

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061421\
 Data File : VU044207.D
 Acq On : 15 Jun 2021 02:00
 Operator : SY/MD
 Sample : M2687-13
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 17-B6(108-112)

Manual Integrations
 APPROVED

MMDadoda
 6/15/2021 9:50:56 AM

Quant Time: Jun 15 05:00:22 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	172971	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.259	114	318074	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	309862	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	148155	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.713	65	133536	48.547	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.100%	
35) Dibromofluoromethane	5.298	113	103516	46.785	ug/l	0.00
Spiked Amount	50.000		Recovery	=	93.560%	
50) Toluene-d8	7.906	98	408172	48.035	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.060%	
62) 4-Bromofluorobenzene	10.639	95	144355	42.390	ug/l	0.00
Spiked Amount	50.000		Recovery	=	84.780%	
Target Compounds						
16) Acetone	2.652	43	11837m	6.246	ug/l	
20) Methylene Chloride	3.057	84	13253	4.249	ug/l	96
43) Isopropyl Acetate	5.922	43	18773	3.122	ug/l	96

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061421\
Data File : VU044207.D
Acq On : 15 Jun 2021 02:00
Operator : SY/MD
Sample : M2687-13
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
17-B6(108-112)

Manual Integrations
APPROVED

MMDadoda
6/15/2021 9:50:56 AM

Quant Time: Jun 15 05:00:22 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

