

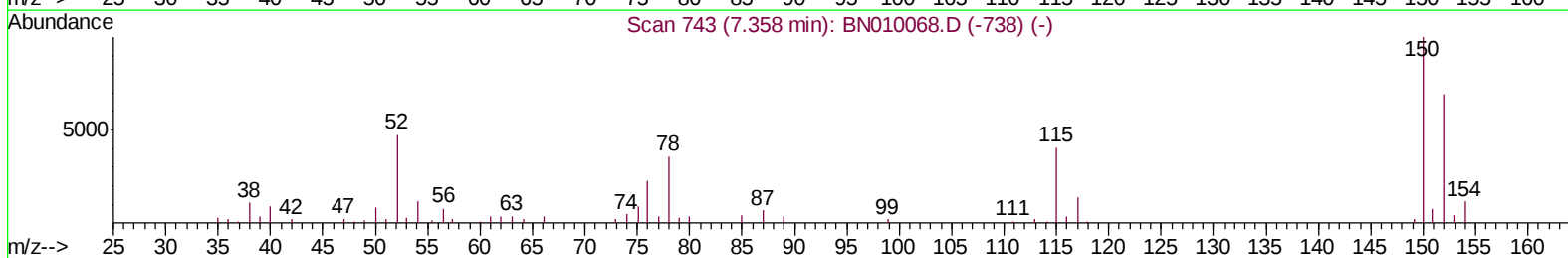
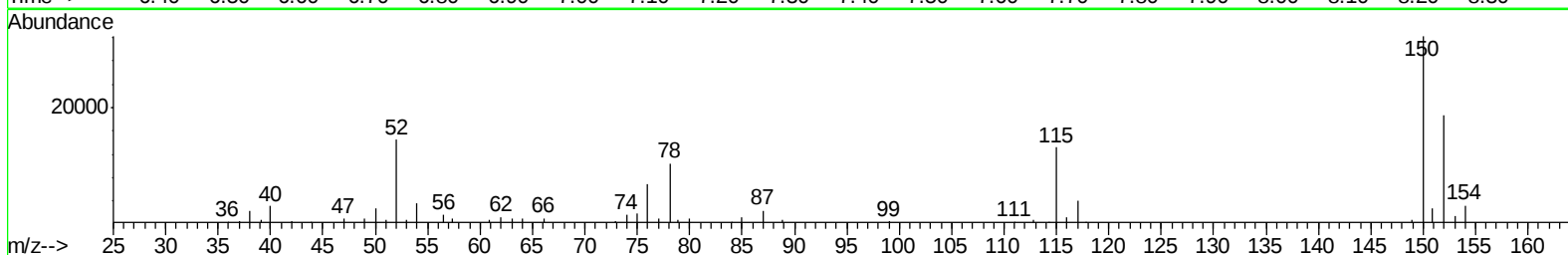
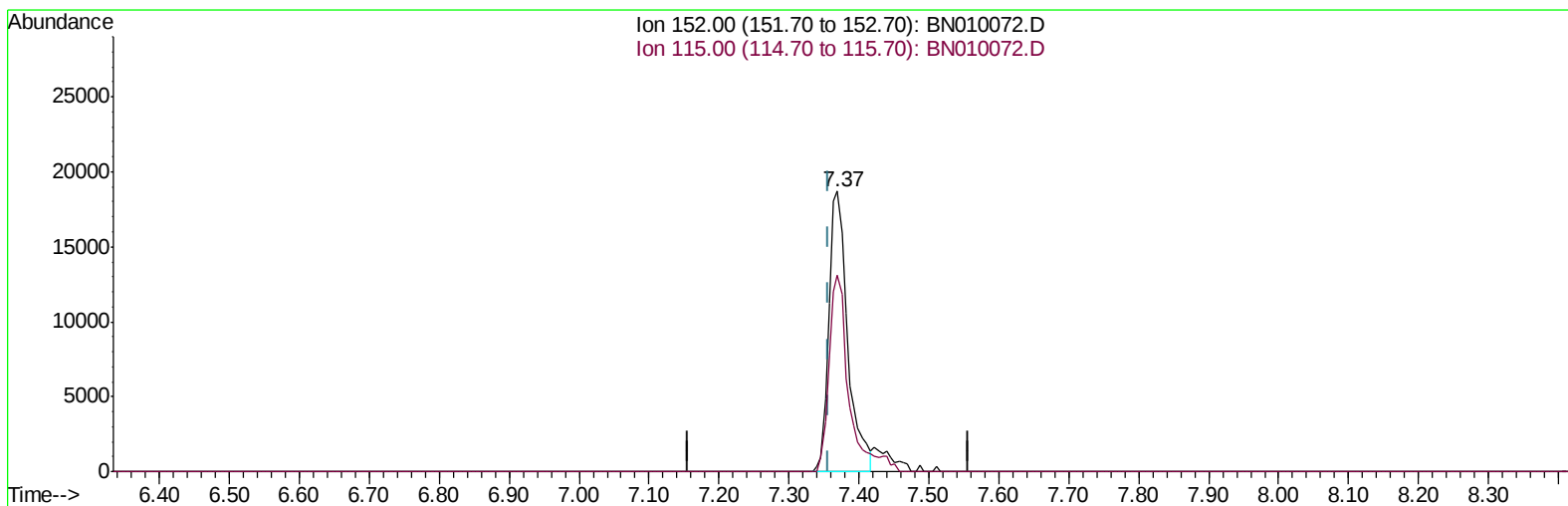
Data File : BN010072.D
Acq On : 03 Mar 2020 21:33
Operator : CG/JU
Sample : L1733-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SVOC-GPC-BLANK

Manual Integrations
APPROVED

mohammad
3/4/2020 7:51:57 AM

Quant Time: Mar 04 01:32:43 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 04 01:30:16 2020
Response via : Initial Calibration



TIC: BN010072.D

(1) 1,4-Dichlorobenzene-d4 (I)

7.369min (+0.012) 20.00ng/ul

response 34963

Ion	Exp%	Act%
152.00	100	100
115.00	65.80	70.16
0.00	0.00	0.00
0.00	0.00	0.00

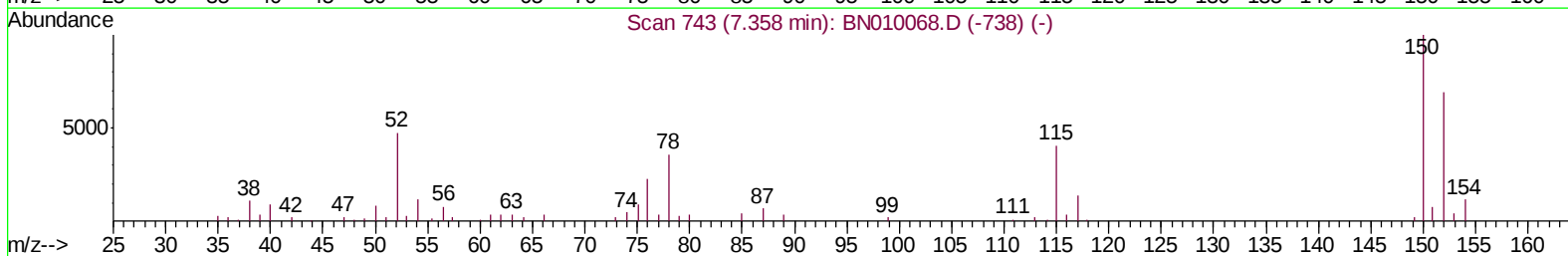
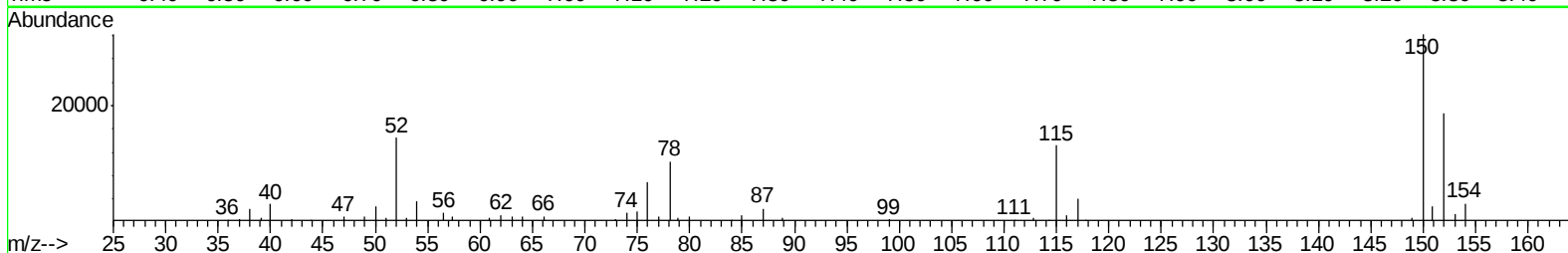
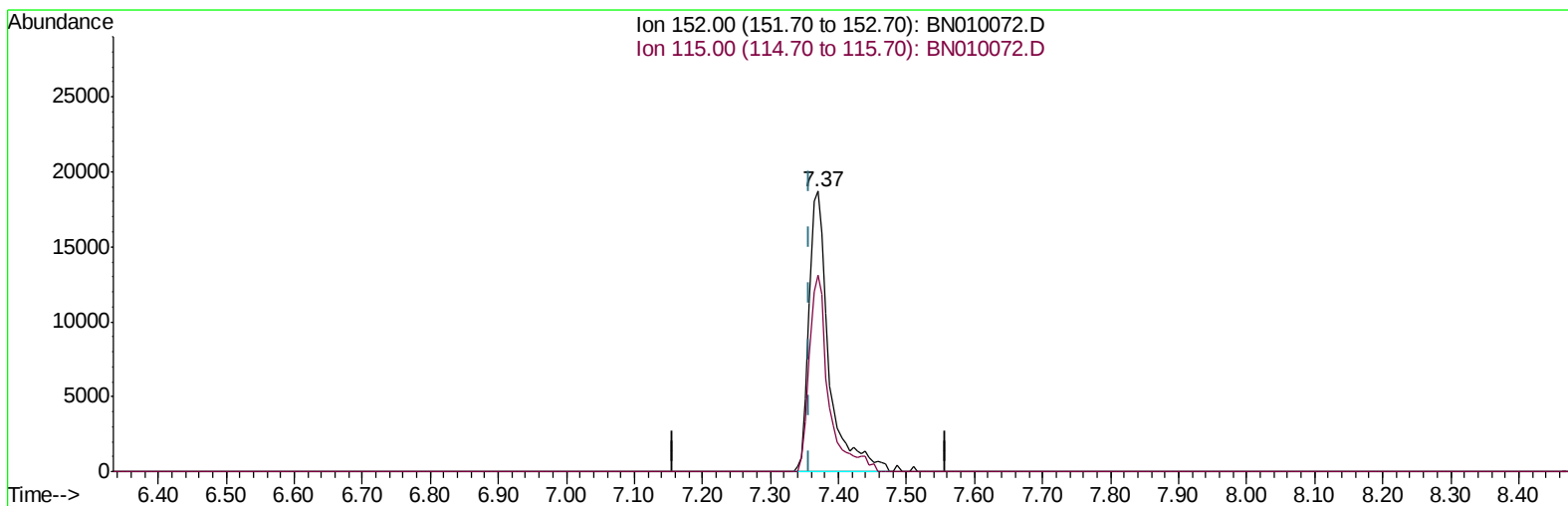
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TIC: BN010072.D

(1) 1,4-Dichlorobenzene-d4 (I)

7.369min (+0.012) 20.00ng/ul m

response 38122

Ion	Exp%	Act%
152.00	100	100
115.00	65.80	70.16
0.00	0.00	0.00
0.00	0.00	0.00

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mohammad
3/4/2020 7:51:57 AM

Quant Time: Mar 04 01:58:20 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	38122m	20.00	ng/ul	0.01
18) Naphthalene-d8	10.12	136	146139	20.00	ng/ul	0.01
36) Acenaphthene-d10	14.01	164	105400	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.77	188	230655	20.00	ng/ul	0.00
78) Chrysene-d12	20.99	240	207359	20.00	ng/ul	0.00
86) Perylene-d12	23.08	264	225000	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	0.00	176	0	0.00	ng/ul	
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
79) Pyrene-d10	0.00	212	0	0.00	ng/ul	
90) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul	

Target Compounds Qvalue

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

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