

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022924\
 Data File : VX040536.D
 Acq On : 29 Feb 2024 14:50
 Operator : JC/MD
 Sample : P1641-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OF-1

Quant Time: Mar 01 03:48:12 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022624W.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 27 05:01:33 2024
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 03/01/2024
 Supervised By : Semsettin Yesilyurt 03/01/2024

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	74829	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	130812	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	119212	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	65161	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	62025	49.558	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	99.120%		
35) Dibromofluoromethane	5.385	113	43638	47.972	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.940%		
50) Toluene-d8	8.647	98	159380	48.269	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	96.540%		
62) 4-Bromofluorobenzene	11.079	95	64006	50.852	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	101.700%		
Target Compounds						
16) Acetone	2.374	43	12938	31.699	ug/l	98
25) 2-Butanone	4.574	43	3482m	4.896	ug/l	
51) 4-Methyl-2-Pentanone	8.574	43	7496	4.711	ug/l	100

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

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