

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062921\
 Data File : VN067563.D
 Acq On : 29 Jun 2021 19:19
 Operator : JC/MD
 Sample : M2827-06
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ERM-7

Manual Integrations
 APPROVED

MMDadoda
 7/5/2021 1:52:22 PM

Quant Time: Jun 30 03:40:54 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062821W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 28 14:31:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	163493	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	307452	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	294155	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	111845	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	123757	55.309	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	110.620%
35) Dibromofluoromethane	8.024	113	90429	47.440	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	94.880%
50) Toluene-d8	10.443	98	411114	53.092	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	106.180%
62) 4-Bromofluorobenzene	12.734	95	132480	48.980	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	97.960%
Target Compounds						Qvalue
16) Acetone	4.299	43	7867	10.659	ug/l #	89
18) Methyl Acetate	4.870	43	3216m	1.634	ug/l	

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

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