

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100719\
 Data File : VN058585.D
 Acq On : 7 Oct 2019 22:18
 Operator : JC/SP
 Sample : K5215-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 13940

Quant Time: Oct 08 15:44:57 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 30 08:07:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	480239	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	743282	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	651288	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	271281	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	324139	46.19	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.38%	
35) Dibromofluoromethane	7.57	113	230545	50.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.22%	
50) Toluene-d8	10.08	98	936815	52.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.98%	
62) 4-Bromofluorobenzene	12.41	95	314231	48.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.40%	
Target Compounds						
11) Tert butyl alcohol	4.76	59	20355	26.108	ug/l	97
16) Acetone	3.80	43	33098	13.835	ug/l	98
19) Methyl tert-butyl Ether	5.02	73	78016	4.881	ug/l	97
22) Diisopropyl ether	5.93	45	80619	4.544	ug/l	93
29) Tetrahydrofuran	7.20	42	26798	12.855	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN100719\
Data File : VN058585.D
Acq On : 7 Oct 2019 22:18
Operator : JC/SP
Sample : K5215-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 30 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
13940

Quant Time: Oct 08 15:44:57 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260
QLast Update : Mon Sep 30 08:07:38 2019
Response via : Initial Calibration

