

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110824\
 Data File : VN084757.D
 Acq On : 08 Nov 2024 18:08
 Operator : JC\MD
 Sample : P4739-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-G2

Quant Time: Nov 08 22:35:29 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	166407	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	298987	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	280845	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	126928	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	120683	50.212	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 100.420%			
35) Dibromofluoromethane	8.165	113	96186	47.528	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 95.060%			
50) Toluene-d8	10.565	98	356681	47.845	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 95.680%			
62) 4-Bromofluorobenzene	12.847	95	139268	49.986	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 99.980%			
Target Compounds						
16) Acetone	4.430	43	18291	23.764	ug/l	91
20) Methylene Chloride	5.271	84	3846	1.819	ug/l #	78
25) 2-Butanone	7.488	43	10354	9.234	ug/l	94
40) Benzene	8.600	78	14181	1.573	ug/l	95
43) Isopropyl Acetate	8.688	43	125706	26.262	ug/l #	84

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

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