

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044155.D
 Acq On : 10 Jun 2021 15:00
 Operator : SY/MD
 Sample : M2637-09
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12(242-246)

Quant Time: Jun 11 01:20:25 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	139146	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	286135	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	291318	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	131944	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	123695	55.900	ug/l	0.00
Spiked Amount	50.000		Recovery	=	111.800%	
35) Dibromofluoromethane	5.298	113	98154	49.313	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.620%	
50) Toluene-d8	7.906	98	388305	50.797	ug/l	0.00
Spiked Amount	50.000		Recovery	=	101.600%	
62) 4-Bromofluorobenzene	10.639	95	140774	45.953	ug/l	0.00
Spiked Amount	50.000		Recovery	=	91.900%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	11347	7.857	ug/l	99
20) Methylene Chloride	3.054	84	12250	5.066	ug/l	98
43) Isopropyl Acetate	5.919	43	17035	3.150	ug/l	97
51) 4-Methyl-2-Pentanone	7.800	43	4150	1.060	ug/l	97

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

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