

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
 Data File : VN082649.D
 Acq On : 19 Jun 2024 18:28
 Operator : JC\MD
 Sample : P2955-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VK-1

Quant Time: Jun 20 03:05:04 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	129962	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	252583	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	263605	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	104046	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	120667	57.898	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 115.800%			
35) Dibromofluoromethane	8.171	113	88964	51.919	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 103.840%			
50) Toluene-d8	10.571	98	343080	52.946	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 105.900%			
62) 4-Bromofluorobenzene	12.853	95	108242	48.452	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 96.900%			
Target Compounds						
16) Acetone	4.430	43	35157	37.092	ug/l	100
20) Methylene Chloride	5.277	84	6369	3.822	ug/l #	83
43) Isopropyl Acetate	8.694	43	206810	33.916	ug/l #	83

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
Data File : VN082649.D
Acq On : 19 Jun 2024 18:28
Operator : JC\MD
Sample : P2955-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VK-1

Quant Time: Jun 20 03:05:04 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

