

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040821\
 Data File : VN066540.D
 Acq On : 8 Apr 2021 14:50
 Operator : JC/MD
 Sample : M1961-10
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TR1-D-TOP

Quant Time: Apr 09 04:34:38 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N040621W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 06 15:38:34 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.587	168	312802	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.515	114	478176	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.350	117	476305	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.289	152	194577	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.949	65	203521	49.41	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.82%
35) Dibromofluoromethane	7.507	113	159524	48.95	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.90%
50) Toluene-d8	10.028	98	616849	49.81	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.62%
62) 4-Bromofluorobenzene	12.348	95	193931	48.29	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.58%
Target Compounds						
					Qvalue	
16) Acetone	3.736	43	23505	13.61	ug/l	99
20) Methylene Chloride	4.460	84	18316	4.94	ug/l	94
43) Isopropyl Acetate	8.094	43	26943	3.03	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040821\
Data File : VN066540.D
Acq On : 8 Apr 2021 14:50
Operator : JC/MD
Sample : M1961-10
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TR1-D-TOP

Quant Time: Apr 09 04:34:38 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N040621W.M
Quant Title : SW846 8260
QLast Update : Tue Apr 06 15:38:34 2021
Response via : Initial Calibration

