

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044151.D
 Acq On : 10 Jun 2021 13:26
 Operator : SY/MD
 Sample : M2626-12
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12(XX)

Quant Time: Jun 11 01:19:46 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U060921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 10 05:11:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	142567	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	297873	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.424	117	305413	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	143627	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	130641	57.623	ug/l	0.00
Spiked Amount	50.000		Recovery	=	115.240%	
35) Dibromofluoromethane	5.298	113	100604	48.552	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.100%	
50) Toluene-d8	7.906	98	400057	50.272	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.540%	
62) 4-Bromofluorobenzene	10.639	95	151562	47.525	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.060%	
Target Compounds						
						Qvalue
16) Acetone	2.633	43	11095	7.400	ug/l	99
20) Methylene Chloride	3.054	84	14646	6.116	ug/l	99
43) Isopropyl Acetate	5.919	43	17656	3.136	ug/l	98
51) 4-Methyl-2-Pentanone	7.800	43	4561	1.119	ug/l	94

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

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