

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110822\
 Data File : VX032655.D
 Acq On : 08 Nov 2022 15:40
 Operator : JC/MD
 Sample : N5514-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-19R

Quant Time: Nov 09 04:39:56 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 08:28:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	109211	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	179687	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	157482	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	72049	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	74917	49.370	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.740%
35) Dibromofluoromethane	5.385	113	59078	46.093	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.180%
50) Toluene-d8	8.647	98	217660	48.837	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.680%
62) 4-Bromofluorobenzene	11.079	95	82134	47.178	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.360%
Target Compounds						
					Qvalue	
16) Acetone	2.373	43	645	1.169	ug/l	92
19) Methyl tert-butyl Ether	3.117	73	41612	10.810	ug/l	99
22) Diisopropyl ether	3.763	45	1413	0.374	ug/l #	90
30) Chloroform	5.098	83	3274	1.374	ug/l	97
42) 1,2-Dichloroethane	6.098	62	4549	2.203	ug/l	98

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

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