

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080218\
 Data File : VN050271.D
 Acq On : 2 Aug 2018 18:14
 Operator : MD\SY
 Sample : PB111674ZHE#09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB111674ZHE#09

Quant Time: Aug 03 01:41:56 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	629848	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1074439	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	914445	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	339938	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	457789	52.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.18%	
35) Dibromofluoromethane	7.60	113	408381	46.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.10%	
50) Toluene-d8	10.09	98	1456918	45.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.34%	
62) 4-Bromofluorobenzene	12.40	95	399835	37.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	75.30%	
Target Compounds						
16) Acetone	3.83	43	35432	13.001	ug/l	99
18) Methyl Acetate	4.34	43	28032	1.915	ug/l	100
43) Isopropyl Acetate	8.18	43	1276285	80.954	ug/l #	71
59) 2-Hexanone	10.77	43	24113	9.048	ug/l #	57

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050271.D
Acq On : 2 Aug 2018 18:14
Operator : MD\SY
Sample : PB111674ZHE#09
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB111674ZHE#09

Quant Time: Aug 03 01:41:56 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

