

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044147.D
 Acq On : 10 Jun 2021 11:51
 Operator : SY/MD
 Sample : M2626-08
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12(222-226)

Quant Time: Jun 11 01:19:08 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	140188	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	290391	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	299799	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	142985	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	127341	57.120	ug/l	0.00
Spiked Amount	50.000		Recovery	=	114.240%	
35) Dibromofluoromethane	5.298	113	99160	49.088	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.180%	
50) Toluene-d8	7.906	98	393777	50.758	ug/l	0.00
Spiked Amount	50.000		Recovery	=	101.520%	
62) 4-Bromofluorobenzene	10.639	95	153010	49.215	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.440%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	11493	7.911	ug/l	100
20) Methylene Chloride	3.054	84	13847	5.833	ug/l	99
25) 2-Butanone	4.723	43	2278	1.224	ug/l	90
43) Isopropyl Acetate	5.919	43	18146	3.306	ug/l	98
51) 4-Methyl-2-Pentanone	7.800	43	5544	1.395	ug/l	94
95) Naphthalene	14.095	128	3096	3.126	ug/l	100

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

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