

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032122\
 Data File : VN071447.D
 Acq On : 22 Mar 2022 03:44
 Operator : JC\MD
 Sample : N2017-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-3

Quant Time: Mar 22 06:31:40 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 06:52:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.080	168	721702	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1233870	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1169575	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	417570	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	476267	48.151	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	96.300%
35) Dibromofluoromethane	8.016	113	360431	46.291	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	92.580%
50) Toluene-d8	10.439	98	1510559	48.185	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.360%
62) 4-Bromofluorobenzene	12.727	95	553468	52.137	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	104.280%
Target Compounds						
					Qvalue	
16) Acetone	4.292	43	9782	2.033	ug/l	98
18) Methyl Acetate	4.857	43	9797	1.221	ug/l	96
30) Chloroform	7.816	83	17338	1.110	ug/l	99
44) Trichloroethene	9.216	130	4419	0.498	ug/l	90
64) Tetrachloroethene	10.980	164	1746	0.214	ug/l #	71

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

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