

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111924\
 Data File : VN084950.D
 Acq On : 19 Nov 2024 16:28
 Operator : JC\MD
 Sample : P4845-16
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FSND-MW-DUP-01-20241111

Quant Time: Nov 20 00:30:57 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 31 18:45:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	168035	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	314106	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	277613	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	124948	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	128607	52.990	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	105.980%		
35) Dibromofluoromethane	8.165	113	102335	48.132	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	96.260%		
50) Toluene-d8	10.565	98	361008	46.094	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	92.180%		
62) 4-Bromofluorobenzene	12.847	95	139048	47.504	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	95.000%		
Target Compounds						
16) Acetone	4.436	43	3856	2.678	ug/l	91
30) Chloroform	7.953	83	1748	0.464	ug/l #	83
38) Carbon Tetrachloride	8.353	117	14173	4.300	ug/l	93
44) Trichloroethene	9.347	130	25282	11.577	ug/l	100
64) Tetrachloroethene	11.100	164	1562	0.846	ug/l #	72

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

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