

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080321\
 Data File : VN067994.D
 Acq On : 3 Aug 2021 20:16
 Operator : JC/MD
 Sample : M3261-08
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 41938

Quant Time: Aug 04 03:43:06 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 29 15:02:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	307661	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	590027	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	594719	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	319558	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	242739	55.732	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 111.460%			
35) Dibromofluoromethane	8.024	113	182935	53.201	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 106.400%			
50) Toluene-d8	10.443	98	762704	48.211	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 96.420%			
62) 4-Bromofluorobenzene	12.733	95	270226	46.868	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 93.740%			
Target Compounds					Qvalue	
16) Acetone	4.291	43	263640	165.895	ug/l	95
17) Carbon Disulfide	4.524	76	13249	1.566	ug/l	99
25) 2-Butanone	7.348	43	49680	21.548	ug/l	92
52) Toluene	10.507	92	2676303	254.917	ug/l	99
67) Ethyl Benzene	11.848	91	9447	0.439	ug/l	97
68) m/p-Xylenes	11.950	106	16354	5.729	ug/l	86
69) o-Xylene	12.283	106	15917	2.010	ug/l	85
79) 2-Chlorotoluene	13.010	91	407894	22.748	ug/l	92
82) 4-Chlorotoluene	13.106	91	496180	27.561	ug/l	94
84) 1,2,4-Trimethylbenzene	13.369	105	51316	2.506	ug/l	95
86) p-Isopropyltoluene	13.618	119	4795473	245.234	ug/l #	68
95) Naphthalene	15.509	128	15487	4.210	ug/l	97

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

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