

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112322\
 Data File : VX033049.D
 Acq On : 24 Nov 2022 09:07
 Operator : JC/MD
 Sample : N5742-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-43

Quant Time: Nov 25 06:38:51 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 22 13:58:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	131348	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	212817	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	189983	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	89312	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	70801	55.509	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	111.020%
35) Dibromofluoromethane	5.385	113	56464	47.062	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.120%
50) Toluene-d8	8.647	98	192706	53.535	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	107.060%
62) 4-Bromofluorobenzene	11.079	95	69743	44.252	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	88.500%
Target Compounds						
						Qvalue
16) Acetone	2.392	43	1606	2.688	ug/l	97
19) Methyl tert-butyl Ether	3.111	73	1845	0.464	ug/l #	1
39) Methylcyclohexane	7.385	83	614	0.281	ug/l #	86
40) Benzene	6.038	78	3266	0.607	ug/l	95
73) Isopropylbenzene	10.964	105	2368	0.371	ug/l	99
78) n-propylbenzene	11.305	91	3326	0.456	ug/l	96
84) 1,2,4-Trimethylbenzene	11.756	105	2446	0.458	ug/l	94

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

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