

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090823\
 Data File : VX037610.D
 Acq On : 08 Sep 2023 18:38
 Operator : JC/MD
 Sample : PB155394ZHE#02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB155394ZHE#02

Manual Integrations
 APPROVED

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

Quant Time: Sep 09 04:58:09 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	5.556	168	138748	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	228616	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	205361	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	86156	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	77548	41.405	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	82.800%		
35) Dibromofluoromethane	5.391	113	69023	43.166	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	86.340%		
50) Toluene-d8	8.646	98	270877	46.187	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	92.380%		
62) 4-Bromofluorobenzene	11.079	95	94417	42.716	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	85.440%		
Target Compounds						
16) Acetone	2.385	43	19489	25.303	ug/l #	84
20) Methylene Chloride	2.788	84	11171	5.410	ug/l	88
37) Ethyl Acetate	4.739	43	3299m	1.523	ug/l	
43) Isopropyl Acetate	6.342	43	95785	29.526	ug/l #	90

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

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