

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032122\
 Data File : VX027591.D
 Acq On : 21 Mar 2022 17:12
 Operator : JC/MD
 Sample : N1998-13
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-139

Quant Time: Mar 22 05:48:06 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 17 15:02:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	344045	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	566972	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	517413	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	232117	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	197640	46.632	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	93.260%
35) Dibromofluoromethane	5.391	113	185693	50.906	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	101.820%
50) Toluene-d8	8.653	98	662990	48.030	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.060%
62) 4-Bromofluorobenzene	11.085	95	231221	45.573	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	91.140%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	5521	3.196	ug/l #	86
19) Methyl tert-butyl Ether	3.123	73	15757	1.462	ug/l #	86
30) Chloroform	5.105	83	3356	0.512	ug/l	99
40) Benzene	6.044	78	15478	1.082	ug/l	97
42) 1,2-Dichloroethane	6.092	62	3790	0.707	ug/l	89

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

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