

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090623\
 Data File : VX037556.D
 Acq On : 06 Sep 2023 15:49
 Operator : JC/MD
 Sample : 04256-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

Quant Time: Sep 06 22:46:53 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081723W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 18 09:30:38 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 09/07/2023
 Supervised By : Mahesh Dadoda 09/07/2023

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	131036	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	217661	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	199007	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	85362	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	76046	40.255	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	80.500%		
35) Dibromofluoromethane	5.385	113	66038	45.811	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	91.620%		
50) Toluene-d8	8.647	98	259598	49.646	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	99.300%		
62) 4-Bromofluorobenzene	11.079	95	93776	45.635	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	91.280%		
Target Compounds						
16) Acetone	2.386	43	18623	25.199	ug/l #	82
20) Methylene Chloride	2.788	84	8856	5.358	ug/l #	82
43) Isopropyl Acetate	6.348	43	84901	25.562	ug/l #	88
51) 4-Methyl-2-Pentanone	8.574	43	9082m	4.503	ug/l	

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

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