

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044169.D
 Acq On : 10 Jun 2021 20:28
 Operator : SY/MD
 Sample : M2658-17
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 G-WALL-2

Quant Time: Jun 11 01:22:36 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	5.385	168	145162	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.259	114	284491	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	304183	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	141855	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	125004	54.151	ug/l	0.00
Spiked Amount	50.000		Recovery	= 108.300%		
35) Dibromofluoromethane	5.298	113	100624	50.846	ug/l	0.00
Spiked Amount	50.000		Recovery	= 101.700%		
50) Toluene-d8	7.906	98	395040	51.977	ug/l	0.00
Spiked Amount	50.000		Recovery	= 103.960%		
62) 4-Bromofluorobenzene	10.639	95	143350	47.065	ug/l	0.00
Spiked Amount	50.000		Recovery	= 94.120%		
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	30398	23.564	ug/l	99
20) Methylene Chloride	3.054	84	13953	5.644	ug/l	99
43) Isopropyl Acetate	5.919	43	14426	2.683	ug/l	98
51) 4-Methyl-2-Pentanone	7.800	43	5826	1.497	ug/l	95

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

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