

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020824\
 Data File : VN080964.D
 Acq On : 08 Feb 2024 15:22
 Operator : JC\MD
 Sample : P1388-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 60135-60136-60137-60138-COMP

Quant Time: Feb 09 01:22:28 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020624W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 07 06:22:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	304657	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	657773	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	737526	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	338507	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	196385	51.099	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 102.200%		
35) Dibromofluoromethane	8.171	113	182464	51.695	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 103.400%		
50) Toluene-d8	10.571	98	653394	51.179	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 102.360%		
62) 4-Bromofluorobenzene	12.853	95	295224	63.655	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 127.300%#		
Target Compounds						
						Qvalue
16) Acetone	4.430	43	67745	53.578	ug/l	95
25) 2-Butanone	7.483	43	37708	18.580	ug/l	99
31) Cyclohexane	8.259	56	44145	8.404	ug/l #	80
51) 4-Methyl-2-Pentanone	10.447	43	23541	5.347	ug/l	99
52) Toluene	10.635	92	24898	2.326	ug/l	100
67) Ethyl Benzene	11.971	91	36580	1.374	ug/l	98
68) m/p-Xylenes	12.071	106	26796	2.649	ug/l	94
69) o-Xylene	12.400	106	21015	2.141	ug/l	92

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

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