

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091523\
 Data File : VX037758.D
 Acq On : 15 Sep 2023 16:28
 Operator : JC/MD
 Sample : 04390-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

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

Quant Time: Sep 16 05:46:33 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091523W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 16 04:55:44 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	208648	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	336829	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	309413	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	138643	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	123732	45.136	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	90.280%		
35) Dibromofluoromethane	5.385	113	102635	45.540	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	91.080%		
50) Toluene-d8	8.647	98	404761	48.335	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	96.680%		
62) 4-Bromofluorobenzene	11.079	95	140950	46.320	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	92.640%		
Target Compounds						
16) Acetone	2.386	43	8576	6.503	ug/l	Qvalue 90
25) 2-Butanone	4.580	43	3732	2.005	ug/l	# 88
27) cis-1,2-Dichloroethene	4.489	96	9771	3.526	ug/l	77
30) Chloroform	5.098	83	5883	1.291	ug/l	93
44) Trichloroethene	7.129	130	8370	3.148	ug/l	100
64) Tetrachloroethene	9.275	164	11940	5.658	ug/l	98

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

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