13:04	4871959

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: IN 5938

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	3.05	mg/L	---	05/13/21 08:09	4897829

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP001

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.21	mg/L	---	05/13/21 11:06	4897830

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP002

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.06	mg/L	---	05/13/21 11:26	4897831

\$ The state of origin does not offer certification for this parameter.

Sampling Point: ETDS

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.04	mg/L	---	05/13/21 11:46	4897832

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: ID5938

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	3.75	mg/L	---	06/11/21 19:34	4924053

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP001

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.37	mg/L	---	06/11/21 19:53	4924054

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP002

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.12	mg/L	---	06/11/21 20:13	4924055

\$ The state of origin does not offer certification for this parameter.

Sampling Point: EP

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.16	mg/L	---	06/11/21 20:33	4924056

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Client Name: Patoka Lake Regional Water & Sewer District

Report #: 524702

Sampling Point: ID5938

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	3.73	mg/L	---	07/28/21 17:46	4959800

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP001

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.54	mg/L	---	07/28/21 18:06	4959801

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP002

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.03	mg/L	---	07/28/21 19:44	4959802

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: ID5938

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	4.34	mg/L	---	08/18/21 18:06	4983272

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP001

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.59	mg/L	---	08/18/21 18:26	4983273

\$ The state of origin does not offer certification for this parameter.

Sampling Point: TP002

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.57	mg/L	---	08/18/21 18:45	4983274

\$ The state of origin does not offer certification for this parameter.

Sampling Point: EP

PWS ID: IN5219012

General Chemistry									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
---	Total Organic Carbon (TOC) \$	5310 C	---	0.500	2.43	mg/L	---	08/18/21 19:05	4983275

\$ The state of origin does not offer certification for this parameter.

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-1948-1

Client Sample ID: ID 5938

Date Collected: 09/13/21 06:00

Date Received: 09/14/21 09:45

Lab Sample ID: 810-1948-1

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	3.6		0.50	mg/L			09/17/21 02:42	1

Client Sample ID: TP001

Date Collected: 09/13/21 10:50

Date Received: 09/14/21 09:45

Lab Sample ID: 810-1948-2

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.4		0.50	mg/L			09/17/21 04:40	1

Client Sample ID: TP002

Date Collected: 09/13/21 11:30

Date Received: 09/14/21 09:45

Lab Sample ID: 810-1948-3

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.2		0.50	mg/L			09/17/21 05:00	1

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-4667-1

Client Sample ID: Raw TP001 & TP002

Date Collected: 10/11/21 05:45

Date Received: 10/12/21 08:30

Lab Sample ID: 810-4667-1

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	3.67		0.500	mg/L			10/19/21 03:07	1

Client Sample ID: TP001

Date Collected: 10/11/21 10:45

Date Received: 10/12/21 08:30

Lab Sample ID: 810-4667-2

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.58		0.500	mg/L			10/19/21 03:27	1

Client Sample ID: TP002

Date Collected: 10/11/21 11:00

Date Received: 10/12/21 08:30

Lab Sample ID: 810-4667-3

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.41		0.500	mg/L			10/19/21 07:46	1

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-7543-1

Client Sample ID: IN5938 TOC Raw TP01+TP02

Date Collected: 11/11/21 05:45

Date Received: 11/12/21 09:30

Lab Sample ID: 810-7543-1

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	3.37		0.500	mg/L			11/24/21 05:43	1

Client Sample ID: TP001 TOC SN-TP01

Date Collected: 11/11/21 10:00

Date Received: 11/12/21 09:30

Lab Sample ID: 810-7543-2

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.16		0.500	mg/L			11/24/21 06:02	1

Client Sample ID: TP002 TOC SN-TP02

Date Collected: 11/11/21 11:40

Date Received: 11/12/21 09:30

Lab Sample ID: 810-7543-3

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	1.97		0.500	mg/L			11/24/21 06:22	1

Client Sample ID: EP

Date Collected: 11/11/21 11:50

Date Received: 11/12/21 09:30

Lab Sample ID: 810-7543-4

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.01		0.500	mg/L			11/24/21 06:42	1

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-9653-1

Client Sample ID: ID 5938 TOC Raw TP1 & TP2

Date Collected: 12/09/21 06:00

Date Received: 12/10/21 08:00

Lab Sample ID: 810-9653-1

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.88		0.500	mg/L			12/16/21 06:39	1

Client Sample ID: TP001 TOC SN-TP01

Date Collected: 12/09/21 11:20

Date Received: 12/10/21 08:00

Lab Sample ID: 810-9653-2

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.12		0.500	mg/L			12/16/21 09:14	1

Client Sample ID: TP002 TOC SN-TP02

Date Collected: 12/09/21 11:45

Date Received: 12/10/21 08:00

Lab Sample ID: 810-9653-3

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.03		0.500	mg/L			12/16/21 09:34	1

Client Sample ID: EP-Non-Compliance

Date Collected: 12/09/21 11:45

Date Received: 12/10/21 08:00

Lab Sample ID: 810-9653-4

Matrix: Drinking Water

PWSID Number: IN5219012

General Chemistry

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Organic Carbon	2.09		0.500	mg/L			12/21/21 23:05	1

Sampling Point: Paoli

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	3.4	ug/L	---	02/13/21 01:45	4831372
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	02/13/21 01:45	4831372
67-66-3	Chloroform	524.2	---	0.5	17	ug/L	---	02/13/21 01:45	4831372
124-48-1	Dibromochloromethane	524.2	---	0.5	< 0.5	ug/L	---	02/13/21 01:45	4831372
---	Total Trihalomethanes	524.2	80 *	0.5	20.4	ug/L	---	02/13/21 01:45	4831372
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	02/12/21 07:41	02/12/21 23:48	4831373
79-43-6	Dichloroacetic acid	552.2	---	1.0	14	ug/L	02/12/21 07:41	02/12/21 23:48	4831373
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	02/12/21 07:41	02/12/21 23:48	4831373
79-11-8	Monochloroacetic acid	552.2	---	2.0	< 2.0	ug/L	02/12/21 07:41	02/12/21 23:48	4831373
76-03-9	Trichloroacetic acid	552.2	---	1.0	11	ug/L	02/12/21 07:41	02/12/21 23:48	4831373
---	Total HAA5	552.2	60 *	2.0	25	ug/L	02/12/21 07:41	02/12/21 23:48	4831373

Sampling Point: Finch Newton

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	3.5	ug/L	---	02/13/21 02:09	4831374
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	02/13/21 02:09	4831374
67-66-3	Chloroform	524.2	---	0.5	18	ug/L	---	02/13/21 02:09	4831374
124-48-1	Dibromochloromethane	524.2	---	0.5	< 0.5	ug/L	---	02/13/21 02:09	4831374
---	Total Trihalomethanes	524.2	80 *	0.5	21.5	ug/L	---	02/13/21 02:09	4831374
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	02/12/21 07:41	02/13/21 01:37	4831375
79-43-6	Dichloroacetic acid	552.2	---	1.0	14	ug/L	02/12/21 07:41	02/13/21 01:37	4831375
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	02/12/21 07:41	02/13/21 01:37	4831375
79-11-8	Monochloroacetic acid	552.2	---	2.0	< 2.0	ug/L	02/12/21 07:41	02/13/21 01:37	4831375
76-03-9	Trichloroacetic acid	552.2	---	1.0	11	ug/L	02/12/21 07:41	02/13/21 01:37	4831375
---	Total HAA5	552.2	60 *	2.0	25	ug/L	02/12/21 07:41	02/13/21 01:37	4831375

Sampling Point: Oakland City

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	3.5	ug/L	---	02/13/21 02:33	4831376
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	02/13/21 02:33	4831376
67-66-3	Chloroform	524.2	---	0.5	18	ug/L	---	02/13/21 02:33	4831376
124-48-1	Dibromochloromethane	524.2	---	0.5	< 0.5	ug/L	---	02/13/21 02:33	4831376
---	Total Trihalomethanes	524.2	80 *	0.5	21.5	ug/L	---	02/13/21 02:33	4831376
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	02/12/21 07:41	02/13/21 02:14	4831377
79-43-6	Dichloroacetic acid	552.2	---	1.0	14	ug/L	02/12/21 07:41	02/13/21 02:14	4831377
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	02/12/21 07:41	02/13/21 02:14	4831377
79-11-8	Monochloroacetic acid	552.2	---	2.0	2.8	ug/L	02/12/21 07:41	02/13/21 02:14	4831377
76-03-9	Trichloroacetic acid	552.2	---	1.0	12	ug/L	02/12/21 07:41	02/13/21 02:14	4831377
---	Total HAA5	552.2	60 *	2.0	28.8	ug/L	02/12/21 07:41	02/13/21 02:14	4831377

Sampling Point: Lynnville

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	3.4	ug/L	---	02/13/21 02:57	4831378
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	02/13/21 02:57	4831378
67-66-3	Chloroform	524.2	---	0.5	18	ug/L	---	02/13/21 02:57	4831378
124-48-1	Dibromochloromethane	524.2	---	0.5	< 0.5	ug/L	---	02/13/21 02:57	4831378
---	Total Trihalomethanes	524.2	80 *	0.5	21.4	ug/L	---	02/13/21 02:57	4831378
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	02/12/21 07:41	02/13/21 02:50	4831379
79-43-6	Dichloroacetic acid	552.2	---	1.0	16	ug/L	02/12/21 07:41	02/13/21 02:50	4831379
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	02/12/21 07:41	02/13/21 02:50	4831379
79-11-8	Monochloroacetic acid	552.2	---	2.0	4.5	ug/L	02/12/21 07:41	02/13/21 02:50	4831379
76-03-9	Trichloroacetic acid	552.2	---	1.0	13	ug/L	02/12/21 07:41	02/13/21 02:50	4831379
---	Total HAA5	552.2	60 *	2.0	33.5	ug/L	02/12/21 07:41	02/13/21 02:50	4831379

Sampling Point: Paoli

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	5.2	ug/L	---	05/06/21 19:04	4892899
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	05/06/21 19:04	4892899
67-66-3	Chloroform	524.2	---	0.5	42	ug/L	---	05/06/21 19:04	4892899
124-48-1	Dibromochloromethane	524.2	---	0.5	< 0.5	ug/L	---	05/06/21 19:04	4892899
---	Total Trihalomethanes	524.2	80 *	0.5	47.2	ug/L	---	05/06/21 19:04	4892899
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	05/11/21 07:34	05/12/21 19:15	4892900
79-43-6	Dichloroacetic acid	552.2	---	1.0	21	ug/L	05/11/21 07:34	05/12/21 19:15	4892900
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	05/11/21 07:34	05/12/21 19:15	4892900
79-11-8	Monochloroacetic acid	552.2	---	2.0	2.6	ug/L	05/11/21 07:34	05/12/21 19:15	4892900
76-03-9	Trichloroacetic acid	552.2	---	1.0	18	ug/L	05/11/21 07:34	05/12/21 19:15	4892900
---	Total HAA5	552.2	60 *	2.0	41.6	ug/L	05/11/21 07:34	05/12/21 19:15	4892900

Sampling Point: Finch Newton

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	5.4	ug/L	---	05/06/21 19:28	4892901
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	05/06/21 19:28	4892901
67-66-3	Chloroform	524.2	---	0.5	44	ug/L	---	05/06/21 19:28	4892901
124-48-1	Dibromochloromethane	524.2	---	0.5	< 0.5	ug/L	---	05/06/21 19:28	4892901
---	Total Trihalomethanes	524.2	80 *	0.5	49.4	ug/L	---	05/06/21 19:28	4892901
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	05/11/21 07:34	05/12/21 19:51	4892902
79-43-6	Dichloroacetic acid	552.2	---	1.0	21	ug/L	05/11/21 07:34	05/12/21 19:51	4892902
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	05/11/21 07:34	05/12/21 19:51	4892902
79-11-8	Monochloroacetic acid	552.2	---	2.0	2.4	ug/L	05/11/21 07:34	05/12/21 19:51	4892902
76-03-9	Trichloroacetic acid	552.2	---	1.0	19	ug/L	05/11/21 07:34	05/12/21 19:51	4892902
---	Total HAA5	552.2	60 *	2.0	42.4	ug/L	05/11/21 07:34	05/12/21 19:51	4892902

Sampling Point: Oakland City

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	5.3	ug/L	---	05/06/21 19:52	4892903
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	05/06/21 19:52	4892903
67-66-3	Chloroform	524.2	---	0.5	45	ug/L	---	05/06/21 19:52	4892903
124-48-1	Dibromochloromethane	524.2	---	0.5	< 0.5	ug/L	---	05/06/21 19:52	4892903
---	Total Trihalomethanes	524.2	80 *	0.5	50.3	ug/L	---	05/06/21 19:52	4892903
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	05/11/21 07:34	05/12/21 20:27	4892904
79-43-6	Dichloroacetic acid	552.2	---	1.0	20	ug/L	05/11/21 07:34	05/12/21 20:27	4892904
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	05/11/21 07:34	05/12/21 20:27	4892904
79-11-8	Monochloroacetic acid	552.2	---	2.0	3.6	ug/L	05/11/21 07:34	05/12/21 20:27	4892904
76-03-9	Trichloroacetic acid	552.2	---	1.0	19	ug/L	05/11/21 07:34	05/12/21 20:27	4892904
---	Total HAA5	552.2	60 *	2.0	42.6	ug/L	05/11/21 07:34	05/12/21 20:27	4892904

Sampling Point: Lynnville

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	5.2	ug/L	---	05/06/21 20:15	4892905
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	05/06/21 20:15	4892905
67-66-3	Chloroform	524.2	---	0.5	43	ug/L	---	05/06/21 20:15	4892905
124-48-1	Dibromochloromethane	524.2	---	0.5	< 0.5	ug/L	---	05/06/21 20:15	4892905
---	Total Trihalomethanes	524.2	80 *	0.5	48.2	ug/L	---	05/06/21 20:15	4892905
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	05/11/21 07:34	05/12/21 21:04	4892906
79-43-6	Dichloroacetic acid	552.2	---	1.0	21	ug/L	05/11/21 07:34	05/12/21 21:04	4892906
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	05/11/21 07:34	05/12/21 21:04	4892906
79-11-8	Monochloroacetic acid	552.2	---	2.0	2.3	ug/L	05/11/21 07:34	05/12/21 21:04	4892906
76-03-9	Trichloroacetic acid	552.2	---	1.0	19	ug/L	05/11/21 07:34	05/12/21 21:04	4892906
---	Total HAA5	552.2	60 *	2.0	42.3	ug/L	05/11/21 07:34	05/12/21 21:04	4892906

† EEA has demonstrated it can achieve these report limits in reagent water, but can not document them in all sample matrices.

Reg Limit Type:	MCL	SMCL	AL
Symbol:	*	^	!

Sampling Point: Paoli

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	5.5	ug/L	---	08/08/21 20:05	4972783
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	08/08/21 20:05	4972783
67-66-3	Chloroform	524.2	---	0.5	48	ug/L	---	08/08/21 20:05	4972783
124-48-1	Dibromochloromethane	524.2	---	0.5	0.5	ug/L	---	08/08/21 20:05	4972783
---	Total Trihalomethanes	524.2	80 *	0.5	54.0	ug/L	---	08/08/21 20:05	4972783
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	08/09/21 07:40	08/09/21 23:32	4972784
79-43-6	Dichloroacetic acid	552.2	---	1.0	15	ug/L	08/09/21 07:40	08/09/21 23:32	4972784
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	08/09/21 07:40	08/09/21 23:32	4972784
79-11-8	Monochloroacetic acid	552.2	---	2.0	< 2.0	ug/L	08/09/21 07:40	08/09/21 23:32	4972784
76-03-9	Trichloroacetic acid	552.2	---	1.0	17	ug/L	08/09/21 07:40	08/09/21 23:32	4972784
---	Total HAA5	552.2	60 *	2.0	32	ug/L	08/09/21 07:40	08/09/21 23:32	4972784

Sampling Point: Finch Newton

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	5.3	ug/L	---	08/08/21 20:28	4972785
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	08/08/21 20:28	4972785
67-66-3	Chloroform	524.2	---	0.5	46	ug/L	---	08/08/21 20:28	4972785
124-48-1	Dibromochloromethane	524.2	---	0.5	0.5	ug/L	---	08/08/21 20:28	4972785
---	Total Trihalomethanes	524.2	80 *	0.5	51.8	ug/L	---	08/08/21 20:28	4972785
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	08/09/21 07:40	08/10/21 00:08	4972786
79-43-6	Dichloroacetic acid	552.2	---	1.0	14	ug/L	08/09/21 07:40	08/10/21 00:08	4972786
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	08/09/21 07:40	08/10/21 00:08	4972786
79-11-8	Monochloroacetic acid	552.2	---	2.0	< 2.0	ug/L	08/09/21 07:40	08/10/21 00:08	4972786
76-03-9	Trichloroacetic acid	552.2	---	1.0	14	ug/L	08/09/21 07:40	08/10/21 00:08	4972786
---	Total HAA5	552.2	60 *	2.0	28	ug/L	08/09/21 07:40	08/10/21 00:08	4972786

Sampling Point: Oakland City

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	5.7	ug/L	---	08/08/21 20:52	4972787
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	08/08/21 20:52	4972787
67-66-3	Chloroform	524.2	---	0.5	52	ug/L	---	08/08/21 20:52	4972787
124-48-1	Dibromochloromethane	524.2	---	0.5	0.5	ug/L	---	08/08/21 20:52	4972787
---	Total Trihalomethanes	524.2	80 *	0.5	58.2	ug/L	---	08/08/21 20:52	4972787
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	08/09/21 07:40	08/10/21 00:44	4972788
79-43-6	Dichloroacetic acid	552.2	---	1.0	23	ug/L	08/09/21 07:40	08/10/21 00:44	4972788
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	08/09/21 07:40	08/10/21 00:44	4972788
79-11-8	Monochloroacetic acid	552.2	---	2.0	< 2.0	ug/L	08/09/21 07:40	08/10/21 00:44	4972788
76-03-9	Trichloroacetic acid	552.2	---	1.0	22	ug/L	08/09/21 07:40	08/10/21 00:44	4972788
---	Total HAA5	552.2	60 *	2.0	45	ug/L	08/09/21 07:40	08/10/21 00:44	4972788

Sampling Point: Lynnville

PWS ID: IN5219012

Disinfection Byproducts									
Analyte ID #	Analyte	Method	Reg Limit	MRL†	Result	Units	Preparation Date	Analyzed Date	EEA ID #
75-27-4	Bromodichloromethane	524.2	---	0.5	6.3	ug/L	---	08/08/21 21:15	4972789
75-25-2	Bromoform	524.2	---	0.5	< 0.5	ug/L	---	08/08/21 21:15	4972789
67-66-3	Chloroform	524.2	---	0.5	54	ug/L	---	08/08/21 21:15	4972789
124-48-1	Dibromochloromethane	524.2	---	0.5	0.6	ug/L	---	08/08/21 21:15	4972789
---	Total Trihalomethanes	524.2	80 *	0.5	60.9	ug/L	---	08/08/21 21:15	4972789
631-64-1	Dibromoacetic acid	552.2	---	1.0	< 1.0	ug/L	08/09/21 07:40	08/10/21 01:20	4972790
79-43-6	Dichloroacetic acid	552.2	---	1.0	22	ug/L	08/09/21 07:40	08/10/21 01:20	4972790
79-08-3	Monobromoacetic acid	552.2	---	1.0	< 1.0	ug/L	08/09/21 07:40	08/10/21 01:20	4972790
79-11-8	Monochloroacetic acid	552.2	---	2.0	< 2.0	ug/L	08/09/21 07:40	08/10/21 01:20	4972790
76-03-9	Trichloroacetic acid	552.2	---	1.0	20	ug/L	08/09/21 07:40	08/10/21 01:20	4972790
---	Total HAA5	552.2	60 *	2.0	42	ug/L	08/09/21 07:40	08/10/21 01:20	4972790

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-6769-1

Client Sample ID: Paoli

Lab Sample ID: 810-6769-1

Date Collected: 11/03/21 07:00

Matrix: Drinking Water

Date Received: 11/04/21 08:00

PWSID Number: IN5219012

Method: 524.2 - Total Trihalomethanes

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Trihalomethanes, Total	38.00		0.5000	ug/L			11/06/21 09:44	1

Method: 524.2 - Volatile Organic Compounds (GC/MS)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Bromodichloromethane	5.0		0.50	ug/L			11/06/21 09:44	1
Bromoform	<0.50		0.50	ug/L			11/06/21 09:44	1
Chloroform	33		0.50	ug/L			11/06/21 09:44	1
Dibromochloromethane	<0.50		0.50	ug/L			11/06/21 09:44	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	106		70 - 130		11/06/21 09:44	1
Toluene-d8 (Surr)	101		70 - 130		11/06/21 09:44	1
4-Bromofluorobenzene (Surr)	102		70 - 130		11/06/21 09:44	1
1,2-Dichlorobenzene-d4 (Surr)	106		70 - 130		11/06/21 09:44	1

Method: 552.2 THAA - Total Haloacetic Acids (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Haloacetic Acids 5	33.90		2.000	ug/L			11/09/21 08:30	1

Method: 552.2 - Haloacetic Acids (HAAs) (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Dibromoacetic acid	<1.0		1.0	ug/L		11/09/21 08:30	11/10/21 22:32	1
Dichloroacetic acid	16		1.0	ug/L		11/09/21 08:30	11/10/21 22:32	1
Monobromoacetic acid	<1.0		1.0	ug/L		11/09/21 08:30	11/10/21 22:32	1
Monochloroacetic acid	2.9		2.0	ug/L		11/09/21 08:30	11/10/21 22:32	1
Trichloroacetic acid	15		1.0	ug/L		11/09/21 08:30	11/10/21 22:32	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
2-Bromopropionic acid (Surr)	109		70 - 130	11/09/21 08:30	11/10/21 22:32	1

Client Sample ID: Lynville

Lab Sample ID: 810-6769-2

Date Collected: 11/03/21 08:00

Matrix: Drinking Water

Date Received: 11/04/21 08:00

PWSID Number: IN5219012

Method: 524.2 - Total Trihalomethanes

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Trihalomethanes, Total	42.80		0.5000	ug/L			11/06/21 10:08	1

Method: 524.2 - Volatile Organic Compounds (GC/MS)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Bromodichloromethane	4.8		0.50	ug/L			11/06/21 10:08	1
Bromoform	<0.50		0.50	ug/L			11/06/21 10:08	1
Chloroform	38		0.50	ug/L			11/06/21 10:08	1
Dibromochloromethane	<0.50		0.50	ug/L			11/06/21 10:08	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	102		70 - 130		11/06/21 10:08	1
Toluene-d8 (Surr)	96		70 - 130		11/06/21 10:08	1
4-Bromofluorobenzene (Surr)	96		70 - 130		11/06/21 10:08	1
1,2-Dichlorobenzene-d4 (Surr)	104		70 - 130		11/06/21 10:08	1

Eurofins Eaton Analytical - South Bend

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-6769-1

Client Sample ID: Lynville

Lab Sample ID: 810-6769-2

Date Collected: 11/03/21 08:00

Matrix: Drinking Water

Date Received: 11/04/21 08:00

PWSID Number: IN5219012

Method: 552.2 THAA - Total Haloacetic Acids (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Haloacetic Acids 5	36.00		2.000	ug/L			11/09/21 08:30	1

Method: 552.2 - Haloacetic Acids (HAAs) (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Dibromoacetic acid	<1.0		1.0	ug/L		11/09/21 08:30	11/10/21 23:09	1
Dichloroacetic acid	19		1.0	ug/L		11/09/21 08:30	11/10/21 23:09	1
Monobromoacetic acid	<1.0		1.0	ug/L		11/09/21 08:30	11/10/21 23:09	1
Monochloroacetic acid	<2.0		2.0	ug/L		11/09/21 08:30	11/10/21 23:09	1
Trichloroacetic acid	17		1.0	ug/L		11/09/21 08:30	11/10/21 23:09	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
2-Bromopropionic acid (Surr)	99		70 - 130	11/09/21 08:30	11/10/21 23:09	1

Client Sample ID: Finch-Newton

Lab Sample ID: 810-6769-3

Date Collected: 11/03/21 10:30

Matrix: Drinking Water

Date Received: 11/04/21 08:00

PWSID Number: IN5219012

Method: 524.2 - Total Trihalomethanes

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Trihalomethanes, Total	41.30		0.5000	ug/L			11/06/21 10:32	1

Method: 524.2 - Volatile Organic Compounds (GC/MS)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Bromodichloromethane	5.3		0.50	ug/L			11/06/21 10:32	1
Bromoform	<0.50		0.50	ug/L			11/06/21 10:32	1
Chloroform	36		0.50	ug/L			11/06/21 10:32	1
Dibromochloromethane	<0.50		0.50	ug/L			11/06/21 10:32	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	100		70 - 130		11/06/21 10:32	1
Toluene-d8 (Surr)	99		70 - 130		11/06/21 10:32	1
4-Bromofluorobenzene (Surr)	100		70 - 130		11/06/21 10:32	1
1,2-Dichlorobenzene-d4 (Surr)	101		70 - 130		11/06/21 10:32	1

Method: 552.2 THAA - Total Haloacetic Acids (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Haloacetic Acids 5	35.40		2.000	ug/L			11/09/21 08:30	1

Method: 552.2 - Haloacetic Acids (HAAs) (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Dibromoacetic acid	<1.0		1.0	ug/L		11/09/21 08:30	11/10/21 23:46	1
Dichloroacetic acid	17		1.0	ug/L		11/09/21 08:30	11/10/21 23:46	1
Monobromoacetic acid	<1.0		1.0	ug/L		11/09/21 08:30	11/10/21 23:46	1
Monochloroacetic acid	2.4		2.0	ug/L		11/09/21 08:30	11/10/21 23:46	1
Trichloroacetic acid	16		1.0	ug/L		11/09/21 08:30	11/10/21 23:46	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
2-Bromopropionic acid (Surr)	101		70 - 130	11/09/21 08:30	11/10/21 23:46	1

Eurofins Eaton Analytical - South Bend

Client Sample Results

Client: Patoka Lake Regional Water & Sewer District
Project/Site: Compliance

Job ID: 810-6769-1

Client Sample ID: Oakland City

Date Collected: 11/03/21 10:47

Date Received: 11/04/21 08:00

Lab Sample ID: 810-6769-4

Matrix: Drinking Water

PWSID Number: IN5219012

Method: 524.2 - Total Trihalomethanes

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Trihalomethanes, Total	40.10		0.5000	ug/L			11/06/21 10:56	1

Method: 524.2 - Volatile Organic Compounds (GC/MS)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Bromodichloromethane	5.1		0.50	ug/L			11/06/21 10:56	1
Bromoform	<0.50		0.50	ug/L			11/06/21 10:56	1
Chloroform	35		0.50	ug/L			11/06/21 10:56	1
Dibromochloromethane	<0.50		0.50	ug/L			11/06/21 10:56	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	98		70 - 130		11/06/21 10:56	1
Toluene-d8 (Surr)	98		70 - 130		11/06/21 10:56	1
4-Bromofluorobenzene (Surr)	97		70 - 130		11/06/21 10:56	1
1,2-Dichlorobenzene-d4 (Surr)	107		70 - 130		11/06/21 10:56	1

Method: 552.2 THAA - Total Haloacetic Acids (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Total Haloacetic Acids 5	25.00		2.000	ug/L			11/09/21 08:30	1

Method: 552.2 - Haloacetic Acids (HAAs) (GC)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Dibromoacetic acid	<1.0		1.0	ug/L		11/09/21 08:30	11/11/21 00:22	1
Dichloroacetic acid	13		1.0	ug/L		11/09/21 08:30	11/11/21 00:22	1
Monobromoacetic acid	<1.0		1.0	ug/L		11/09/21 08:30	11/11/21 00:22	1
Monochloroacetic acid	<2.0		2.0	ug/L		11/09/21 08:30	11/11/21 00:22	1
Trichloroacetic acid	12		1.0	ug/L		11/09/21 08:30	11/11/21 00:22	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
2-Bromopropionic acid (Surr)	98		70 - 130	11/09/21 08:30	11/11/21 00:22	1

Client Sample ID: EPDS

Date Collected: 11/03/21 11:30

Date Received: 11/04/21 08:00

Lab Sample ID: 810-6769-5

Matrix: Drinking Water

PWSID Number: IN5219012

Method: 524.2 - Total Trihalomethanes

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Trihalomethanes, Total	32.70		0.5000	ug/L			11/06/21 11:20	1

Method: 524.2 - Volatile Organic Compounds (GC/MS)

Analyte	Result	Qualifier	RL	Unit	D	Prepared	Analyzed	Dil Fac
Bromodichloromethane	4.7		0.50	ug/L			11/06/21 11:20	1
Bromoform	<0.50		0.50	ug/L			11/06/21 11:20	1
Chloroform	28		0.50	ug/L			11/06/21 11:20	1
Dibromochloromethane	<0.50		0.50	ug/L			11/06/21 11:20	1

Surrogate	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
1,2-Dichloroethane-d4 (Surr)	103		70 - 130		11/06/21 11:20	1
Toluene-d8 (Surr)	106		70 - 130		11/06/21 11:20	1
4-Bromofluorobenzene (Surr)	105		70 - 130		11/06/21 11:20	1
1,2-Dichlorobenzene-d4 (Surr)	113		70 - 130		11/06/21 11:20	1

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MONTHLY REPORTS OF OPERATION FOR # 1 WATER TREATMENT PLANT

Report for Patoka Lake Regional Water District
I.D.E.M. Field-Rep. Karla Goodman
Public Water Supply I.D. Number 5219012
Signed [Signature]

For the month of Dec 2021
Title Water Plant Superintendent
Certification No. WT058254

Date	Turbidity		Alkalinity		Ph		Hardness		Iron		Manganese		Aluminum	Temp. C
	Raw	Finish	Raw	Finish	Raw	Finish	Raw	Finish	Raw	Finish	Raw	Finish	Finish	Finish
1	3.8	.13	72	61	7.9	7.4	84	86	.015	.010	.099	.005	.028	11
2	4.5	.13	77	57	7.7	7.4	90	86	.065	.001	.179	.002	.021	13
3	4.4	.10	70	60	7.7	7.3	78	86	.080	.001	.171	.006	.021	11
4	3.5	.13	72	60	7.9	7.2	84	84	.044	.001	.132	.010	.039	11
5	4.4	.13	77	61	7.7	7.2	87	86	.059	.001	.183	.006	.023	13
6	5.5	.13	70	58	7.7	7.3	83	86	.114	.001	.241	.010	.023	14
7	4.1	.13	79	56	7.7	7.3	82	84	.049	.007	.148	.008	.021	12
8	4.3	.13	62	56	7.3	7.3	82	84	.058	.001	.111	.006	.026	11
9	4.2	.13	70	57	7.9	7.3	82	84	.048	.008	.148	.006	.028	10
10	5.7	.13	71	58	8.0	7.3	85	84	.099	.005	.225	.012	.022	10
11	4.5	.13	70	52	7.8	7.3	82	83	.008	.001	.130	.001	.021	12
12	2.6	.14	68	56	7.9	7.4	82	83	.040	.007	.080	.005	.023	13
13	3.3	.14	64	52	7.4	7.3	80	81	.048	.001	.127	.007	.023	12
14	2.5	.14	68	54	7.9	7.3	84	84	.037	.003	.112	.005	.028	10
15	3.8	.14	68	57	7.9	7.3	80	82	.036	.001	.134	.006	.028	10
16	3.9	.14	73	60	7.9	7.3	82	80	.068	.001	.141	.010	.016	14
17	4.7	.14	66	52	7.5	7.3	82	83	.070	.002	.132	.001	.017	11
18	4.7	.14	66	52	7.8/	7.2	84	82	.079	.010	.127	.006	.012	10
19	5.6	.14	65	51	7.8	7.2	82	84	.084	.005	.174	.013	.017	10
20	3.7	.14	63	51	7.9	7.3	84	84	.036	.002	.130	.014	.015	10
21	3.1	.12	62	56	7.8	7.3	82	83	.050	.010	.122	.017	.018	15
22	3.4	.12	68	52	7.8	7.3	82	82	.057	.001	.120	.009	.016	11
23	5.2	.12	71	52	7.9	7.2	82	84	.073	.001	.167	.015	.015	9
24	5.3	.12	67	54	7.9	7.3	81	82	.081	.010	.157	.005	.026	9
25	3.0	.12	76	56	7.8	7.4	80	80	.072	.008	.074	.001	.016	12
26	3.8	.103	74	52	7.9	7.3	78	80	.061	.005	.118	.005	.017	10
27	3.2	.13	77	52	7.9	7.3	78	78	.053	.010	.111	.001	.019	12
28	3.0	.13	65	54	8.0	7.3	79	79	.037	.001	.097	.001	.025	10
29	3.2	.09	0	0	7.9	7.3	87	84	.040	.005	.110	.008	.026	10
30	4.3	.08	70	52	7.9	7.3	81	83	.060	.001	.144	.011	.028	11
31	3.0	.09	72	86	7.8	7.3	81	82	.050	.001	.111	.010	.025	12

* All parameters are to be expressed in mg / l except pH and turbidity

MONTHLY REPORTS OF OPERATION FOR # 2 WATER TREATMENT PLANT

Report for Patoka Lake Regional Water District
I.D.E.M. Anna Readle
Signed [Signature]
Public Water Supply I.D. Number 5219012

For the month of December-2021
Title Water Plant Supt.
Certification No. WT058254

Date	Turbidity		Alkalinity		Ph		Hardness		Iron		Manganese		Aluminum	Fluoride	Temp. C	Ammonia
	Raw	Finish	Raw	Finish	Raw	Finish	Raw	Finish	Raw	Finish	Raw	Finish	Finish	Finish	Finish	Finish
1	4.9	.11	64	52	7.7	7.4	80	80	.081	.001	.182	.010	.042	0.56	12	0.92
2	5.8	.11	64	52	7.8	7.4	80	80	.057	.008	.212	.009	.031	0.69	12	0.88
3	3.7	.08	71	62	7.7	7.5	80	75	.080	.001	.156	.017	.011	0.58	12	0.74
4	3.5	.08	70	54	7.8	7.4	72	70	.079	.001	.188	.014	.017	0.65	12	0.7
5	5.2	.08	71	50	7.7	7.3	80	80	.103	.001	.120	.009	.012	0.66	13	0.75
6	4.5	.10	60	50	7.8	7.3	80	80	.083	.001	.173	.004	.022	0.73	11	0.89
7	8.2	.11	68	44	7.7	7.4	80	80	.147	.003	.251	.012	.028	0.69	11	0.88
8	6.1	.08	64	44	7.8	7.4	84	84	.103	.001	.205	.007	.040	0.62	11	0.85
9	7.6	.08	64	44	7.8	7.4	84	84	.142	.001	.246	.013	.025	0.62	11	0.89
10	6.0	.12	64	48	7.7	7.4	84	80	.065	.004	.162	.011	.016	0.67	12	0.87
11	5.0	.12	68	52	7.8	7.4	80	80	.071	.001	.144	.004	.031	0.34	11	0.87
12	9.2	.11	68	48	7.8	7.4	80	80	.154	.002	.293	.011	.034	0.79	11	0.88
13	4.8	.08	75	52	7.6	7.4	80	79	.090	.001	.180	.012	.009	0.64	12	0.89
14	4.3	.09	79	50	7.7	7.4	75	80	.087	.001	.167	.002	.017	0.48	12	0.68
15	5.2	.12	64	52	7.8	7.4	80	80	.079	.001	.170	.007	.009	0.78	11	0.88
16	4.9	.10	64	52	7.9	7.4	80	80	.079	.002	.153	.006	.018	0.79	12	0.89
17	4.4	.09	76	67	7.7	7.4	78	80	.089	.001	.154	.007	.016	0.50	12	0.87
18	4.7	.09	80	74	7.7	7.4	80	80	.112	.009	.146	.012	.009	0.84	12	0.9
19	4.2	.09	79	58	7.8	7.4	79	81	.086	.001	.141	.001	.014	0.90	11	0.88
20	5.2	.12	68	52	7.7	7.3	80	80	.087	.001	.167	.012	.022	0.80	11	0.88
21	4.9	.11	68	52	7.8	7.4	82	80	.079	.002	.162	.014	.019	0.80	11	0.87
22	3.3	.09	75	75	7.7	7.4	80	80	.071	.001	.128	.012	.012	0.70	11	0.88
23	4.3	.10	76	63	7.7	7.4	78	78	.093	.002	.161	.016	.012	0.65	11	0.88
24	5.4	.13	64	56	7.9	7.4	80	80	.084	.004	.181	.010	.017	0.75	11	0.87
25	7.7	.10	68	56	7.9	7.5	80	80	.124	.004	.256	.015	.034	0.90	12	0.85
26	3.3	.12	80	70	7.8	7.4	76	80	.082	.001	.127	.013	.016	0.64	12	0.87
27	3.4	.11	78	69	7.8	7.4	80	78	.080	.001	.131	.010	.018	0.91	12	0.85
28	3.6	.11	76	64	7.8	7.5	76	80	.083	.002	.134	.017	.013	0.64	12	0.9
29	3.1	.12	78	70	7.9	7.4	76	86	.075	.004	.077	.013	.014	0.51	12	0.89
30	3.6	.12	80	70	7.8	7.4	76	84	.075	.001	.144	.001	.018	0.64	12	0.85
31	3.3	.10	74	68	7.9	7.5	76	80	.064	.001	.141	.013	.013	0.62	12	0.88

* All parameters are to be expressed in mg / l except pH and turbidity