

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062421\
 Data File : VN067490.D
 Acq On : 24 Jun 2021 18:07
 Operator : JC/MD
 Sample : M2804-05
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DUP08-20210621

Manual Integrations
 APPROVED

MMDadoda
 6/25/2021 6:23:00 PM

Quant Time: Jun 25 02:15:37 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062421W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 24 12:18:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	241244	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	518095	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	449426	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	142522	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	244514	49.524	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	99.040%		
35) Dibromofluoromethane	8.027	113	157038	49.200	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	98.400%		
50) Toluene-d8	10.446	98	685173	50.330	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	100.660%		
62) 4-Bromofluorobenzene	12.733	95	225869	46.568	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	93.140%		
Target Compounds						Qvalue
12) 1,1-Dichloroethene	4.186	96	2113m	0.640	ug/l	
16) Acetone	4.293	43	6260m	3.313	ug/l	
27) cis-1,2-Dichloroethene	7.327	96	5638	1.344	ug/l	81
30) Chloroform	7.823	83	5105	0.689	ug/l	96
44) Trichloroethene	9.217	130	5268	1.556	ug/l	98
64) Tetrachloroethene	10.979	164	3755	1.285	ug/l	90

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

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