

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101620\
 Data File : VN064169.D
 Acq On : 16 Oct 2020 17:08
 Operator : JC/MD
 Sample : PB132400TB
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB132400TB

Quant Time: Oct 17 02:03:05 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N101420W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 15 01:16:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	310322	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	595733	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	645118	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	229493	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	262093	59.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	118.60%	
35) Dibromofluoromethane	7.55	113	210049	50.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.88%	
50) Toluene-d8	10.06	98	816800	52.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.00%	
62) 4-Bromofluorobenzene	12.38	95	272416	50.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.88%	
Target Compounds						
16) Acetone	3.79	43	14598	11.136	ug/l	96
18) Methyl Acetate	4.29	43	4668	1.420	ug/l #	81
20) Methylene Chloride	4.51	84	16020	3.949	ug/l	93
43) Isopropyl Acetate	8.13	43	118380	13.999	ug/l #	88

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN101620\
Data File : VN064169.D
Acq On : 16 Oct 2020 17:08
Operator : JC/MD
Sample : PB132400TB
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB132400TB

Quant Time: Oct 17 02:03:05 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N101420W.M
Quant Title : SW846 8260
QLast Update : Thu Oct 15 01:16:44 2020
Response via : Initial Calibration

