

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX123124\
 Data File : VX044558.D
 Acq On : 30 Dec 2024 20:48
 Operator : JC/MD
 Sample : P5376-14
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT161S1-20241219

Quant Time: Dec 31 00:38:09 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	105027	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	186168	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	163359	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	66627	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	84530	45.904	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	91.800%	
35) Dibromofluoromethane	5.385	113	65178	50.550	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	101.100%	
50) Toluene-d8	8.647	98	233818	53.751	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	107.500%	
62) 4-Bromofluorobenzene	11.079	95	73565	48.398	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	96.800%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.300	50	654	0.421	ug/l	89
12) 1,1-Dichloroethene	2.306	96	599	0.509	ug/l	92
16) Acetone	2.386	43	1084	1.444	ug/l #	77
19) Methyl tert-butyl Ether	3.117	73	2758	0.600	ug/l	94
20) Methylene Chloride	2.794	84	604	0.428	ug/l #	64
24) 1,1-Dichloroethane	3.599	63	6222	2.528	ug/l #	91
30) Chloroform	5.092	83	1222	0.442	ug/l	92
44) Trichloroethene	7.129	130	1465	1.161	ug/l	88
64) Tetrachloroethene	9.269	164	722	0.666	ug/l	84

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

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