

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091319\
 Data File : VN057969.D
 Acq On : 12 Sep 2019 14:31
 Operator : JC/SP
 Sample : K4644-23
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T11-D4-(4)

Quant Time: Sep 13 05:39:39 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	262855	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	452230	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	406440	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	204148	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	322799	60.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.26%	
35) Dibromofluoromethane	7.58	113	221299	53.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.04%	
50) Toluene-d8	10.09	98	811679	50.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.18%	
62) 4-Bromofluorobenzene	12.41	95	271935	44.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.06%	
Target Compounds						
16) Acetone	3.81	43	10102	5.788	ug/l	98
18) Methyl Acetate	4.32	43	7819	1.680	ug/l #	86
20) Methylene Chloride	4.53	84	9376	1.631	ug/l	93
43) Isopropyl Acetate	8.16	43	22709	2.492	ug/l #	90

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091319\
Data File : VN057969.D
Acq On : 12 Sep 2019 14:31
Operator : JC/SP
Sample : K4644-23
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T11-D4-(4)

Quant Time: Sep 13 05:39:39 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 10 03:02:13 2019
Response via : Initial Calibration

