

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031320\
 Data File : VN060502.D
 Acq On : 13 Mar 2020 15:04
 Operator : JC/MD
 Sample : PB127506ZHE#03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB127506ZHE#03

Quant Time: Mar 16 08:55:32 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N022720W.M
 Quant Title : SW846 8260
 QLast Update : Thu Feb 27 13:52:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	206014	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	344187	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	325224	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	153024	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	133968	50.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.36%	
35) Dibromofluoromethane	7.57	113	102265	64.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	128.02%	
50) Toluene-d8	10.08	98	424155	52.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.92%	
62) 4-Bromofluorobenzene	12.40	95	162583	53.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.08%	
Target Compounds						
16) Acetone	3.79	43	6260	7.334	ug/l	95
20) Methylene Chloride	4.52	84	4033	1.720	ug/l #	95
43) Isopropyl Acetate	8.14	43	67448	12.404	ug/l #	90
51) 4-Methyl-2-Pentanone	9.97	43	5183	1.681	ug/l	94

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN031320\
Data File : VN060502.D
Acq On : 13 Mar 2020 15:04
Operator : JC/MD
Sample : PB127506ZHE#03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB127506ZHE#03

Quant Time: Mar 16 08:55:32 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N022720W.M
Quant Title : SW846 8260
QLast Update : Thu Feb 27 13:52:50 2020
Response via : Initial Calibration

