

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080321\
 Data File : VN067993.D
 Acq On : 3 Aug 2021 19:52
 Operator : JC/MD
 Sample : M3261-07
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 41937

Quant Time: Aug 04 03:42:44 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	284371	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	546886	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	544899	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	212972	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	225088	55.912	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 111.820%			
35) Dibromofluoromethane	8.024	113	171840	53.916	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 107.840%			
50) Toluene-d8	10.446	98	700388	47.800	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 95.600%			
62) 4-Bromofluorobenzene	12.733	95	243599	45.719	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 91.440%			
Target Compounds						
16) Acetone	4.299	43	48025	32.695	ug/l	92
17) Carbon Disulfide	4.535	76	9122	1.166	ug/l	99
25) 2-Butanone	7.351	43	27745	13.019	ug/l	96
52) Toluene	10.507	92	1410864	144.986	ug/l	98
67) Ethyl Benzene	11.848	91	8076	0.409	ug/l #	75
68) m/p-Xylenes	11.956	106	14905	5.721	ug/l	93
69) o-Xylene	12.283	106	10988	1.514	ug/l	99
79) 2-Chlorotoluene	13.010	91	263541	22.053	ug/l	94
82) 4-Chlorotoluene	13.109	91	463031	38.591	ug/l	93
84) 1,2,4-Trimethylbenzene	13.369	105	30158	2.210	ug/l	96
86) p-Isopropyltoluene	13.613	119	5407	0.415	ug/l #	70
95) Naphthalene	15.507	128	8828	4.052	ug/l #	91

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

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