

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080218\
 Data File : VN050262.D
 Acq On : 2 Aug 2018 14:29
 Operator : MD\SY
 Sample : PB111674TB
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB111674TB

Quant Time: Aug 03 01:25:03 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 26 18:10:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	687716	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1102375	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	937789	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	339715	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	477899	49.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.62%	
35) Dibromofluoromethane	7.59	113	419916	46.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.30%	
50) Toluene-d8	10.09	98	1522721	46.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.04%	
62) 4-Bromofluorobenzene	12.40	95	417693	38.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	76.68%	
Target Compounds						
16) Acetone	3.83	43	37900	12.646	ug/l	99
43) Isopropyl Acetate	8.17	43	303747	17.882	ug/l	99
59) 2-Hexanone	10.77	43	24812	9.059	ug/l #	60
95) Naphthalene	15.14	128	16458	5.382	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050262.D
Acq On : 2 Aug 2018 14:29
Operator : MD\SY
Sample : PB111674TB
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB111674TB

Quant Time: Aug 03 01:25:03 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260
QLast Update : Thu Jul 26 18:10:10 2018
Response via : Initial Calibration

