

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044161.D
 Acq On : 10 Jun 2021 17:21
 Operator : SY/MD
 Sample : M2651-09
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 17-B6(0-4)

Quant Time: Jun 11 01:21:20 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	141291	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	284430	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	300744	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	140758	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	124872	55.576	ug/l	0.00
Spiked Amount	50.000		Recovery	=	111.160%	
35) Dibromofluoromethane	5.298	113	99277	50.176	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.360%	
50) Toluene-d8	7.906	98	391608	51.536	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.080%	
62) 4-Bromofluorobenzene	10.639	95	144842	47.565	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.120%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	11683	7.997	ug/l	96
20) Methylene Chloride	3.054	84	11492	4.587	ug/l	97
43) Isopropyl Acetate	5.919	43	16093	2.993	ug/l	99
51) 4-Methyl-2-Pentanone	7.796	43	4418	1.135	ug/l	92

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

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