

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090923\
 Data File : VX037638.D
 Acq On : 09 Sep 2023 19:27
 Operator : JC/MD
 Sample : 04338-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP2

Quant Time: Sep 11 01:28:15 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 09 06:38:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	143297	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	240292	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	218346	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	94868	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	79770	41.239	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	82.480%
35) Dibromofluoromethane	5.385	113	72199	42.958	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	85.920%
50) Toluene-d8	8.647	98	284383	46.134	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	92.260%
62) 4-Bromofluorobenzene	11.079	95	101341	43.621	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.240%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	64558	81.158	ug/l #	81
20) Methylene Chloride	2.788	84	9794	4.442	ug/l #	84
25) 2-Butanone	4.568	43	15484	12.835	ug/l #	85
29) Tetrahydrofuran	5.013	42	4091	5.293	ug/l #	84
43) Isopropyl Acetate	6.342	43	88597	25.983	ug/l #	87
84) 1,2,4-Trimethylbenzene	11.756	105	6887	1.140	ug/l	100

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

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