

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121824\
 Data File : VX044404.D
 Acq On : 19 Dec 2024 01:24
 Operator : JC/MD
 Sample : P5317-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT191D2-20241216

Quant Time: Dec 19 05:07:54 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121124W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 12 04:28:14 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.550	168	108502	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	198842	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	177204	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	73178	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	84577	44.458	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	88.920%
35) Dibromofluoromethane	5.385	113	65208	47.350	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.700%
50) Toluene-d8	8.647	98	239506	51.549	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.100%
62) 4-Bromofluorobenzene	11.079	95	77874	47.967	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	95.940%

Target Compounds						Qvalue
3) Chloromethane	1.300	50	1231	0.767	ug/l	96
9) 1,1,2-Trichlorotrifluo...	2.312	101	6377	4.816	ug/l	95
12) 1,1-Dichloroethene	2.306	96	1878	1.544	ug/l #	81
16) Acetone	2.392	43	1191	1.536	ug/l #	80
20) Methylene Chloride	2.788	84	587	0.403	ug/l #	73
27) cis-1,2-Dichloroethene	4.489	96	1185	0.722	ug/l	77
30) Chloroform	5.092	83	1305	0.457	ug/l	92
38) Carbon Tetrachloride	5.659	117	1167	0.609	ug/l #	89
44) Trichloroethene	7.123	130	129798	96.301	ug/l	99
55) 1,1,2-Trichloroethane	9.159	97	462	0.343	ug/l #	78

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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