

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091720\
 Data File : VN063417.D
 Acq On : 17 Sep 2020 14:26
 Operator : JC/MD
 Sample : PB131674TB
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB131674TB

Quant Time: Sep 18 03:48:05 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	255148	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	473943	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	473418	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	208519	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	207186	47.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.40%	
35) Dibromofluoromethane	7.55	113	151276	46.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.22%	
50) Toluene-d8	10.06	98	611587	49.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.14%	
62) 4-Bromofluorobenzene	12.38	95	245707	54.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.66%	
Target Compounds						
16) Acetone	3.78	43	14416	10.328	ug/l	99
18) Methyl Acetate	4.29	43	5200	1.467	ug/l #	70
20) Methylene Chloride	4.51	84	14125	3.994	ug/l	87
43) Isopropyl Acetate	8.13	43	92600	10.450	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091720\
Data File : VN063417.D
Acq On : 17 Sep 2020 14:26
Operator : JC/MD
Sample : PB131674TB
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB131674TB

Quant Time: Sep 18 03:48:05 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 16 13:05:20 2020
Response via : Initial Calibration

