

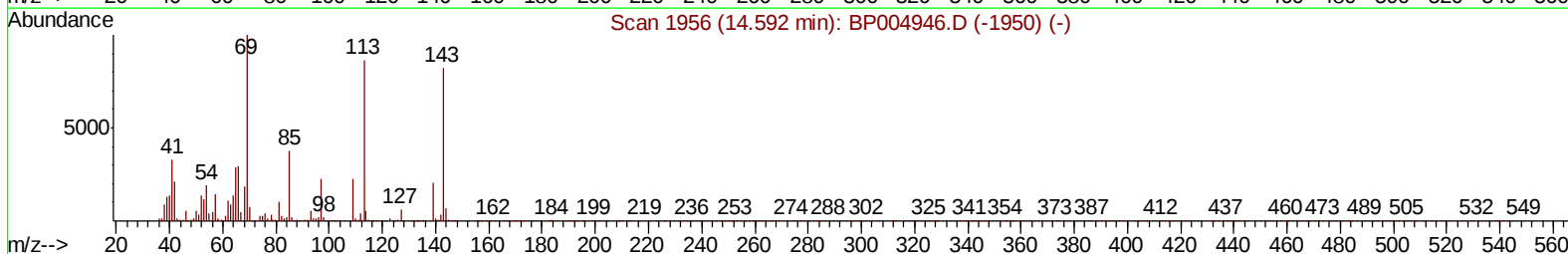
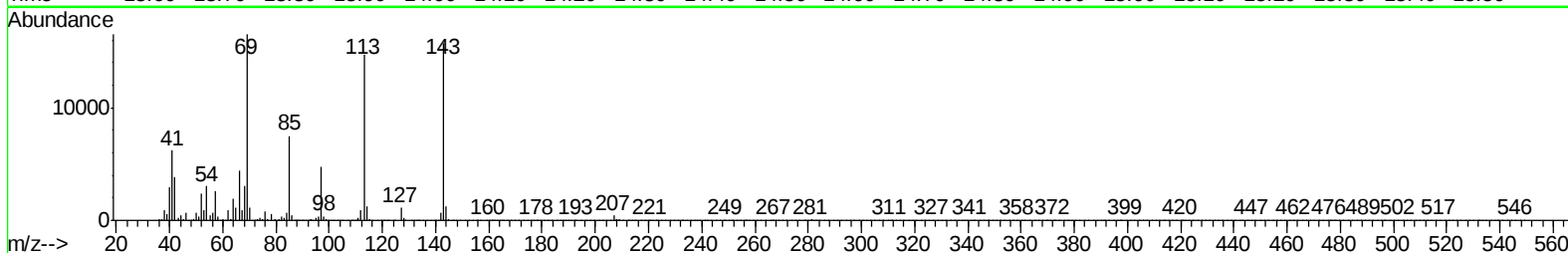
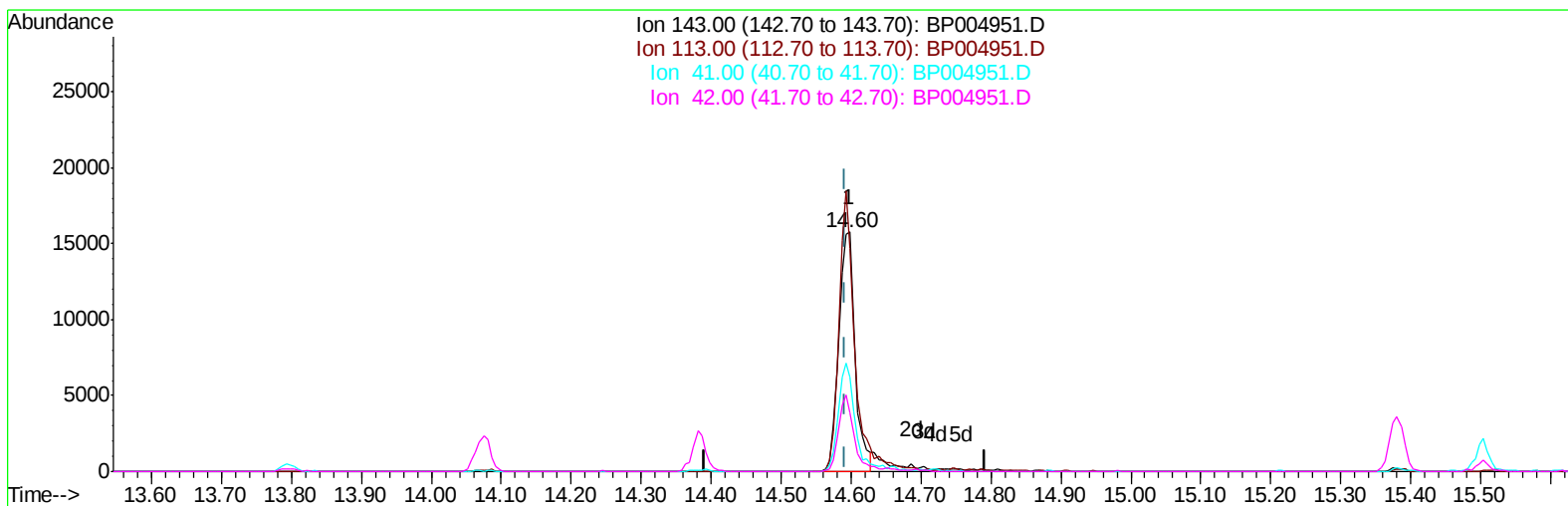
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP030921\
Data File : BP004951.D
Acq On : 09 Mar 2021 19:51
Operator : CG/JU
Sample : PB134968BL
Misc : SOM-MDL-S-PB-07
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK68

Manual Integrations
APPROVED

mohammad
3/10/2021 11:54:35 AM

Quant Time: Mar 10 00:55:53 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP030921MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 09 15:13:30 2021
Response via : Initial Calibration



TIC: BP004951.D

(52) 4-Nitrophenol-d4 (S)

14.598min (+0.006) 27.46ng/ul

response 26043

Ion	Exp%	Act%
143.00	100	100
113.00	98.40	93.34
41.00	40.30	39.20
42.00	28.70	24.55

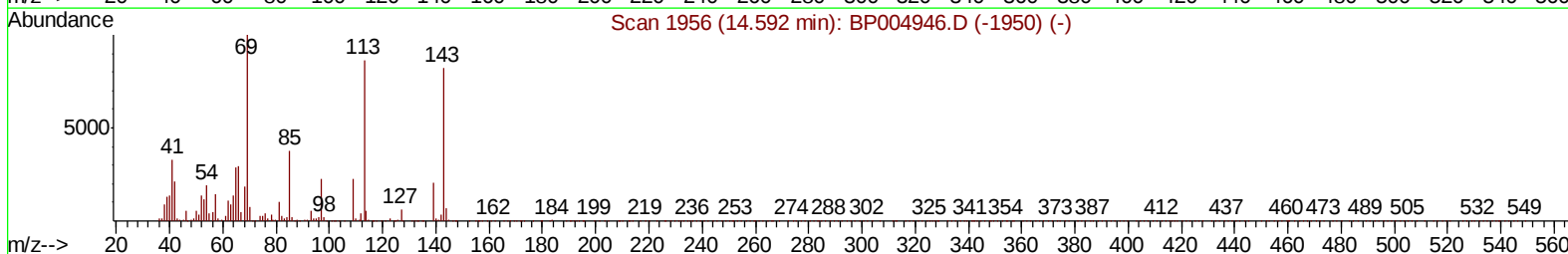
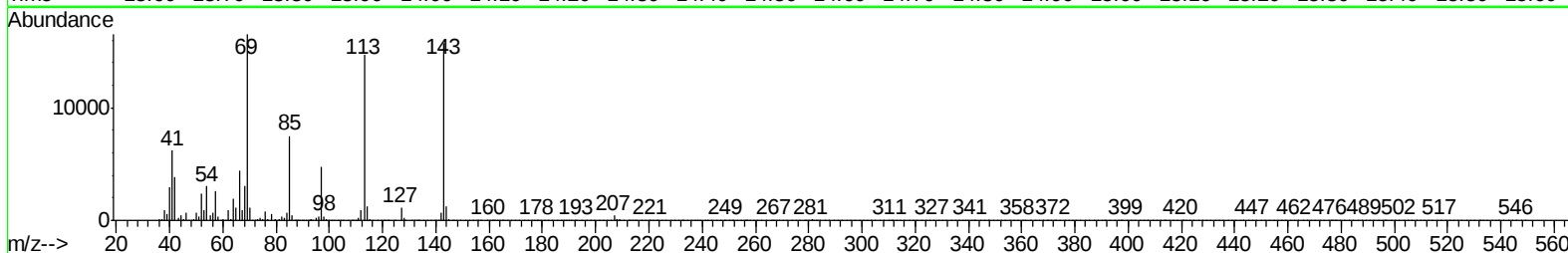
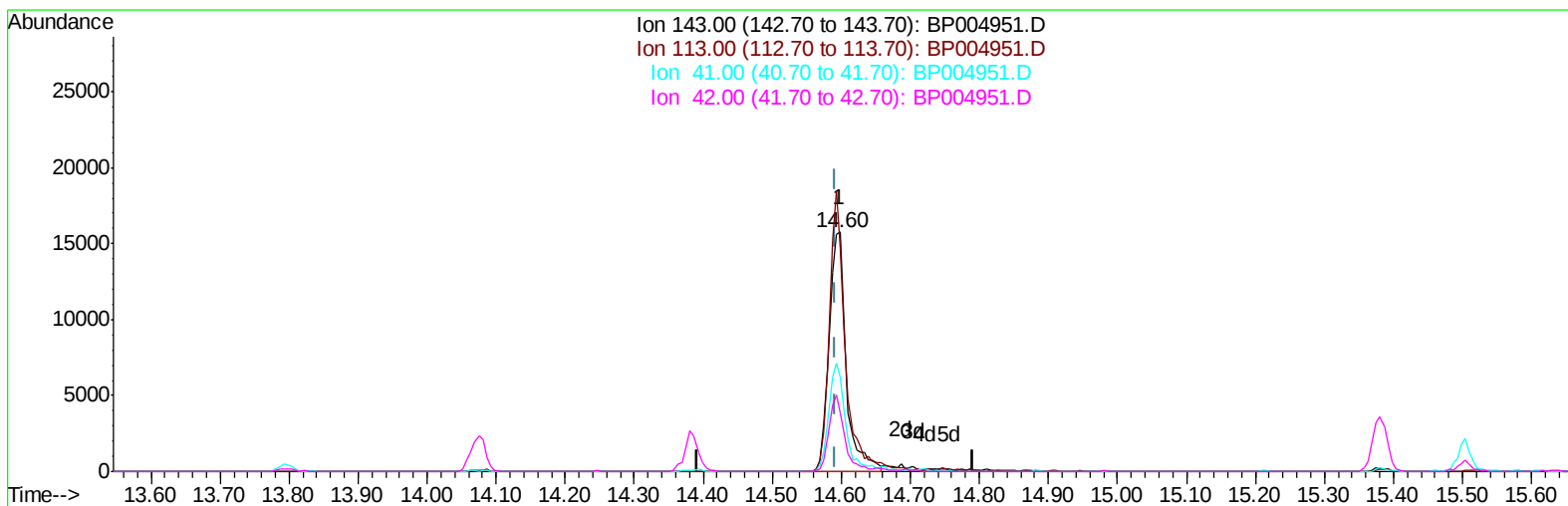
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TIC: BP004951.D

(52) 4-Nitrophenol-d4 (S)

14.598min (+0.006) 28.80ng/ul m

response 27317

Ion	Exp%	Act%
143.00	100	100
113.00	98.40	93.34
41.00	40.30	39.20
42.00	28.70	24.55

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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 09 15:13:30 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	29845	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	115583	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.38	164	75021	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.14	188	172164	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	199734	20.00	ng/ul	0.00
86) Perylene-d12	23.52	264	209762	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	4822	6.03	ng/uL	0.00
5) Phenol-d5	6.90	99	61247	26.13	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.07	67	35000	28.10	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	49986	26.99	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	49637	25.14	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	23389	27.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	26592	27.84	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	45607	24.82	ng/ul	0.00
29) 4-Chloroaniline-d4	10.66	131	64447	32.03	ng/ul	0.00
44) Dimethylphthalate-d6	13.80	166	183720	33.28	ng/ul	0.00
47) Acenaphthylene-d8	14.07	160	192127	28.51	ng/ul	0.00
52) 4-Nitrophenol-d4	14.60	143	27317m	28.80	ng/ul	0.00
58) Fluorene-d10	15.38	176	150307	31.67	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.50	200	31595	28.61	ng/ul	0.00
71) Anthracene-d10	17.24	188	258993	33.56	ng/ul	0.00
79) Pyrene-d10	19.49	212	318587	34.39	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	373462	34.82	ng/ul	0.00

Target Compounds Ovalue

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

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