

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101719\
 Data File : VN058840.D
 Acq On : 17 Oct 2019 19:56
 Operator : JC/SP
 Sample : K5394-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T-15-E1-(20)

Quant Time: Oct 18 04:49:51 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N101319W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 12 00:00:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	352699	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	570300	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	525877	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	199398	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	265255	53.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.44%	
35) Dibromofluoromethane	7.57	113	188678	53.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.12%	
50) Toluene-d8	10.09	98	731885	51.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.92%	
62) 4-Bromofluorobenzene	12.40	95	235905	44.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.48%	
Target Compounds						
16) Acetone	3.80	43	10811	5.289	ug/l	93
18) Methyl Acetate	4.31	43	9676	1.866	ug/l	99
20) Methylene Chloride	4.53	84	18749	4.389	ug/l	92
43) Isopropyl Acetate	8.15	43	78067	7.858	ug/l #	91

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN101719\
Data File : VN058840.D
Acq On : 17 Oct 2019 19:56
Operator : JC/SP
Sample : K5394-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T-15-E1-(20)

Quant Time: Oct 18 04:49:51 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N101319W.M
Quant Title : SW846 8260
QLast Update : Sat Oct 12 00:00:30 2019
Response via : Initial Calibration

