

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090823\
 Data File : VX037611.D
 Acq On : 08 Sep 2023 19:03
 Operator : JC/MD
 Sample : PB155394ZHE#03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB155394ZHE#03

Manual Integrations
 APPROVED

Reviewed By : John Carlone 09/09/2023
 Supervised By : Mahesh Dadoda 09/11/2023

Quant Time: Sep 09 04:58:38 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090823W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 09 03:56:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	131217	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	219697	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	199765	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	88062	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	75295	42.509	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	85.020%		
35) Dibromofluoromethane	5.385	113	66648	43.372	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	86.740%		
50) Toluene-d8	8.647	98	259859	46.107	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	92.220%		
62) 4-Bromofluorobenzene	11.079	95	94040	44.272	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	88.540%		
Target Compounds						
16) Acetone	2.386	43	18576	25.502	ug/l #	83
20) Methylene Chloride	2.788	84	11663	6.076	ug/l #	83
25) 2-Butanone	4.574	43	5669	5.132	ug/l #	84
37) Ethyl Acetate	4.745	43	2984m	1.434	ug/l	
43) Isopropyl Acetate	6.348	43	92255	29.592	ug/l #	88

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

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