

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110824\
 Data File : VN084758.D
 Acq On : 08 Nov 2024 18:32
 Operator : JC\MD
 Sample : P4739-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-G2

Quant Time: Nov 08 22:36:02 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	8.218	168	166644	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	302134	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	279048	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	127135	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	119846	49.793	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	99.580%	
35) Dibromofluoromethane	8.159	113	97653	47.750	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	95.500%	
50) Toluene-d8	10.565	98	362890	48.171	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	96.340%	
62) 4-Bromofluorobenzene	12.847	95	141293	50.184	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	100.360%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	18587	24.157	ug/l	97
20) Methylene Chloride	5.271	84	4031	1.904	ug/l #	75
25) 2-Butanone	7.483	43	9941	8.853	ug/l	98
40) Benzene	8.606	78	9511	1.044	ug/l	95
43) Isopropyl Acetate	8.688	43	133630	27.678	ug/l #	81

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

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