

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044158.D
 Acq On : 10 Jun 2021 16:10
 Operator : SY/MD
 Sample : M2651-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 CON-ED-EXCAVATION

Quant Time: Jun 11 01:20:52 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	144016	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	298161	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	309159	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	145049	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	129261	56.440	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.880%	
35) Dibromofluoromethane	5.298	113	101988	49.172	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.340%	
50) Toluene-d8	7.906	98	402458	50.525	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.060%	
62) 4-Bromofluorobenzene	10.639	95	151958	47.603	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.200%	
Target Compounds					Qvalue	
16) Acetone	2.633	43	23551	17.928	ug/l	99
20) Methylene Chloride	3.054	84	12837	5.144	ug/l	98
43) Isopropyl Acetate	5.919	43	16660	2.956	ug/l	98
51) 4-Methyl-2-Pentanone	7.800	43	4376	1.073	ug/l	93

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

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