

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080619\
 Data File : VN057099.D
 Acq On : 6 Aug 2019 13:34
 Operator : JC/SP
 Sample : K4152-37
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 2S-E1-(30)

Quant Time: Aug 07 04:12:23 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	582880	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	850526	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	701808	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	234554	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	292153	48.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.92%	
35) Dibromofluoromethane	7.58	113	255514	46.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.86%	
50) Toluene-d8	10.08	98	977433	47.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.14%	
62) 4-Bromofluorobenzene	12.40	95	277800	41.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	82.20%	
Target Compounds						
8) Diethyl Ether	3.39	74	3610	1.114	ug/l	84
16) Acetone	3.82	43	12468	5.768	ug/l	93
18) Methyl Acetate	4.32	43	18852	2.734	ug/l #	75
20) Methylene Chloride	4.52	84	16577	2.662	ug/l	91
43) Isopropyl Acetate	8.16	43	44682	4.510	ug/l #	69

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080619\
Data File : VN057099.D
Acq On : 6 Aug 2019 13:34
Operator : JC/SP
Sample : K4152-37
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
2S-E1-(30)

Quant Time: Aug 07 04:12:23 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

