

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060821\
 Data File : VU044110.D
 Acq On : 08 Jun 2021 13:05
 Operator : SY/MD
 Sample : M2612-33
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12(54-58)

Quant Time: Jun 08 23:12:55 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U060721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 08 17:38:51 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.382	168	206250	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	412557	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	433410	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	188172	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	189872	52.860	ug/l	0.00
Spiked Amount	50.000		Recovery	= 105.720%		
35) Dibromofluoromethane	5.298	113	148774	50.533	ug/l	0.00
Spiked Amount	50.000		Recovery	= 101.060%		
50) Toluene-d8	7.906	98	583024	51.341	ug/l	0.00
Spiked Amount	50.000		Recovery	= 102.680%		
62) 4-Bromofluorobenzene	10.639	95	201637	44.848	ug/l	0.00
Spiked Amount	50.000		Recovery	= 89.700%		
Target Compounds						
						Qvalue
16) Acetone	2.633	43	19357	9.337	ug/l	96
20) Methylene Chloride	3.054	84	8775	1.163	ug/l	96
43) Isopropyl Acetate	5.919	43	26328	3.180	ug/l	99

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

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