

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022322\
 Data File : VN071039.D
 Acq On : 23 Feb 2022 17:44
 Operator : JC\MD
 Sample : N1642-02 50X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31454

Quant Time: Feb 24 00:41:04 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Feb 21 07:57:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.080	168	853968	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.969	114	1356841	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1202969	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	422725	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	508575	52.135	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 104.260%			
35) Dibromofluoromethane	8.022	113	395039	48.949	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 97.900%			
50) Toluene-d8	10.439	98	1636308	50.487	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 100.980%			
62) 4-Bromofluorobenzene	12.727	95	559473	48.293	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 96.580%			
Target Compounds					Qvalue	
16) Acetone	4.287	43	603174	112.275	ug/l	91
51) 4-Methyl-2-Pentanone	10.333	43	7134	0.525	ug/l	88
52) Toluene	10.504	92	616130	25.070	ug/l	95
67) Ethyl Benzene	11.839	91	98109	2.162	ug/l	96
68) m/p-Xylenes	11.951	106	140806	8.035	ug/l	99
69) o-Xylene	12.274	106	45041	2.631	ug/l	98

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

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