

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044148.D
 Acq On : 10 Jun 2021 12:15
 Operator : SY/MD
 Sample : M2626-09
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12(226-230)

Quant Time: Jun 11 01:19:19 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	143620	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	297825	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	302092	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	143558	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	130654	57.206	ug/l	0.00
Spiked Amount	50.000		Recovery	=	114.420%	
35) Dibromofluoromethane	5.298	113	100289	48.408	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.820%	
50) Toluene-d8	7.906	98	399782	50.246	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.500%	
62) 4-Bromofluorobenzene	10.639	95	152511	47.830	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.660%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	11477	7.656	ug/l	96
20) Methylene Chloride	3.054	84	13408	5.446	ug/l	98
43) Isopropyl Acetate	5.919	43	18178	3.229	ug/l	98
51) 4-Methyl-2-Pentanone	7.800	43	4629	1.136	ug/l	99

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

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