

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040621\
 Data File : VN066503.D
 Acq On : 6 Apr 2021 19:49
 Operator : JC/MD
 Sample : M1934-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-2

Quant Time: Apr 07 04:49:37 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	7.585	168	309149	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.515	114	462106	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.350	117	461800	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.289	152	201741	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.947	65	193763	47.59	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	95.18%
35) Dibromofluoromethane	7.504	113	154726	49.13	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.26%
50) Toluene-d8	10.028	98	596998	49.89	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.78%
62) 4-Bromofluorobenzene	12.345	95	190195	49.01	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.02%
Target Compounds						
					Qvalue	
16) Acetone	3.739	43	22786	13.35	ug/l	98
20) Methylene Chloride	4.452	84	12029	3.28	ug/l	90
43) Isopropyl Acetate	8.086	43	29442	3.42	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040621\
Data File : VN066503.D
Acq On : 6 Apr 2021 19:49
Operator : JC/MD
Sample : M1934-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-2

Quant Time: Apr 07 04:49:37 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

