

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060222\
 Data File : VU048811.D
 Acq On : 02 Jun 2022 18:09
 Operator : SY/MD
 Sample : N3003-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RB04-20220523

Quant Time: Jun 03 01:17:57 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	363792	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	588556	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	534423	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	249441	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	236622	43.310	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	86.620%
35) Dibromofluoromethane	5.288	113	187531	44.662	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.320%
50) Toluene-d8	7.899	98	718973	47.767	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.540%
62) 4-Bromofluorobenzene	10.632	95	245487	43.126	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	86.260%
Target Compounds						
					Qvalue	
16) Acetone	2.626	43	23609	8.479	ug/l	95
20) Methylene Chloride	3.044	84	3987	0.842	ug/l	96
25) 2-Butanone	4.719	43	7771	1.943	ug/l	92
95) Naphthalene	14.089	128	28048	3.889	ug/l	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060222\
Data File : VU048811.D
Acq On : 02 Jun 2022 18:09
Operator : SY/MD
Sample : N3003-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
RB04-20220523

Quant Time: Jun 03 01:17:57 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

