

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121824\
 Data File : VX044392.D
 Acq On : 18 Dec 2024 18:42
 Operator : JC/MD
 Sample : P5281-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE115D1-20241213

Quant Time: Dec 19 00:51:43 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	115938	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	206968	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	178386	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	70750	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	87364	42.978	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	85.960%	
35) Dibromofluoromethane	5.385	113	67237	46.906	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	93.820%	
50) Toluene-d8	8.647	98	244633	50.586	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	101.180%	
62) 4-Bromofluorobenzene	11.079	95	77794	46.037	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	92.080%	
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.166	85	2308	1.287	ug/l	100
9) 1,1,2-Trichlorotrifluo...	2.319	101	33134	23.419	ug/l	96
12) 1,1-Dichloroethene	2.313	96	13073	10.060	ug/l	82
16) Acetone	2.392	43	788	0.951	ug/l #	88
20) Methylene Chloride	2.782	84	641	0.412	ug/l #	83
24) 1,1-Dichloroethane	3.605	63	3460	1.273	ug/l #	93
27) cis-1,2-Dichloroethene	4.489	96	4801	2.737	ug/l	92
30) Chloroform	5.093	83	5081	1.664	ug/l	95
38) Carbon Tetrachloride	5.666	117	5756	2.885	ug/l #	83
44) Trichloroethene	7.123	130	757632	540.040	ug/l	99
55) 1,1,2-Trichloroethane	9.159	97	1528	1.089	ug/l #	80
64) Tetrachloroethene	9.281	164	530	0.448	ug/l #	44

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

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