

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061724\
 Data File : VN082617.D
 Acq On : 17 Jun 2024 18:38
 Operator : JC\MD
 Sample : P2889-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPJ

Quant Time: Jun 18 05:02:04 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	8.224	168	130772	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	247442	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	254092	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	103301	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	117675	56.113	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 112.220%		
35) Dibromofluoromethane	8.171	113	85461	50.911	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.820%		
50) Toluene-d8	10.571	98	329640	51.929	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 103.860%		
62) 4-Bromofluorobenzene	12.853	95	107957	49.328	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	= 98.660%		
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	20963	21.980	ug/l	98
20) Methylene Chloride	5.283	84	8133	4.850	ug/l	91
25) 2-Butanone	7.495	43	10749	7.929	ug/l	96
43) Isopropyl Acetate	8.694	43	261341m	43.749	ug/l	

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061724\
Data File : VN082617.D
Acq On : 17 Jun 2024 18:38
Operator : JC\MD
Sample : P2889-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WCS-TPJ

Quant Time: Jun 18 05:02:04 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

