

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061324\
 Data File : VX041858.D
 Acq On : 13 Jun 2024 16:51
 Operator : JC/MD
 Sample : PB161467ZHE#13
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB161467ZHE#13

Quant Time: Jun 14 01:49:50 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	5.556	168	172707	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	266878	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	262095	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	127641	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	88773	41.910	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	83.820%
35) Dibromofluoromethane	5.385	113	83111	45.042	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.080%
50) Toluene-d8	8.647	98	315839	48.988	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.980%
62) 4-Bromofluorobenzene	11.079	95	116201	48.833	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.660%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	88725	93.624	ug/l	# 87
18) Methyl Acetate	2.715	43	7145	3.533	ug/l	96
20) Methylene Chloride	2.788	84	9581	5.054	ug/l	# 89
25) 2-Butanone	4.580	43	21669	16.460	ug/l	96
37) Ethyl Acetate	4.751	43	3320	1.177	ug/l	# 89
43) Isopropyl Acetate	6.348	43	120641	28.339	ug/l	# 95

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

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