

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112224\
 Data File : VX043966.D
 Acq On : 22 Nov 2024 17:39
 Operator : JC/MD
 Sample : P4846-14
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-MW-EVAL-02S-20241113

Quant Time: Nov 22 22:27:08 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 22 03:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	121293	50.000	ug/l	0.01
34) 1,4-Difluorobenzene	6.757	114	231454	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	196600	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	81016	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	107041	51.453	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.900%			
35) Dibromofluoromethane	5.379	113	76784	46.427	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 92.860%			
50) Toluene-d8	8.647	98	276241	48.455	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 96.900%			
62) 4-Bromofluorobenzene	11.079	95	92396	47.622	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 95.240%			
Target Compounds						
3) Chloromethane	1.295	50	607	0.375	ug/l	91
16) Acetone	2.386	43	1530	1.601	ug/l #	87
27) cis-1,2-Dichloroethene	4.489	96	3306	1.853	ug/l	90
32) 1,1,1-Trichloroethane	5.379	97	12125	4.640	ug/l	99
44) Trichloroethene	7.117	130	146202	80.450	ug/l	96
64) Tetrachloroethene	9.269	164	28564	22.034	ug/l	96

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

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