

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121721\
 Data File : VN070173.D
 Acq On : 17 Dec 2021 19:09
 Operator : JC/MD
 Sample : M5039-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-29

Quant Time: Dec 18 05:47:45 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

Manual Integrations APPROVED

Reviewed By : John Carlone 12/20/2021
 Supervised By : Mahesh Dadoda 12/23/2021

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	941555	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1582608	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	1520118	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	716106	50.000	ug/l	# 0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	731945	50.348	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 100.700%	
35) Dibromofluoromethane	8.024	113	480296	48.451	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 96.900%	
50) Toluene-d8	10.443	98	1987935	49.943	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 99.880%	
62) 4-Bromofluorobenzene	12.731	95	816949	56.653	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 113.300%	
Target Compounds						
						Qvalue
3) Chloromethane	2.303	50	4974	0.362	ug/l	91
11) Tert butyl alcohol	5.385	59	8970784	3264.520	ug/l	99
16) Acetone	4.291	43	2111410	241.441	ug/l	97
17) Carbon Disulfide	4.521	76	20157	0.870	ug/l	96
19) Methyl tert-butyl Ether	5.618	73	29113982	834.866	ug/l	# 85
22) Diisopropyl ether	6.501	45	676135	17.117	ug/l	88
25) 2-Butanone	7.340	43	1739955	146.437	ug/l	92
31) Cyclohexane	8.096	56	1742544	86.131	ug/l	89
39) Methylcyclohexane	9.462	83	1405777	78.445	ug/l	89
40) Benzene	8.461	78	19107991	414.933	ug/l	97
51) 4-Methyl-2-Pentanone	10.333	43	239543	11.128	ug/l	93
52) Toluene	10.505	92	5061642	180.035	ug/l	98
67) Ethyl Benzene	11.827	91	26452973m	472.224	ug/l	
68) m/p-Xylenes	11.940	106	28392075m	1392.006	ug/l	
69) o-Xylene	12.280	106	14370779	699.483	ug/l	# 1
73) Isopropylbenzene	12.578	105	7424846	116.535	ug/l	98
78) n-propylbenzene	12.913	91	17583159m	248.003	ug/l	
80) 1,3,5-Trimethylbenzene	13.047	105	20518973m	400.122	ug/l	
84) 1,2,4-Trimethylbenzene	13.350	105	27976498m	572.836	ug/l	
85) sec-Butylbenzene	13.501	105	602402	9.780	ug/l	# 64
86) p-Isopropyltoluene	13.616	119	307346m	6.384	ug/l	

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Badoda

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

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