

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091423\
 Data File : VX037733.D
 Acq On : 14 Sep 2023 17:03
 Operator : JC/MD
 Sample : 04390-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-RW10A-GW-58-60

Quant Time: Sep 14 17:43:01 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	5.550	168	244856	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	420744	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	385834	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	171047	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	162625	41.050	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	82.100%
35) Dibromofluoromethane	5.385	113	131848	44.210	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	88.420%
50) Toluene-d8	8.647	98	503657	44.780	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	89.560%#
62) 4-Bromofluorobenzene	11.079	95	180937	44.590	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	89.180%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	9557	5.967	ug/l #	87
25) 2-Butanone	4.587	43	4370	1.673	ug/l	94
27) cis-1,2-Dichloroethene	4.489	96	8637	2.444	ug/l	80
30) Chloroform	5.093	83	6375	1.054	ug/l	99
44) Trichloroethene	7.129	130	6967	2.267	ug/l	91
64) Tetrachloroethene	9.275	164	11168	4.467	ug/l	98

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

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