

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
 Data File : VN071757.D
 Acq On : 01 Apr 2022 16:55
 Operator : JC\MD
 Sample : N2090-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-51

Manual Integrations
 APPROVED

Reviewed By : John Carlone 04/02/2022
 Supervised By : Mahesh Dadoda 04/02/2022

Quant Time: Apr 02 00:55:42 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	8.087	168	656715	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1113740	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1146303	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.669	152	424091	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	379695	50.645	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 101.300%			
35) Dibromofluoromethane	8.016	113	320060	48.836	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 97.680%			
50) Toluene-d8	10.439	98	1411450	52.647	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 105.300%			
62) 4-Bromofluorobenzene	12.727	95	486467	56.105	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 112.200%			
Target Compounds					Qvalue	
16) Acetone	4.287	43	14908m	4.653	ug/l	
19) Methyl tert-butyl Ether	5.616	73	18371	0.912	ug/l	98
40) Benzene	8.457	78	36083	1.248	ug/l #	89
73) Isopropylbenzene	12.574	105	21412	0.599	ug/l	98
78) n-propylbenzene	12.916	91	13494	0.359	ug/l	99

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

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