

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN073120\
 Data File : VN062680.D
 Acq On : 31 Jul 2020 13:05
 Operator : JC/MD
 Sample : PB130640ZHE#03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB130640ZHE#03

Quant Time: Aug 01 06:41:20 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 21 18:48:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	134683	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	247116	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	253461	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	92334	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	119194	56.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.94%	
35) Dibromofluoromethane	7.55	113	83031	51.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.30%	
50) Toluene-d8	10.06	98	330332	51.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.02%	
62) 4-Bromofluorobenzene	12.38	95	122125	51.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.14%	
Target Compounds						
8) Diethyl Ether	3.37	74	1248	1.203	ug/l	96
16) Acetone	3.78	43	15266	19.390	ug/l	100
18) Methyl Acetate	4.29	43	4328	2.171	ug/l #	79
20) Methylene Chloride	4.50	84	15551	8.235	ug/l	94
43) Isopropyl Acetate	8.13	43	48273	10.546	ug/l #	90

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN073120\
Data File : VN062680.D
Acq On : 31 Jul 2020 13:05
Operator : JC/MD
Sample : PB130640ZHE#03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB130640ZHE#03

Quant Time: Aug 01 06:41:20 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
Quant Title : SW846 8260
QLast Update : Tue Jul 21 18:48:59 2020
Response via : Initial Calibration

