

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120721\
 Data File : VN069811.D
 Acq On : 7 Dec 2021 13:42
 Operator : JC/MD
 Sample : M4956-22
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EB02-20211206

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/08/2021
 Supervised By : Amit Patel 12/14/2021

Quant Time: Dec 08 04:34:52 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1300814	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2340062	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2455798	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	934925	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	991820	49.382	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	98.760%		
35) Dibromofluoromethane	8.021	113	693934	47.343	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	94.680%		
50) Toluene-d8	10.443	98	2976163	50.567	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	101.140%		
62) 4-Bromofluorobenzene	12.731	95	1156764	54.252	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	108.500%		
Target Compounds					Qvalue	
16) Acetone	4.291	43	75356	6.237	ug/l	95
20) Methylene Chloride	5.092	84	17936m	1.117	ug/l	
25) 2-Butanone	7.351	43	13050	0.795	ug/l	93
68) m/p-Xylenes	11.953	106	17837	0.541	ug/l	86
69) o-Xylene	12.283	106	11579	0.349	ug/l	90

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

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