

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042821\
 Data File : VN066850.D
 Acq On : 29 Apr 2021 7:17
 Operator : JC/MD
 Sample : M2199-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SAMPLE-3-COMP

Quant Time: Apr 29 07:47:43 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 27 16:02:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.594	168	313803	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.519	114	478038	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.348	117	450908	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.290	152	184523	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.956	65	173691	43.08	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 86.16%			
35) Dibromofluoromethane	7.513	113	156338	52.98	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 105.96%			
50) Toluene-d8	10.031	98	598333	51.66	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 103.32%			
62) 4-Bromofluorobenzene	12.349	95	178404	45.52	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 91.04%			
Target Compounds						
						Qvalue
16) Acetone	3.758	43	16623	10.93	ug/l	92
17) Carbon Disulfide	3.989	76	20876	2.17	ug/l	99
18) Methyl Acetate	4.263	43	4443	2.17	ug/l #	65
20) Methylene Chloride	4.472	84	13134	3.41	ug/l	98
43) Isopropyl Acetate	8.095	43	24119	3.22	ug/l #	90

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

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