

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN021720\
 Data File : VN060120.D
 Acq On : 17 Feb 2020 15:30
 Operator : JC/MD
 Sample : PB126834ZHE#05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB126834ZHE#05

Manual Integrations
 APPROVED

MMDadoda
 2/18/2020 9:28:14 AM

Quant Time: Feb 17 17:25:58 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 12 14:12:21 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	182774	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	290852	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	275596	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	124052	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	113434	51.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.00%	
35) Dibromofluoromethane	7.57	113	88852	49.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.40%	
50) Toluene-d8	10.08	98	358058	50.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.72%	
62) 4-Bromofluorobenzene	12.40	95	128857	49.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.62%	
Target Compounds						
16) Acetone	3.79	43	6003m	6.188	ug/l	
20) Methylene Chloride	4.52	84	7809	4.035	ug/l	96
43) Isopropyl Acetate	8.14	43	55662	13.374	ug/l #	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN021720\
Data File : VN060120.D
Acq On : 17 Feb 2020 15:30
Operator : JC/MD
Sample : PB126834ZHE#05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB126834ZHE#05

Manual Integrations
APPROVED

MMDadoda
2/18/2020 9:28:14 AM

Quant Time: Feb 17 17:25:58 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
Quant Title : SW846 8260
QLast Update : Wed Feb 12 14:12:21 2020
Response via : Initial Calibration

