

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111220\
 Data File : VN064576.D
 Acq On : 12 Nov 2020 13:10
 Operator : JC/MD
 Sample : L4697-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-2-S

Manual Integrations
 APPROVED

MMDadoda
 11/13/2020 11:42:20 AM

Quant Time: Nov 13 03:26:58 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 04 12:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	225992	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	462549	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	517829	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	183177	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	224102	56.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.86%	
35) Dibromofluoromethane	7.55	113	166032	52.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.44%	
50) Toluene-d8	10.06	98	628218	51.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.22%	
62) 4-Bromofluorobenzene	12.38	95	235306	51.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.24%	
Target Compounds						
16) Acetone	3.79	43	11585	7.726	ug/l #	83
18) Methyl Acetate	4.29	43	4354m	1.477	ug/l	
20) Methylene Chloride	4.51	84	13542	4.552	ug/l	91
43) Isopropyl Acetate	8.13	43	42148	5.322	ug/l #	93

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111220\
Data File : VN064576.D
Acq On : 12 Nov 2020 13:10
Operator : JC/MD
Sample : L4697-04
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-2-S

Manual Integrations
APPROVED

MMDadoda
11/13/2020 11:42:20 AM

Quant Time: Nov 13 03:26:58 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260
QLast Update : Wed Nov 04 12:09:03 2020
Response via : Initial Calibration

