

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100722\
 Data File : VX031995.D
 Acq On : 08 Oct 2022 05:48
 Operator : JC/MD
 Sample : N4937-21
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TB-093022

Quant Time: Oct 10 04:51:30 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	58716	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	93882	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	98240	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	60110	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	64117	52.875	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	105.760%
35) Dibromofluoromethane	5.385	113	51679	48.630	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.260%
50) Toluene-d8	8.653	98	185901	48.615	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	97.240%
62) 4-Bromofluorobenzene	11.079	95	63663	47.093	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.180%
Target Compounds						
						Qvalue
16) Acetone	2.398	43	1056	2.562	ug/l #	70
20) Methylene Chloride	2.788	84	435	0.402	ug/l #	85
40) Benzene	6.043	78	5936	1.439	ug/l #	82
52) Toluene	8.720	92	4575	1.696	ug/l	94
67) Ethyl Benzene	10.195	91	2564	0.509	ug/l	100
68) m/p-Xylenes	10.305	106	4773	2.385	ug/l	92
69) o-Xylene	10.646	106	1456	0.749	ug/l	95
80) 1,3,5-Trimethylbenzene	11.451	105	1193	0.317	ug/l	89
84) 1,2,4-Trimethylbenzene	11.756	105	3502	0.917	ug/l	93

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

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