

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX020922\
 Data File : VX026878.D
 Acq On : 09 Feb 2022 18:53
 Operator : JC/MD
 Sample : N1446-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-107

Quant Time: Feb 10 01:09:27 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020822W.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 08 07:07:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	127732	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	215489	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	189862	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	79311	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	81824	53.365	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 106.720%	
35) Dibromofluoromethane	5.391	113	66778	52.014	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.020%	
50) Toluene-d8	8.653	98	248556	52.743	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 105.480%	
62) 4-Bromofluorobenzene	11.079	95	89012	53.649	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 107.300%	
Target Compounds						Qvalue
16) Acetone	2.385	43	882	1.310	ug/l	100
19) Methyl tert-butyl Ether	3.123	73	3741	0.809	ug/l	96
27) cis-1,2-Dichloroethene	4.489	96	2451	1.552	ug/l	86
30) Chloroform	5.104	83	1326	0.506	ug/l	95
42) 1,2-Dichloroethane	6.092	62	2515	1.209	ug/l	99

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

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