

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044149.D
 Acq On : 10 Jun 2021 12:39
 Operator : SY/MD
 Sample : M2626-10
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12-DUP-01

Quant Time: Jun 11 01:19:28 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	140676	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	290608	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	300620	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	142499	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	127337	56.920	ug/l	0.00
Spiked Amount	50.000		Recovery	=	113.840%	
35) Dibromofluoromethane	5.298	113	99799	49.368	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.740%	
50) Toluene-d8	7.906	98	393093	50.632	ug/l	0.00
Spiked Amount	50.000		Recovery	=	101.260%	
62) 4-Bromofluorobenzene	10.639	95	151231	48.607	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.220%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	12246	8.533	ug/l	100
20) Methylene Chloride	3.054	84	14022	5.898	ug/l	96
43) Isopropyl Acetate	5.919	43	17864	3.252	ug/l	98
51) 4-Methyl-2-Pentanone	7.800	43	5260	1.323	ug/l	87

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
Data File : VU044149.D
Acq On : 10 Jun 2021 12:39
Operator : SY/MD
Sample : M2626-10
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
18-B12-DUP-01

Quant Time: Jun 11 01:19:28 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

