

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111221\
 Data File : VU045777.D
 Acq On : 12 Nov 2021 18:16
 Operator : SY/MD
 Sample : PB140714TB
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB140714TB

Quant Time: Nov 15 05:23:08 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	5.379	168	115609	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	217995	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	224857	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	109494	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	93009	53.720	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 107.440%	
35) Dibromofluoromethane	5.292	113	73236	50.319	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.640%	
50) Toluene-d8	7.903	98	289885	53.563	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 107.120%	
62) 4-Bromofluorobenzene	10.636	95	107121	49.360	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 98.720%	
Target Compounds						
						Qvalue
16) Acetone	2.626	43	16614	18.850	ug/l	95
20) Methylene Chloride	3.047	84	7717	5.039	ug/l	96
43) Isopropyl Acetate	5.916	43	17307	4.634	ug/l	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111221\
Data File : VU045777.D
Acq On : 12 Nov 2021 18:16
Operator : SY/MD
Sample : PB140714TB
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
PB140714TB

Quant Time: Nov 15 05:23:08 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

