

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN071619\
 Data File : VN056779.D
 Acq On : 16 Jul 2019 19:16
 Operator : JC/SP
 Sample : K3627-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TR-04-071519

Quant Time: Jul 17 03:31:58 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N071519W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 16 07:03:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	263416	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	470194	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	500510	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	191740	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	152795	52.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.88%	
35) Dibromofluoromethane	7.59	113	139090	47.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.06%	
50) Toluene-d8	10.09	98	596299	52.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.52%	
62) 4-Bromofluorobenzene	12.40	95	218914	57.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.88%	
Target Compounds						
16) Acetone	3.82	43	19438	14.477	ug/l	98
18) Methyl Acetate	4.34	43	9020	2.058	ug/l #	77
20) Methylene Chloride	4.55	84	288330	98.263	ug/l	98
43) Isopropyl Acetate	8.17	43	13753	2.040	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN071619\
Data File : VN056779.D
Acq On : 16 Jul 2019 19:16
Operator : JC/SP
Sample : K3627-04
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TR-04-071519

Quant Time: Jul 17 03:31:58 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N071519W.M
Quant Title : SW846 8260
QLast Update : Tue Jul 16 07:03:34 2019
Response via : Initial Calibration

