

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111021\
 Data File : VU045690.D
 Acq On : 10 Nov 2021 16:08
 Operator : SY/MD
 Sample : M4497-08
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 346

Quant Time: Nov 10 16:38:40 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U110521W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 05 13:35:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.378	168	124787	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	239141	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	242446	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	118252	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	102562	54.880	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 109.760%	
35) Dibromofluoromethane	5.295	113	80916	50.680	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.360%	
50) Toluene-d8	7.899	98	312015	52.554	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 105.100%	
62) 4-Bromofluorobenzene	10.635	95	118926	49.954	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 99.900%	
Target Compounds						
						Qvalue
16) Acetone	2.626	43	13847	14.555	ug/l	94
20) Methylene Chloride	3.047	84	6574	3.977	ug/l	98
43) Isopropyl Acetate	5.912	43	18948	4.625	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111021\
Data File : VU045690.D
Acq On : 10 Nov 2021 16:08
Operator : SY/MD
Sample : M4497-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
346

Quant Time: Nov 10 16:38:40 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U110521W.M
Quant Title : SW846 8260
QLast Update : Fri Nov 05 13:35:16 2021
Response via : Initial Calibration

