

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN022219\
 Data File : VN053912.D
 Acq On : 22 Feb 2019 13:32
 Operator : JC/SP
 Sample : K1546-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RBR-200034

Quant Time: Feb 22 22:48:55 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N022019W.M
 Quant Title : SW846 8260
 QLast Update : Thu Feb 21 03:36:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	613458	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1038098	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	995490	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	374471	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	331405	52.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.72%	
35) Dibromofluoromethane	7.59	113	331163	48.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.22%	
50) Toluene-d8	10.09	98	1272346	51.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.42%	
62) 4-Bromofluorobenzene	12.40	95	410887	50.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.16%	
Target Compounds						
16) Acetone	3.82	43	19463	11.288	ug/l #	83
20) Methylene Chloride	4.55	84	178474	27.648	ug/l	99
43) Isopropyl Acetate	8.17	43	27478	2.707	ug/l #	88
95) Naphthalene	15.14	128	4709	4.900	ug/l #	93

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN022219\
Data File : VN053912.D
Acq On : 22 Feb 2019 13:32
Operator : JC/SP
Sample : K1546-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
RBR-200034

Quant Time: Feb 22 22:48:55 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N022019W.M
Quant Title : SW846 8260
QLast Update : Thu Feb 21 03:36:07 2019
Response via : Initial Calibration

