

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101022\
 Data File : VX032017.D
 Acq On : 10 Oct 2022 17:04
 Operator : JC/MD
 Sample : N4937-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-102

Quant Time: Oct 11 05:13:12 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 06 14:28:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	54200	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	87893	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	95668	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	58408	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	62459	55.800	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.600%
35) Dibromofluoromethane	5.391	113	49867	50.123	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.240%
50) Toluene-d8	8.647	98	182008	50.840	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	101.680%
62) 4-Bromofluorobenzene	11.079	95	62341	49.257	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.520%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.983	59	6545	46.997	ug/l	96
16) Acetone	2.386	43	3202	8.417	ug/l #	82
19) Methyl tert-butyl Ether	3.117	73	2108	0.761	ug/l #	79
40) Benzene	6.044	78	10933	2.830	ug/l	99
42) 1,2-Dichloroethane	6.092	62	3329	2.365	ug/l	97
73) Isopropylbenzene	10.964	105	2371	0.538	ug/l	97

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

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