

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111724\
 Data File : VN084901.D
 Acq On : 17 Nov 2024 17:36
 Operator : JC\MD
 Sample : P4845-16
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

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

Quant Time: Nov 18 03:27:25 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	156883	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	268721	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	236467	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	105675	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.571	65	113675	50.167	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	100.340%	
35) Dibromofluoromethane	8.165	113	92147	50.660	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	101.320%	
50) Toluene-d8	10.565	98	309556	46.200	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	92.400%	
62) 4-Bromofluorobenzene	12.847	95	116917	46.690	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	93.380%	
Target Compounds						
30) Chloroform	7.965	83	1630	0.463	ug/l #	71
38) Carbon Tetrachloride	8.353	117	11853	4.204	ug/l	98
44) Trichloroethene	9.347	130	22205	11.885	ug/l	96
64) Tetrachloroethene	11.100	164	964	0.613	ug/l	95

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

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