

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN062818\
 Data File : VN049595.D
 Acq On : 28 Jun 2018 17:34
 Operator : MD\SY
 Sample : J3677-14
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 93-95-B1(10-15)

Quant Time: Jun 29 03:38:51 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061618W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 16 01:02:01 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1074182	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1721118	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1408397	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	558833	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	716758	50.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.62%	
35) Dibromofluoromethane	7.59	113	640473	46.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.00%	
50) Toluene-d8	10.09	98	2293598	43.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.52%	
62) 4-Bromofluorobenzene	12.40	95	655277	37.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	74.76%	
Target Compounds						
8) Diethyl Ether	3.41	74	7870	1.06	ug/l	92
16) Acetone	3.82	43	78886	22.80	ug/l	90
18) Methyl Acetate	4.33	43	173639	16.46	ug/l	91
20) Methylene Chloride	4.55	84	89852	6.22	ug/l	90
25) 2-Butanone	6.85	43	8746	1.88	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062818\
Data File : VN049595.D
Acq On : 28 Jun 2018 17:34
Operator : MD\SY
Sample : J3677-14
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
93-95-B1(10-15)

Quant Time: Jun 29 03:38:51 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N061618W.M
Quant Title : SW846 8260
QLast Update : Sat Jun 16 01:02:01 2018
Response via : Initial Calibration

