

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081919\
 Data File : VN057401.D
 Acq On : 19 Aug 2019 15:32
 Operator : JC/SP
 Sample : K4339-23
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T8-B1-(0)

Quant Time: Aug 20 08:19:57 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081519W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 15 11:33:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	1041004	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	1441393	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	1221801	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	327060	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	476874	42.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	84.28%	
35) Dibromofluoromethane	7.58	113	441961	46.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.56%	
50) Toluene-d8	10.09	98	1657502	47.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.92%	
62) 4-Bromofluorobenzene	12.40	95	449221	36.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	73.56%	
Target Compounds						
16) Acetone	3.81	43	25363	6.461	ug/l	89
18) Methyl Acetate	4.32	43	36041	1.068	ug/l	99
20) Methylene Chloride	4.55	84	20955	1.908	ug/l	90
43) Isopropyl Acetate	8.16	43	214107	10.337	ug/l #	80

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN081919\
Data File : VN057401.D
Acq On : 19 Aug 2019 15:32
Operator : JC/SP
Sample : K4339-23
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T8-B1-(0)

Quant Time: Aug 20 08:19:57 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N081519W.M
Quant Title : SW846 8260
QLast Update : Thu Aug 15 11:33:14 2019
Response via : Initial Calibration

