

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100918\
 Data File : VN051747.D
 Acq On : 9 Oct 2018 13:17
 Operator : MD\SY
 Sample : J5235-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 6 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 CLEARWELL

Quant Time: Oct 10 10:06:39 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA N\METHODS\624N100518W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Oct 05 11:11:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.20	128	112040	30.00	ug/l	0.00
28) 1,4-Difluorobenzene	8.59	114	639140	30.00	ug/l	0.00
57) Chlorobenzene-d5	11.41	117	527240	30.00	ug/l	0.00

System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.03	65	228180	29.34	ug/l	0.00
Spiked Amount	30.000	Range	50 - 169	Recovery	=	97.80%
60) 4-Bromofluorobenzene	12.40	95	201617	25.06	ug/l	0.00
Spiked Amount	30.000	Range	56 - 143	Recovery	=	83.53%
63) Toluene-d8	10.09	98	745733	31.86	ug/l	0.00
Spiked Amount	30.000	Range	66 - 137	Recovery	=	106.20%

Target Compounds					Ovalue
15) Acetone	3.82	58	6622	16.444 ug/l	88
25) Chloroform	7.38	83	37508	3.102 ug/l	99
41) Bromodichloromethane	9.40	83	9841	1.013 ug/l #	98
53) Dibromochloromethane	10.90	129	5689	0.705 ug/l	96

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100918\
Data File : VN051747.D
Acq On : 9 Oct 2018 13:17
Operator : MD\SY
Sample : J5235-09
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 6 Sample Multiplier: 28

Instrument :
MSVOA_N
ClientSampleId :
CLEARWELL

Quant Time: Oct 10 10:06:39 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA N\METHODS\624N100518W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
QLast Update : Fri Oct 05 11:11:26 2018
Response via : Initial Calibration

