

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061324\
 Data File : VX041856.D
 Acq On : 13 Jun 2024 16:05
 Operator : JC/MD
 Sample : PB161467ZHE#11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB161467ZHE#11

Quant Time: Jun 14 01:00:27 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 30 07:36:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	160245	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	252721	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	250084	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	116656	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	83224	42.346	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	84.700%
35) Dibromofluoromethane	5.385	113	77325	44.254	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	88.500%
50) Toluene-d8	8.647	98	298628	48.913	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.820%
62) 4-Bromofluorobenzene	11.079	95	108568	48.181	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.360%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	89819	102.500	ug/l	# 88
18) Methyl Acetate	2.709	43	7036	3.749	ug/l	99
20) Methylene Chloride	2.788	84	10514	5.978	ug/l	90
25) 2-Butanone	4.574	43	22686	18.572	ug/l	95
43) Isopropyl Acetate	6.348	43	125332	31.090	ug/l	96

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

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