

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091619\
 Data File : VN058125.D
 Acq On : 16 Sep 2019 17:35
 Operator : JC/SP
 Sample : K4878-19
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CSA-2-7

Quant Time: Sep 17 09:14:00 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	336489	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	566594	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	529509	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	274940	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	355982	52.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.46%	
35) Dibromofluoromethane	7.58	113	242375	46.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.70%	
50) Toluene-d8	10.08	98	970328	47.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.58%	
62) 4-Bromofluorobenzene	12.41	95	325755	42.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	84.20%	
Target Compounds						
16) Acetone	3.80	43	20593	9.217	ug/l #	89
18) Methyl Acetate	4.31	43	12403	2.082	ug/l	91
20) Methylene Chloride	4.53	84	12274	1.697	ug/l #	91
43) Isopropyl Acetate	8.16	43	107595	9.424	ug/l #	86

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091619\
Data File : VN058125.D
Acq On : 16 Sep 2019 17:35
Operator : JC/SP
Sample : K4878-19
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CSA-2-7

Quant Time: Sep 17 09:14:00 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

