

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061421\
 Data File : VU044212.D
 Acq On : 15 Jun 2021 03:57
 Operator : SY/MD
 Sample : M2687-18
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DUP-01

Manual Integrations
 APPROVED

MMDadoda
 6/15/2021 9:51:01 AM

Quant Time: Jun 15 05:01:46 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	174559	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	318118	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	310287	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	147426	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.713	65	132363	47.682	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.360%	
35) Dibromofluoromethane	5.301	113	103483	46.763	ug/l	0.00
Spiked Amount	50.000		Recovery	=	93.520%	
50) Toluene-d8	7.906	98	407781	47.982	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.960%	
62) 4-Bromofluorobenzene	10.639	95	143895	42.250	ug/l	0.00
Spiked Amount	50.000		Recovery	=	84.500%	
Target Compounds						
16) Acetone	2.655	43	17680m	10.282	ug/l	
20) Methylene Chloride	3.057	84	15233	5.011	ug/l	94
43) Isopropyl Acetate	5.922	43	18866	3.137	ug/l	98

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

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