

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091823\
 Data File : VX037775.D
 Acq On : 18 Sep 2023 14:30
 Operator : JC/MD
 Sample : 04437-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-RW10A-GW-240-242

Quant Time: Sep 18 17:57:02 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	5.556	168	297023	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	492054	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	456957	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	222070	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	179927	46.106	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	92.220%
35) Dibromofluoromethane	5.391	113	153368	46.583	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.160%
50) Toluene-d8	8.653	98	583573	47.704	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.400%
62) 4-Bromofluorobenzene	11.079	95	214899	48.343	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.680%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	9253	4.929	ug/l	# 88
17) Carbon Disulfide	2.514	76	4406	0.606	ug/l	# 95
22) Diisopropyl ether	3.763	45	6122	0.609	ug/l	# 68
24) 1,1-Dichloroethane	3.611	63	2611	0.433	ug/l	# 91
25) 2-Butanone	4.587	43	3104	1.172	ug/l	# 89

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

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