

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100219\
 Data File : VN058479.D
 Acq On : 2 Oct 2019 17:01
 Operator : JC/SP
 Sample : PB123550TB
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB123550TB

Quant Time: Oct 03 05:34:07 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 30 08:07:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	327530	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	547664	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	468095	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	167187	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	245589	51.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.62%	
35) Dibromofluoromethane	7.58	113	170296	50.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.48%	
50) Toluene-d8	10.09	98	693133	52.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.42%	
62) 4-Bromofluorobenzene	12.41	95	210477	44.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.54%	
Target Compounds						
16) Acetone	3.81	43	10973	6.725	ug/l	98
18) Methyl Acetate	4.31	43	4331	1.055	ug/l #	70
20) Methylene Chloride	4.53	84	89480	23.764	ug/l	99
43) Isopropyl Acetate	8.15	43	99723	11.944	ug/l	100
95) Naphthalene	15.14	128	10153	1.214	ug/l #	95

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN100219\
Data File : VN058479.D
Acq On : 2 Oct 2019 17:01
Operator : JC/SP
Sample : PB123550TB
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB123550TB

Quant Time: Oct 03 05:34:07 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260
QLast Update : Mon Sep 30 08:07:38 2019
Response via : Initial Calibration

