

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091820\
 Data File : VN063467.D
 Acq On : 18 Sep 2020 19:01
 Operator : JC/MD
 Sample : L3871-16
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TR-05-091720

Manual Integrations
 APPROVED

MMDadoda
 9/21/2020 11:49:53 AM

Quant Time: Sep 21 02:29:10 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	233294	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	438565	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	442438	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	190915	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	191979	48.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.68%	
35) Dibromofluoromethane	7.55	113	138977	45.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.54%	
50) Toluene-d8	10.06	98	567118	49.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.36%	
62) 4-Bromofluorobenzene	12.38	95	226009	54.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.00%	
Target Compounds						
16) Acetone	3.78	43	42718	33.470	ug/l	99
18) Methyl Acetate	4.30	43	4440m	1.370	ug/l	
20) Methylene Chloride	4.51	84	13998	4.329	ug/l	95
43) Isopropyl Acetate	8.13	43	91033	11.102	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091820\
Data File : VN063467.D
Acq On : 18 Sep 2020 19:01
Operator : JC/MD
Sample : L3871-16
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TR-05-091720

Manual Integrations
APPROVED

MMDadoda
9/21/2020 11:49:53 AM

Quant Time: Sep 21 02:29:10 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 16 13:05:20 2020
Response via : Initial Calibration

