

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061820\
 Data File : VN061932.D
 Acq On : 19 Jun 2020 5:18
 Operator : JC/MD
 Sample : PB129596ZHE#09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB129596ZHE#09

Quant Time: Jun 19 07:53:17 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 17 01:19:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	142576	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	275505	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	270480	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	101949	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	98236	51.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.74%	
35) Dibromofluoromethane	7.55	113	85460	48.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.98%	
50) Toluene-d8	10.06	98	345263	49.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.66%	
62) 4-Bromofluorobenzene	12.38	95	112386	47.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.56%	
Target Compounds						
16) Acetone	3.78	43	10580	18.732	ug/l	92
18) Methyl Acetate	4.28	43	3550	2.183	ug/l #	77
20) Methylene Chloride	4.50	84	4497	2.444	ug/l	92
43) Isopropyl Acetate	8.13	43	34223	10.349	ug/l #	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN061820\
Data File : VN061932.D
Acq On : 19 Jun 2020 5:18
Operator : JC/MD
Sample : PB129596ZHE#09
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB129596ZHE#09

Quant Time: Jun 19 07:53:17 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N061620W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 17 01:19:44 2020
Response via : Initial Calibration

