

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020222\
 Data File : VN070912.D
 Acq On : 3 Feb 2022 00:31
 Operator : JC/MD
 Sample : N1342-01DL 20X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31523DL

Quant Time: Feb 03 06:00:44 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020222W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 02 15:30:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1080371	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1689130	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1502634	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	625862	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	669444	45.777	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	91.560%		
35) Dibromofluoromethane	8.024	113	485773	48.007	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	96.020%		
50) Toluene-d8	10.440	98	2063453	49.666	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	99.340%		
62) 4-Bromofluorobenzene	12.731	95	726256	48.513	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	97.020%		
Target Compounds						
					Qvalue	
8) Diethyl Ether	3.824	74	29957	3.657	ug/l	87
11) Tert butyl alcohol	5.382	59	209364	68.132	ug/l	100
16) Acetone	4.293	43	224151	25.634	ug/l	98
18) Methyl Acetate	4.862	43	35350	1.315	ug/l #	84
25) 2-Butanone	7.343	43	87784	6.955	ug/l	92
39) Methylcyclohexane	9.459	83	97638	4.832	ug/l	91
40) Benzene	8.461	78	11831062	229.083	ug/l	98
52) Toluene	10.505	92	3805796	118.933	ug/l	96
67) Ethyl Benzene	11.843	91	201437	3.514	ug/l	96
68) m/p-Xylenes	11.948	106	389968	18.358	ug/l	93
69) o-Xylene	12.280	106	98005	4.610	ug/l	95

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

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