

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU112221\
 Data File : VU045939.D
 Acq On : 23 Nov 2021 01:45
 Operator : SY/MD
 Sample : M4775-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PK-10I-20211118

Quant Time: Nov 23 06:51:12 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U110521W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 05 13:35:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.379	168	117298	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	212060	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	222801	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	110973	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	89943	51.201	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	102.400%
35) Dibromofluoromethane	5.292	113	71885	50.773	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.540%
50) Toluene-d8	7.899	98	280001	53.185	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	106.360%
62) 4-Bromofluorobenzene	10.632	95	107718	51.024	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	102.040%
Target Compounds						
						Qvalue
16) Acetone	2.668	43	1742	1.948	ug/l	98
19) Methyl tert-butyl Ether	3.366	73	3131	0.674	ug/l	97
65) Chlorobenzene	9.449	112	11408	2.607	ug/l	99
88) 1,4-Dichlorobenzene	11.838	146	2713	0.755	ug/l #	63
91) 1,2-Dichlorobenzene	12.211	146	1019	0.295	ug/l	92

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

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