

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU052722\
 Data File : VU048696.D
 Acq On : 27 May 2022 16:15
 Operator : SY/MD
 Sample : N3056-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-11

Quant Time: May 28 00:20:30 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	5.375	168	204128	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	335412	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	352182	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	198876	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	161239	52.596	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	105.200%
35) Dibromofluoromethane	5.292	113	138260	57.779	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	115.560%
50) Toluene-d8	7.899	98	396534	46.228	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	92.460%
62) 4-Bromofluorobenzene	10.632	95	186024	57.344	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	114.680%
Target Compounds						
					Qvalue	
16) Acetone	2.629	43	3069	1.964	ug/l	91
19) Methyl tert-butyl Ether	3.369	73	5990	0.760	ug/l	99
30) Chloroform	5.095	83	2351	0.504	ug/l	100
84) 1,2,4-Trimethylbenzene	11.471	105	1544	1.312	ug/l	95

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU052722\
Data File : VU048696.D
Acq On : 27 May 2022 16:15
Operator : SY/MD
Sample : N3056-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-11

Quant Time: May 28 00:20:30 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

