

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN071919\
 Data File : VN056836.D
 Acq On : 19 Jul 2019 16:50
 Operator : JC/SP
 Sample : K3615-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HR-06-071719-C

Quant Time: Jul 20 00:15:02 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	249056	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	450354	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	489998	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	177231	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	145609	52.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.70%	
35) Dibromofluoromethane	7.59	113	132800	47.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.76%	
50) Toluene-d8	10.09	98	571681	52.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.62%	
62) 4-Bromofluorobenzene	12.40	95	210483	57.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.32%	
Target Compounds						
16) Acetone	3.82	43	17427	13.727	ug/l	97
18) Methyl Acetate	4.33	43	6019	1.160	ug/l #	90
20) Methylene Chloride	4.55	84	499121	179.907	ug/l	98
43) Isopropyl Acetate	8.17	43	15602	2.416	ug/l #	91

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN071919\
Data File : VN056836.D
Acq On : 19 Jul 2019 16:50
Operator : JC/SP
Sample : K3615-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HR-06-071719-C

Quant Time: Jul 20 00:15:02 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

