

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101823\
 Data File : VX038365.D
 Acq On : 18 Oct 2023 16:12
 Operator : JC/MD
 Sample : 04940-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GW-19-10-13

Quant Time: Oct 19 01:42:07 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 05 15:27:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	110626	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	201020	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	210720	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	111212	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	80252	45.102	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.200%
35) Dibromofluoromethane	5.391	113	63700	43.977	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	87.960%
50) Toluene-d8	8.647	98	255244	46.115	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	92.220%
62) 4-Bromofluorobenzene	11.079	95	98805	47.087	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.180%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	1958	2.637	ug/l	100
30) Chloroform	5.093	83	2471	0.981	ug/l	96
44) Trichloroethene	7.129	130	2210	1.512	ug/l	87
48) Methyl methacrylate	7.708	41	686	0.406	ug/l #	85
52) Toluene	8.720	92	162393	44.623	ug/l	93
64) Tetrachloroethene	9.275	164	4682	3.432	ug/l #	87
67) Ethyl Benzene	10.201	91	2232	0.284	ug/l	91
68) m/p-Xylenes	10.305	106	1861	0.611	ug/l	95
69) o-Xylene	10.646	106	1014	0.341	ug/l	94

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

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