

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080619\
 Data File : VN057102.D
 Acq On : 6 Aug 2019 14:44
 Operator : JC/SP
 Sample : K4152-22
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 2S-H1-(0-1)

Quant Time: Aug 07 04:18:43 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	7.65	168	494728	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	719101	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	616517	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	208460	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	249449	49.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.52%	
35) Dibromofluoromethane	7.58	113	217814	47.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.64%	
50) Toluene-d8	10.08	98	849560	48.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.78%	
62) 4-Bromofluorobenzene	12.40	95	253892	44.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.86%	
Target Compounds						
8) Diethyl Ether	3.40	74	3730	1.356	ug/l #	36
16) Acetone	3.82	43	14272	7.779	ug/l	93
18) Methyl Acetate	4.32	43	19322	3.301	ug/l	94
20) Methylene Chloride	4.52	84	15523	2.937	ug/l	89
43) Isopropyl Acetate	8.16	43	51384	6.134	ug/l #	67

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080619\
Data File : VN057102.D
Acq On : 6 Aug 2019 14:44
Operator : JC/SP
Sample : K4152-22
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
2S-H1-(0-1)

Quant Time: Aug 07 04:18:43 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

