

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080519\
 Data File : VN057086.D
 Acq On : 5 Aug 2019 16:48
 Operator : JC/SP
 Sample : K4130-38
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1S-M-(0-1)

Quant Time: Aug 06 02:23:44 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072719W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jul 27 01:33:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	573894	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	833849	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	689494	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	178980	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	284952	48.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.02%	
35) Dibromofluoromethane	7.58	113	255850	47.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.86%	
50) Toluene-d8	10.08	98	974876	47.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.78%	
62) 4-Bromofluorobenzene	12.40	95	255934	38.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	77.24%	
Target Compounds						
8) Diethyl Ether	3.40	74	4838	1.516	ug/l	98
16) Acetone	3.81	43	24131	11.338	ug/l	98
18) Methyl Acetate	4.32	43	8539	1.258	ug/l #	76
20) Methylene Chloride	4.54	84	17270	2.817	ug/l	94
43) Isopropyl Acetate	8.17	43	25313	2.606	ug/l #	82

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080519\
Data File : VN057086.D
Acq On : 5 Aug 2019 16:48
Operator : JC/SP
Sample : K4130-38
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
1S-M-(0-1)

Quant Time: Aug 06 02:23:44 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072719W.M
Quant Title : SW846 8260
QLast Update : Sat Jul 27 01:33:29 2019
Response via : Initial Calibration

