

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060722\
 Data File : VU048909.D
 Acq On : 07 Jun 2022 15:02
 Operator : SY/MD
 Sample : N3165-11
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 64BR-20220531

Quant Time: Jun 08 04:52:37 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
 Quant Title : SW846 8260
 QLast Update : Fri May 27 06:42:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.372	168	254325	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	417746	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	388894	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	174336	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	190878	49.975	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	99.940%		
35) Dibromofluoromethane	5.288	113	143949	48.300	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	96.600%		
50) Toluene-d8	7.899	98	511355	47.865	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	95.720%		
62) 4-Bromofluorobenzene	10.632	95	185497	45.911	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	91.820%		
Target Compounds						Qvalue
16) Acetone	2.626	43	5128	2.634	ug/l	97
27) cis-1,2-Dichloroethene	4.668	96	6494	1.895	ug/l	82
37) Ethyl Acetate	4.803	43	78540	14.283	ug/l	99
44) Trichloroethene	6.542	130	2029	0.613	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060722\
Data File : VU048909.D
Acq On : 07 Jun 2022 15:02
Operator : SY/MD
Sample : N3165-11
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
64BR-20220531

Quant Time: Jun 08 04:52:37 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
Quant Title : SW846 8260
QLast Update : Fri May 27 06:42:31 2022
Response via : Initial Calibration

