

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110421\
 Data File : VU045542.D
 Acq On : 03 Nov 2021 23:59
 Operator : SY/MD
 Sample : M4470-05
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SP-2

Quant Time: Nov 05 13:57:50 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U102821W.M
 Quant Title : SW846 8260
 QLast Update : Fri Oct 29 11:09:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.378	168	174699	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	345160	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	352936	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	177087	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	112588	39.586	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	79.180%
35) Dibromofluoromethane	5.295	113	102217	46.575	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.140%
50) Toluene-d8	7.902	98	418178	48.590	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.180%
62) 4-Bromofluorobenzene	10.635	95	181613	55.378	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.760%
Target Compounds						
					Qvalue	
16) Acetone	2.649	43	14529	11.252	ug/l #	83
20) Methylene Chloride	3.051	84	8094	3.146	ug/l	91
43) Isopropyl Acetate	5.919	43	19526	3.256	ug/l	96

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110421\
Data File : VU045542.D
Acq On : 03 Nov 2021 23:59
Operator : SY/MD
Sample : M4470-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SP-2

Quant Time: Nov 05 13:57:50 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U102821W.M
Quant Title : SW846 8260
QLast Update : Fri Oct 29 11:09:42 2021
Response via : Initial Calibration

