

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102521\
 Data File : VU045393.D
 Acq On : 25 Oct 2021 19:03
 Operator : SY/MD
 Sample : M4300-11
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TH-CCD

Quant Time: Oct 26 05:53:40 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 28 12:32:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.378	168	211600	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	406845	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	371027	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	165753	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	173244	49.184	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.360%
35) Dibromofluoromethane	5.295	113	126581	49.320	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.640%
50) Toluene-d8	7.902	98	502914	49.594	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.180%
62) 4-Bromofluorobenzene	10.636	95	178675	42.677	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	85.360%
Target Compounds						
					Qvalue	
16) Acetone	2.636	43	12774	6.110	ug/l	99
20) Methylene Chloride	3.051	84	11930	3.788	ug/l	91
43) Isopropyl Acetate	5.915	43	27384	3.233	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102521\
Data File : VU045393.D
Acq On : 25 Oct 2021 19:03
Operator : SY/MD
Sample : M4300-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
TH-CCD

Quant Time: Oct 26 05:53:40 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 28 12:32:13 2021
Response via : Initial Calibration

