

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091323\
 Data File : VX037690.D
 Acq On : 13 Sep 2023 13:16
 Operator : JC/MD
 Sample : 04373-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1044-MW-11(15)

Quant Time: Sep 14 02:19:07 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091123W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 12 10:10:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.549	168	218374	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	380680	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	350792	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	147548	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	154543	43.741	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	87.480%	
35) Dibromofluoromethane	5.385	113	122579	45.428	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	90.860%	
50) Toluene-d8	8.652	98	463052	45.502	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	91.000%#	
62) 4-Bromofluorobenzene	11.079	95	157675	42.947	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	85.900%	
Target Compounds						
					Qvalue	
16) Acetone	2.385	43	13250	9.276	ug/l	99
25) 2-Butanone	4.580	43	14743	6.330	ug/l	100
27) cis-1,2-Dichloroethene	4.489	96	2023	0.642	ug/l	77
44) Trichloroethene	7.134	130	3969	1.428	ug/l	92
51) 4-Methyl-2-Pentanone	8.579	43	21172	4.812	ug/l	97
64) Tetrachloroethene	9.274	164	18346	8.072	ug/l	99

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

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