

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092718\
 Data File : VN051602.D
 Acq On : 27 Sep 2018 23:25
 Operator : MD\SY
 Sample : J5096-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MT-T-5-2

Quant Time: Sep 28 03:04:47 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092318W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 24 06:41:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	628347	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1039969	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	860222	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	319203	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	457345	58.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	116.20%	
35) Dibromofluoromethane	7.59	113	398322	47.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.80%	
50) Toluene-d8	10.09	98	1430890	45.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.38%	
62) 4-Bromofluorobenzene	12.40	95	392668	38.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	77.52%	
Target Compounds						
18) Methyl Acetate	4.34	43	20690	3.451	ug/l	95
20) Methylene Chloride	4.55	84	50527	6.487	ug/l	98
30) Chloroform	7.37	83	20754	1.548	ug/l	99
43) Isopropyl Acetate	8.17	43	269911	21.557	ug/l #	85
95) Naphthalene	15.14	128	20434	9.489	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092718\
Data File : VN051602.D
Acq On : 27 Sep 2018 23:25
Operator : MD\SY
Sample : J5096-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 67 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MT-T-5-2

Quant Time: Sep 28 03:04:47 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N092318W.M
Quant Title : SW846 8260
QLast Update : Mon Sep 24 06:41:32 2018
Response via : Initial Calibration

