

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN012920\
 Data File : VN059913.D
 Acq On : 29 Jan 2020 15:33
 Operator : JC/MD
 Sample : L1289-38
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-3-COMP

Quant Time: Jan 30 01:07:12 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jan 18 03:09:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	206394	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	311713	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	284785	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	118285	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	110994	48.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.86%	
35) Dibromofluoromethane	7.57	113	94568	50.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.74%	
50) Toluene-d8	10.08	98	371219	51.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.08%	
62) 4-Bromofluorobenzene	12.40	95	125165	48.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.42%	
Target Compounds						
18) Methyl Acetate	4.30	43	7237	2.703	ug/l	95
20) Methylene Chloride	4.52	84	8822	3.039	ug/l	92
43) Isopropyl Acetate	8.14	43	38093	8.699	ug/l #	92
95) Naphthalene	15.13	128	18255	2.831	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN012920\
Data File : VN059913.D
Acq On : 29 Jan 2020 15:33
Operator : JC/MD
Sample : L1289-38
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-3-COMP

Quant Time: Jan 30 01:07:12 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
Quant Title : SW846 8260
QLast Update : Sat Jan 18 03:09:32 2020
Response via : Initial Calibration

