

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092420\
 Data File : VN063637.D
 Acq On : 24 Sep 2020 18:40
 Operator : JC/MD
 Sample : L4119-14
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RE123D2-20200922

Quant Time: Sep 25 04:58:01 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	7.63	168	196415	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	372930	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	383228	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	159485	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	171040	51.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.32%	
35) Dibromofluoromethane	7.55	113	126552	49.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.04%	
50) Toluene-d8	10.06	98	478343	49.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.54%	
62) 4-Bromofluorobenzene	12.38	95	184471	51.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.68%	
Target Compounds						
3) Chloromethane	2.04	50	2849	0.968	ug/l	96
16) Acetone	3.79	43	6327	5.888	ug/l #	83
44) Trichloroethene	8.80	130	8222	2.989	ug/l	93
64) Tetrachloroethene	10.60	164	4461	1.710	ug/l	96

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092420\
Data File : VN063637.D
Acq On : 24 Sep 2020 18:40
Operator : JC/MD
Sample : L4119-14
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
RE123D2-20200922

Quant Time: Sep 25 04:58:01 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

