

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060821\
 Data File : VU044092.D
 Acq On : 08 Jun 2021 06:02
 Operator : SY/MD
 Sample : M2613-15
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 18-B12(90-94)

Quant Time: Jun 08 07:15:16 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U060721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 08 05:21:34 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.382	168	216023	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	425279	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	448107	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	207086	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	194228	51.626	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.260%	
35) Dibromofluoromethane	5.298	113	149089	49.125	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.260%	
50) Toluene-d8	7.906	98	587707	50.205	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.400%	
62) 4-Bromofluorobenzene	10.639	95	218970	47.246	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.500%	
Target Compounds						
						Qvalue
16) Acetone	2.636	43	14756	6.796	ug/l	90
20) Methylene Chloride	3.054	84	20339	4.707	ug/l	95
43) Isopropyl Acetate	5.919	43	28005	3.281	ug/l	99
95) Naphthalene	14.095	128	4897	3.173	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060821\
Data File : VU044092.D
Acq On : 08 Jun 2021 06:02
Operator : SY/MD
Sample : M2613-15
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
18-B12(90-94)

Quant Time: Jun 08 07:15:16 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U060721W.M
Quant Title : SW846 8260
QLast Update : Tue Jun 08 05:21:34 2021
Response via : Initial Calibration

