

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092319\
 Data File : VN058324.D
 Acq On : 24 Sep 2019 3:43
 Operator : JC/SP
 Sample : K4997-20
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-12-091919

Manual Integrations
 APPROVED

MMDadoda
 9/25/2019 12:49:49 AM

Quant Time: Sep 24 09:17:02 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091819W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 19 09:27:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	464533	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	791796	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	687860	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	281938	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	360743	55.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.52%	
35) Dibromofluoromethane	7.57	113	244028	50.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.40%	
50) Toluene-d8	10.08	98	998009	53.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.44%	
62) 4-Bromofluorobenzene	12.41	95	328588	47.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.34%	
Target Compounds						
16) Acetone	3.81	43	12850m	5.851	ug/l	
20) Methylene Chloride	4.53	84	65004	16.758	ug/l	95
43) Isopropyl Acetate	8.16	43	131044	11.841	ug/l #	82

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092319\
Data File : VN058324.D
Acq On : 24 Sep 2019 3:43
Operator : JC/SP
Sample : K4997-20
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 53 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-12-091919

Manual Integrations
APPROVED

MMDadoda
9/25/2019 12:49:49 AM

Quant Time: Sep 24 09:17:02 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091819W.M
Quant Title : SW846 8260
QLast Update : Thu Sep 19 09:27:53 2019
Response via : Initial Calibration

