

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112624\
 Data File : VN085046.D
 Acq On : 26 Nov 2024 16:35
 Operator : JC\MD
 Sample : P5000-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-733

Quant Time: Nov 27 05:17:10 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112224W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 23 02:54:10 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.218	168	156639	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	285590	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	259624	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	105556	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.571	65	118192	52.972	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 105.940%			
35) Dibromofluoromethane	8.165	113	94793	49.220	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 98.440%			
50) Toluene-d8	10.565	98	338992	49.115	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 98.220%			
62) 4-Bromofluorobenzene	12.847	95	120921	46.848	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 93.700%			
Target Compounds						
16) Acetone	4.430	43	65433	100.695	ug/l	96
18) Methyl Acetate	5.024	43	6650	2.103	ug/l #	82
20) Methylene Chloride	5.259	84	2573	1.298	ug/l #	80
25) 2-Butanone	7.489	43	7464	7.026	ug/l	92
43) Isopropyl Acetate	8.688	43	316730	64.296	ug/l #	69
51) 4-Methyl-2-Pentanone	10.441	43	11776	5.234	ug/l #	72

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

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