

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100923\
 Data File : VN079467.D
 Acq On : 09 Oct 2023 13:11
 Operator : JC\MD
 Sample : 04731-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DTP-2

Quant Time: Oct 09 23:13:57 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100523W.M
 Quant Title : SW846 8260
 QLast Update : Fri Oct 06 03:34:49 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	283919	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	567367	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	565215	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	177774	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	238843	51.883	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	103.760%
35) Dibromofluoromethane	8.171	113	202588	49.687	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.380%
50) Toluene-d8	10.571	98	698418	47.341	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.680%
62) 4-Bromofluorobenzene	12.853	95	236107	49.869	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.740%
Target Compounds						
						Qvalue
16) Acetone	4.442	43	24220	13.265	ug/l	91
20) Methylene Chloride	5.271	84	13900	2.909	ug/l #	91
43) Isopropyl Acetate	8.689	43	203104	19.039	ug/l #	79

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100923\
Data File : VN079467.D
Acq On : 09 Oct 2023 13:11
Operator : JC\MD
Sample : 04731-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DTP-2

Quant Time: Oct 09 23:13:57 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100523W.M
Quant Title : SW846 8260
QLast Update : Fri Oct 06 03:34:49 2023
Response via : Initial Calibration

