

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021825\
 Data File : VX044989.D
 Acq On : 18 Feb 2025 15:46
 Operator : JC/MD
 Sample : PB166736ZHE#02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB166736ZHE#02

Manual Integrations
 APPROVED

Reviewed By : John Carlone 02/19/2025
 Supervised By : Mahesh Dadoda 02/19/2025

Quant Time: Feb 19 06:13:59 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X021025W.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 11 03:41:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	84519	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	173182	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	159204	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	70222	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	70309	56.829	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 113.660%			
35) Dibromofluoromethane	5.379	113	56855	50.494	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.980%			
50) Toluene-d8	8.647	98	216663	50.887	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 101.780%			
62) 4-Bromofluorobenzene	11.079	95	75103	52.323	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 104.640%			
Target Compounds						
16) Acetone	2.380	43	13347	26.835	ug/l	95
18) Methyl Acetate	2.709	43	2596	1.654	ug/l #	89
20) Methylene Chloride	2.782	84	1229	1.009	ug/l	90
37) Ethyl Acetate	4.733	43	2666m	1.475	ug/l	
43) Isopropyl Acetate	6.336	43	100465	34.431	ug/l	99

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

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