

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN111518\
 Data File : VN052347.D
 Acq On : 15 Nov 2018 17:33
 Operator : MD\SY
 Sample : J5767-10
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 CL-01-111418-A

Quant Time: Nov 16 04:11:56 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N111418W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 15 03:02:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	309533	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	520852	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	447380	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	161381	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	222782	50.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.92%	
35) Dibromofluoromethane	7.59	113	167437	49.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.68%	
50) Toluene-d8	10.09	98	665070	50.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.66%	
62) 4-Bromofluorobenzene	12.40	95	197180	45.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.70%	
Target Compounds						
16) Acetone	3.82	43	8583	7.13	ug/l	94
20) Methylene Chloride	4.55	84	2051964	567.59	ug/l	97
43) Isopropyl Acetate	8.17	43	20150	2.76	ug/l #	97
95) Naphthalene	15.13	128	24199	6.62	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN111518\
Data File : VN052347.D
Acq On : 15 Nov 2018 17:33
Operator : MD\SY
Sample : J5767-10
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 28

Instrument :
MSVOA_N
ClientSampleId :
CL-01-111418-A

Quant Time: Nov 16 04:11:56 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N111418W.M
Quant Title : SW846 8260
QLast Update : Thu Nov 15 03:02:37 2018
Response via : Initial Calibration

