

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111220\
 Data File : VN064577.D
 Acq On : 12 Nov 2020 13:34
 Operator : JC/MD
 Sample : L4697-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-2-D

Manual Integrations
 APPROVED

MMDadoda
 11/13/2020 11:42:24 AM

Quant Time: Nov 13 03:29:04 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	231131	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	487029	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	548038	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	193727	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	231870	57.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.18%	
35) Dibromofluoromethane	7.55	113	170095	51.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.60%	
50) Toluene-d8	10.06	98	659210	51.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.88%	
62) 4-Bromofluorobenzene	12.38	95	243946	50.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.66%	
Target Compounds						
16) Acetone	3.79	43	11780	7.681	ug/l #	89
18) Methyl Acetate	4.28	43	3138m	1.041	ug/l	
20) Methylene Chloride	4.51	84	12988	4.269	ug/l	96
43) Isopropyl Acetate	8.13	43	39925	4.788	ug/l #	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111220\
Data File : VN064577.D
Acq On : 12 Nov 2020 13:34
Operator : JC/MD
Sample : L4697-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-2-D

Manual Integrations
APPROVED

MMDadoda
11/13/2020 11:42:24 AM

Quant Time: Nov 13 03:29:04 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260
QLast Update : Wed Nov 04 12:09:03 2020
Response via : Initial Calibration

