

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101022\
 Data File : VX032018.D
 Acq On : 10 Oct 2022 17:27
 Operator : JC/MD
 Sample : N4937-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-104S

Quant Time: Oct 11 05:59:55 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	49443	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	84775	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	91202	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	54918	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	61806	60.529	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 121.060%	
35) Dibromofluoromethane	5.385	113	50101	52.210	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.420%	
50) Toluene-d8	8.647	98	171186	49.576	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 99.160%	
62) 4-Bromofluorobenzene	11.079	95	58521	47.939	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.880%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	681	1.962	ug/l #	83
19) Methyl tert-butyl Ether	3.111	73	21096	8.347	ug/l	94
30) Chloroform	5.093	83	3669	2.281	ug/l	89

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

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