

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022422\
 Data File : VN071057.D
 Acq On : 24 Feb 2022 15:56
 Operator : JC\MD
 Sample : N1642-08 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31409

Quant Time: Feb 25 02:44:27 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	843845	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1346440	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1185249	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	418786	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	513506	53.272	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 106.540%	
35) Dibromofluoromethane	8.016	113	395067	49.330	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 98.660%	
50) Toluene-d8	10.439	98	1620322	50.380	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 100.760%	
62) 4-Bromofluorobenzene	12.727	95	550990	47.928	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 95.860%	
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	268129	50.508	ug/l	96
18) Methyl Acetate	4.857	43	25069	1.573	ug/l #	89
25) 2-Butanone	7.333	43	24700	3.498	ug/l	94
52) Toluene	10.504	92	561990	23.044	ug/l	95
67) Ethyl Benzene	11.839	91	27273	0.610	ug/l	94
68) m/p-Xylenes	11.951	106	39158	2.268	ug/l	93
69) o-Xylene	12.274	106	13161	0.780	ug/l	92

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

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