

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044150.D
 Acq On : 10 Jun 2021 13:02
 Operator : SY/MD
 Sample : M2626-11
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12(X)

Quant Time: Jun 11 01:19:37 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	140748	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	292548	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	300718	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	141885	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	127450	56.942	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.880%	
35) Dibromofluoromethane	5.298	113	97527	47.924	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.840%	
50) Toluene-d8	7.906	98	389215	49.800	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.600%	
62) 4-Bromofluorobenzene	10.639	95	151412	48.342	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.680%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	11573	7.940	ug/l	99
20) Methylene Chloride	3.051	84	14008	5.887	ug/l	94
43) Isopropyl Acetate	5.922	43	17662	3.194	ug/l	98
51) 4-Methyl-2-Pentanone	7.803	43	4913	1.228	ug/l	92

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

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