

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042022\
 Data File : VN072101.D
 Acq On : 21 Apr 2022 01:29
 Operator : JC\MD
 Sample : N2482-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COP-1R

Quant Time: Apr 21 03:41:01 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 31 01:57:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	615693	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1030513	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1127842	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	424077	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	324506	46.168	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	92.340%
35) Dibromofluoromethane	8.022	113	292420	48.222	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	96.440%
50) Toluene-d8	10.439	98	1329672	53.602	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	107.200%
62) 4-Bromofluorobenzene	12.727	95	481119	59.969	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	119.940%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	76852	25.584	ug/l	94
25) 2-Butanone	7.351	43	7250	1.674	ug/l	100
37) Ethyl Acetate	7.410	43	60054	6.479	ug/l	96
40) Benzene	8.457	78	34139	1.276	ug/l	99
52) Toluene	10.498	92	142489	8.426	ug/l	100
67) Ethyl Benzene	11.839	91	52222	1.366	ug/l	96
68) m/p-Xylenes	11.945	106	86980	5.921	ug/l	99
69) o-Xylene	12.274	106	68572	4.703	ug/l	96
80) 1,3,5-Trimethylbenzene	13.057	105	28200	0.996	ug/l	98
84) 1,2,4-Trimethylbenzene	13.363	105	32160	1.164	ug/l	98
95) Naphthalene	15.509	128	10139	4.288	ug/l	99

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

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