

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044154.D
 Acq On : 10 Jun 2021 14:36
 Operator : SY/MD
 Sample : M2637-08
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12(238-242)

Quant Time: Jun 11 01:20:16 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	141849	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	296737	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	304118	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	143440	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	128879	57.133	ug/l	0.00
Spiked Amount	50.000		Recovery	=	114.260%	
35) Dibromofluoromethane	5.298	113	100539	48.706	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.420%	
50) Toluene-d8	7.906	98	395932	49.944	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.880%	
62) 4-Bromofluorobenzene	10.639	95	150100	47.247	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.500%	
Target Compounds						
					Qvalue	
16) Acetone	2.632	43	10951	7.323	ug/l	95
20) Methylene Chloride	3.054	84	10741	4.185	ug/l	96
43) Isopropyl Acetate	5.919	43	16365	2.918	ug/l	99
51) 4-Methyl-2-Pentanone	7.800	43	4729	1.165	ug/l	92

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

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