

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044159.D
 Acq On : 10 Jun 2021 16:34
 Operator : SY/MD
 Sample : M2651-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 17-B2(0-5)

Quant Time: Jun 11 01:21:01 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	135680	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	274714	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	297587	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	144757	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	120170	55.695	ug/l	0.00
Spiked Amount	50.000		Recovery	=	111.380%	
35) Dibromofluoromethane	5.298	113	94730	49.571	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.140%	
50) Toluene-d8	7.906	98	377034	51.373	ug/l	0.00
Spiked Amount	50.000		Recovery	=	102.740%	
62) 4-Bromofluorobenzene	10.639	95	147928	50.296	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.600%	
Target Compounds						
						Qvalue
16) Acetone	2.636	43	16252	12.554	ug/l	99
20) Methylene Chloride	3.050	84	12198	5.199	ug/l	98
43) Isopropyl Acetate	5.919	43	14955	2.880	ug/l	96
51) 4-Methyl-2-Pentanone	7.799	43	4450	1.184	ug/l	90

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

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