

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020823\
 Data File : VN076584.D
 Acq On : 08 Feb 2023 16:58
 Operator : JC\MD
 Sample : 01418-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 JPP-20-020223

Quant Time: Feb 09 00:51:17 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020723W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 08 09:04:43 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.233	168	706095	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.104	114	1180993	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.869	117	1048867	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.792	152	417463	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.586	65	394101	47.288	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	94.580%
35) Dibromofluoromethane	8.175	113	352063	47.956	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.920%
50) Toluene-d8	10.569	98	1383666	50.099	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	100.200%
62) 4-Bromofluorobenzene	12.851	95	466487	47.944	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.880%
Target Compounds						
					Qvalue	
16) Acetone	4.451	43	79704	29.008	ug/l	95
20) Methylene Chloride	5.293	84	29007	3.425	ug/l	97
37) Ethyl Acetate	7.575	43	61511	7.833	ug/l	98
43) Isopropyl Acetate	8.698	43	317033	24.261	ug/l #	87

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020823\
Data File : VN076584.D
Acq On : 08 Feb 2023 16:58
Operator : JC\MD
Sample : 01418-10
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
JPP-20-020223

Quant Time: Feb 09 00:51:17 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020723W.M
Quant Title : SW846 8260
QLast Update : Wed Feb 08 09:04:43 2023
Response via : Initial Calibration

