

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062121\
 Data File : VN067425.D
 Acq On : 21 Jun 2021 18:40
 Operator : JC/MD
 Sample : M2795-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 B3GW

Manual Integrations
 APPROVED

MMDadoda
 6/22/2021 6:24:56 PM

Quant Time: Jun 22 03:12:42 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 15 05:19:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	278838	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	576003	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	490288	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.680	152	161325	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.442	65	273969	49.806	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	99.620%		
35) Dibromofluoromethane	8.027	113	178976	46.719	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	93.440%		
50) Toluene-d8	10.446	98	742993	50.602	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	101.200%		
62) 4-Bromofluorobenzene	12.733	95	245866	40.575	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	81.160%		
Target Compounds						Qvalue
3) Chloromethane	2.303	50	2488	0.588	ug/l	96
16) Acetone	4.293	43	15153	7.178	ug/l	96
17) Carbon Disulfide	4.527	76	12194	1.129	ug/l #	94
19) Methyl tert-butyl Ether	5.621	73	4734m	0.353	ug/l	
25) 2-Butanone	7.351	43	6320	2.095	ug/l	96
27) cis-1,2-Dichloroethene	7.335	96	2051m	0.470	ug/l	
39) Methylcyclohexane	9.464	83	3858m	0.564	ug/l	
40) Benzene	8.461	78	5632	0.312	ug/l	96

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

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