

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032122\
 Data File : VN071448.D
 Acq On : 22 Mar 2022 04:08
 Operator : JC\MD
 Sample : N2017-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-4

Manual Integrations
 APPROVED

Reviewed By : John Carlone 03/22/2022
 Supervised By : Mahesh Dadoda 03/22/2022

Quant Time: Mar 22 06:32:04 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	8.081	168	771626	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1307300	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1273847	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	478862	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	504409	47.696	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	95.400%		
35) Dibromofluoromethane	8.022	113	383630	46.503	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	93.000%		
50) Toluene-d8	10.439	98	1625106	48.927	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	97.860%		
62) 4-Bromofluorobenzene	12.727	95	613265	54.525	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	109.040%		
Target Compounds					Qvalue	
16) Acetone	4.304	43	8982m	1.746	ug/l	
18) Methyl Acetate	4.869	43	9638	1.161	ug/l #	85
30) Chloroform	7.810	83	25396	1.520	ug/l	91
44) Trichloroethene	9.216	130	7444	0.792	ug/l	97
64) Tetrachloroethene	10.980	164	7638	0.860	ug/l	93

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

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