

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091620\
 Data File : VN063404.D
 Acq On : 17 Sep 2020 7:51
 Operator : JC/MD
 Sample : L4019-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VNJ-220

Quant Time: Sep 17 08:12:33 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.62	168	228832	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	436260	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	430506	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	186669	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	195745	50.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.50%	
35) Dibromofluoromethane	7.55	113	140618	46.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.12%	
50) Toluene-d8	10.06	98	566134	49.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.70%	
62) 4-Bromofluorobenzene	12.38	95	224223	53.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.72%	
Target Compounds						
16) Acetone	3.78	43	19281	15.401	ug/l	97
18) Methyl Acetate	4.28	43	4465	1.405	ug/l #	67
20) Methylene Chloride	4.51	84	12273	3.869	ug/l	96
43) Isopropyl Acetate	8.13	43	80059	9.815	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091620\
Data File : VN063404.D
Acq On : 17 Sep 2020 7:51
Operator : JC/MD
Sample : L4019-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 54 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VNJ-220

Quant Time: Sep 17 08:12:33 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 16 13:05:20 2020
Response via : Initial Calibration

