

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090418\
 Data File : VN051026.D
 Acq On : 4 Sep 2018 17:59
 Operator : MD\SY
 Sample : J4761-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RBR-200031

Quant Time: Sep 05 02:34:47 2018
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 22 05:38:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	578700	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	946956	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	791543	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	260707	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	372225	53.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.86%	
35) Dibromofluoromethane	7.59	113	360544	47.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.72%	
50) Toluene-d8	10.09	98	1273944	45.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.46%	
62) 4-Bromofluorobenzene	12.40	95	340709	36.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	72.80%	
Target Compounds						
16) Acetone	3.83	43	44067	20.22	ug/l	94
18) Methyl Acetate	4.34	43	15114	2.30	ug/l	99
20) Methylene Chloride	4.55	84	93816	11.95	ug/l	97
43) Isopropyl Acetate	8.17	43	263024	23.83	ug/l	93
95) Naphthalene	15.14	128	9100	6.18	ug/l	98

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090418\
Data File : VN051026.D
Acq On : 4 Sep 2018 17:59
Operator : MD\SY
Sample : J4761-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
RBR-200031

Quant Time: Sep 05 02:34:47 2018
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
Quant Title : SW846 8260
QLast Update : Wed Aug 22 05:38:12 2018
Response via : Initial Calibration

