

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044167.D
 Acq On : 10 Jun 2021 19:41
 Operator : SY/MD
 Sample : M2658-11
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 G-FLOOR-2

Quant Time: Jun 11 01:22:16 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	142601	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	280009	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	294027	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	152810	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	120981	53.349	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.700%	
35) Dibromofluoromethane	5.298	113	89417	45.906	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.820%	
50) Toluene-d8	7.906	98	372945	49.855	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.720%	
62) 4-Bromofluorobenzene	10.639	95	150484	50.198	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.400%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	43961	35.709	ug/l	99
20) Methylene Chloride	3.051	84	13865	5.723	ug/l	99
29) Tetrahydrofuran	5.067	42	1148	1.116	ug/l #	85
51) 4-Methyl-2-Pentanone	7.800	43	4402	1.149	ug/l	100

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

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