

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091019\
 Data File : VN057838.D
 Acq On : 10 Sep 2019 6:08
 Operator : JC/SP
 Sample : K4622-23
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAR-11-A2-(0)

Manual Integrations
 APPROVED

MMDadoda
 9/13/2019 5:16:27 AM

Quant Time: Sep 10 08:41:04 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	313619	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	529423	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	489526	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	249818	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	339681	53.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.96%	
35) Dibromofluoromethane	7.58	113	234400	47.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.94%	
50) Toluene-d8	10.09	98	962932	50.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.52%	
62) 4-Bromofluorobenzene	12.40	95	314725	43.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.06%	
Target Compounds						
16) Acetone	3.80	43	9596	4.608	ug/l	93
18) Methyl Acetate	4.31	43	8055m	1.451	ug/l	
43) Isopropyl Acetate	8.15	43	101169	9.483	ug/l #	85

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091019\
Data File : VN057838.D
Acq On : 10 Sep 2019 6:08
Operator : JC/SP
Sample : K4622-23
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleID :
PAR-11-A2-(0)

Manual Integrations
APPROVED

MMDadoda
9/13/2019 5:16:27 AM

Quant Time: Sep 10 08:41:04 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 10 03:02:13 2019
Response via : Initial Calibration

