

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061419\
 Data File : VN056322.D
 Acq On : 14 Jun 2019 19:13
 Operator : JC/SP
 Sample : K3352-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 20190612-LOC-12

Quant Time: Jun 15 00:44:08 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 04 11:02:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	328641	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	566522	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	580092	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	186884	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	195278	59.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	119.68%	
35) Dibromofluoromethane	7.59	113	175877	50.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.20%	
50) Toluene-d8	10.09	98	716222	53.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.10%	
62) 4-Bromofluorobenzene	12.40	95	240494	52.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.74%	
Target Compounds						
16) Acetone	3.82	43	27523	17.636	ug/l	96
18) Methyl Acetate	4.33	43	5819	1.336	ug/l #	78
20) Methylene Chloride	4.54	84	80246	24.957	ug/l	98
30) Chloroform	7.37	83	7597	1.454	ug/l	96
43) Isopropyl Acetate	8.17	43	11274	1.605	ug/l #	90

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN061419\
Data File : VN056322.D
Acq On : 14 Jun 2019 19:13
Operator : JC/SP
Sample : K3352-04
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
20190612-LOC-12

Quant Time: Jun 15 00:44:08 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
Quant Title : SW846 8260
QLast Update : Tue Jun 04 11:02:09 2019
Response via : Initial Calibration

