

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN103024\
 Data File : VN084601.D
 Acq On : 31 Oct 2024 05:21
 Operator : JC\MD
 Sample : P4611-18
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-6

Quant Time: Nov 01 05:21:28 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	177732	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	326468	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	294514	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	126703	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	129504	50.449	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	100.900%	
35) Dibromofluoromethane	8.159	113	104423	47.255	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	94.500%	
50) Toluene-d8	10.565	98	388088	47.676	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	95.360%	
62) 4-Bromofluorobenzene	12.847	95	142013	46.680	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	93.360%	
Target Compounds						
16) Acetone	4.430	43	12096	13.615	ug/l	99
18) Methyl Acetate	5.024	43	18301	3.004	ug/l	97
43) Isopropyl Acetate	8.688	43	197375	38.195	ug/l #	78

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

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