

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
 Data File : VN082644.D
 Acq On : 19 Jun 2024 16:25
 Operator : JC\MD
 Sample : P2918-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 POLE

Quant Time: Jun 19 16:42:49 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061724W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 18 04:52:44 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	131187	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	253057	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	263288	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	104312	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	121847	57.919	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	115.840%
35) Dibromofluoromethane	8.171	113	90950	52.979	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	105.960%
50) Toluene-d8	10.571	98	344299	53.034	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	106.060%
62) 4-Bromofluorobenzene	12.853	95	110624	49.425	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	=	98.860%
Target Compounds						
						Qvalue
16) Acetone	4.442	43	21626	22.603	ug/l	96
20) Methylene Chloride	5.277	84	10242	6.089	ug/l	90
25) 2-Butanone	7.500	43	12843	9.444	ug/l #	90
43) Isopropyl Acetate	8.694	43	271966	44.517	ug/l #	80

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
Data File : VN082644.D
Acq On : 19 Jun 2024 16:25
Operator : JC\MD
Sample : P2918-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
POLE

Quant Time: Jun 19 16:42:49 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061724W.M
Quant Title : SW846 8260
QLast Update : Tue Jun 18 04:52:44 2024
Response via : Initial Calibration

