

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX011224\
 Data File : VX039740.D
 Acq On : 12 Jan 2024 16:23
 Operator : JC/MD
 Sample : P1117-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 S1-TCLP-VOC-GRAB-7

Quant Time: Jan 12 22:40:58 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jan 11 03:51:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	96123	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	184387	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	178511	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	78301	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	73394	45.581	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	91.160%
35) Dibromofluoromethane	5.385	113	57771	46.517	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.040%
50) Toluene-d8	8.647	98	228094	48.696	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.400%
62) 4-Bromofluorobenzene	11.079	95	89719	48.982	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.960%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	20840	23.013	ug/l	93
18) Methyl Acetate	2.703	43	9965	5.260	ug/l	100
20) Methylene Chloride	2.788	84	4663	3.404	ug/l	92
37) Ethyl Acetate	4.733	43	2690	1.043	ug/l #	89
43) Isopropyl Acetate	6.342	43	89101	23.552	ug/l	100

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

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