

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
 Data File : VN082645.D
 Acq On : 19 Jun 2024 16:50
 Operator : JC\MD
 Sample : P2916-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPF

Quant Time: Jun 20 03:03:08 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	8.229	168	130640	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	249229	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	258191	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	100794	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	118520	56.573	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	113.140%
35) Dibromofluoromethane	8.171	113	88510	52.350	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	104.700%
50) Toluene-d8	10.570	98	332025	51.929	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.860%
62) 4-Bromofluorobenzene	12.853	95	106610	48.363	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	=	96.720%
Target Compounds						
						Qvalue
16) Acetone	4.436	43	20479	21.494	ug/l	93
20) Methylene Chloride	5.277	84	10882	6.496	ug/l	95
25) 2-Butanone	7.494	43	12677	9.361	ug/l	95
43) Isopropyl Acetate	8.694	43	254737	42.338	ug/l #	82

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
Data File : VN082645.D
Acq On : 19 Jun 2024 16:50
Operator : JC\MD
Sample : P2916-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WCS-TPF

Quant Time: Jun 20 03:03:08 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

