

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091419\
 Data File : VN058059.D
 Acq On : 14 Sep 2019 8:11
 Operator : JC/SP
 Sample : K4825-17
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CSA-4-3

Quant Time: Sep 16 02:16:00 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	324762	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	567632	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	516795	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	272051	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	374203	56.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.78%	
35) Dibromofluoromethane	7.58	113	250358	47.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.58%	
50) Toluene-d8	10.09	98	1032360	50.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.50%	
62) 4-Bromofluorobenzene	12.40	95	335720	43.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.62%	
Target Compounds						
16) Acetone	3.80	43	13464	6.244	ug/l #	84
18) Methyl Acetate	4.31	43	10894	1.895	ug/l #	89
20) Methylene Chloride	4.53	84	84395	19.816	ug/l	97
43) Isopropyl Acetate	8.15	43	130754	11.432	ug/l #	83

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091419\
Data File : VN058059.D
Acq On : 14 Sep 2019 8:11
Operator : JC/SP
Sample : K4825-17
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 62 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CSA-4-3

Quant Time: Sep 16 02:16:00 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

