

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX020922\
 Data File : VX026902.D
 Acq On : 10 Feb 2022 05:15
 Operator : JC/MD
 Sample : N1447-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-139A

Quant Time: Feb 10 07:30:22 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	5.556	168	128012	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	218327	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	198160	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	85992	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	84979	55.301	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	110.600%
35) Dibromofluoromethane	5.391	113	69024	53.065	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	106.120%
50) Toluene-d8	8.653	98	255445	53.500	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	107.000%
62) 4-Bromofluorobenzene	11.079	95	96712	57.532	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	115.060%
Target Compounds						
						Qvalue
16) Acetone	2.386	43	1850	2.741	ug/l	96
19) Methyl tert-butyl Ether	3.117	73	5535	1.194	ug/l	98
30) Chloroform	5.099	83	1250	0.476	ug/l	95
40) Benzene	6.044	78	6079	1.016	ug/l	93
42) 1,2-Dichloroethane	6.092	62	1189	0.564	ug/l	85

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

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