

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN110620\
 Data File : VN064492.D
 Acq On : 6 Nov 2020 18:51
 Operator : JC/MD
 Sample : L4651-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB-2020-WS-000-03

Quant Time: Nov 07 04:34:20 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 04 12:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	244977	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	529054	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	561044	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	203401	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	244537	57.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.62%	
35) Dibromofluoromethane	7.55	113	176948	49.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.24%	
50) Toluene-d8	10.06	98	680478	48.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.76%	
62) 4-Bromofluorobenzene	12.38	95	261708	49.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.42%	
Target Compounds						
16) Acetone	3.78	43	32091	19.743	ug/l	94
18) Methyl Acetate	4.29	43	6142	1.922	ug/l	98
20) Methylene Chloride	4.51	84	12739	3.950	ug/l	95
43) Isopropyl Acetate	8.13	43	95732	10.569	ug/l	97
44) Trichloroethene	8.80	130	6678	1.612	ug/l	89
64) Tetrachloroethene	10.60	164	61598	14.253	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN110620\
Data File : VN064492.D
Acq On : 6 Nov 2020 18:51
Operator : JC/MD
Sample : L4651-01
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB-2020-WS-000-03

Quant Time: Nov 07 04:34:20 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260
QLast Update : Wed Nov 04 12:09:03 2020
Response via : Initial Calibration

