

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090219\
 Data File : VN057638.D
 Acq On : 1 Sep 2019 12:40
 Operator : JC/SP
 Sample : K4618-15
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-4-28-30

Quant Time: Sep 02 02:12:44 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N090119W.M
 Quant Title : SW846 8260
 QLast Update : Sat Aug 31 15:54:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	188499	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	327412	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	281041	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	89891	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	229827	53.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.94%	
35) Dibromofluoromethane	7.58	113	137223	47.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.98%	
50) Toluene-d8	10.09	98	540290	49.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.44%	
62) 4-Bromofluorobenzene	12.40	95	158648	38.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	76.02%	
Target Compounds						
16) Acetone	3.81	43	17381	14.565	ug/l	100
18) Methyl Acetate	4.32	43	9712	2.370	ug/l #	71
20) Methylene Chloride	4.53	84	4688	1.587	ug/l	89
43) Isopropyl Acetate	8.16	43	77188	10.376	ug/l	98
95) Naphthalene	15.13	128	4302	7.863	ug/l #	93

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN090219\
Data File : VN057638.D
Acq On : 1 Sep 2019 12:40
Operator : JC/SP
Sample : K4618-15
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SB-4-28-30

Quant Time: Sep 02 02:12:44 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N090119W.M
Quant Title : SW846 8260
QLast Update : Sat Aug 31 15:54:32 2019
Response via : Initial Calibration

