

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120721\
 Data File : VN069841.D
 Acq On : 8 Dec 2021 3:41
 Operator : JC/MD
 Sample : M4957-07
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RE125D1-20211206

Quant Time: Dec 08 06:38:28 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	8.088	168	1315870	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2371300	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2543476	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	986917	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	1013939	49.906	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	99.820%
35) Dibromofluoromethane	8.021	113	707774	47.651	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	95.300%
50) Toluene-d8	10.440	98	3055750	51.236	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	102.480%
62) 4-Bromofluorobenzene	12.731	95	1205922	55.813	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	111.620%
Target Compounds						
					Qvalue	
9) 1,1,2-Trichlorotrifluo...	4.215	101	43716	3.361	ug/l	97
12) 1,1-Dichloroethene	4.178	96	10830	0.864	ug/l	87
16) Acetone	4.293	43	13592	1.112	ug/l	96
19) Methyl tert-butyl Ether	5.618	73	20852	0.428	ug/l	98
30) Chloroform	7.823	83	14896	0.516	ug/l	98
44) Trichloroethene	9.215	130	2162309	131.297	ug/l	91
64) Tetrachloroethene	10.977	164	88801	5.149	ug/l	97

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

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