

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042022\
 Data File : VN072111.D
 Acq On : 21 Apr 2022 05:31
 Operator : JC\MD
 Sample : N2482-16
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SPN-7

Quant Time: Apr 21 07:16:07 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	619953	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1052668	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1100198	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	417240	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	329191	46.512	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	93.020%
35) Dibromofluoromethane	8.016	113	297263	47.989	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	95.980%
50) Toluene-d8	10.439	98	1314259	51.866	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	103.740%
62) 4-Bromofluorobenzene	12.727	95	449955	54.904	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	109.800%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	64156	21.211	ug/l	97
25) 2-Butanone	7.345	43	5920	1.358	ug/l #	89
37) Ethyl Acetate	7.416	43	17406	1.838	ug/l	96
40) Benzene	8.457	78	7346	0.269	ug/l #	89
68) m/p-Xylenes	11.945	106	6409	0.447	ug/l	92
73) Isopropylbenzene	12.574	105	18829	0.536	ug/l	99
78) n-propylbenzene	12.915	91	12884	0.348	ug/l	99

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

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