

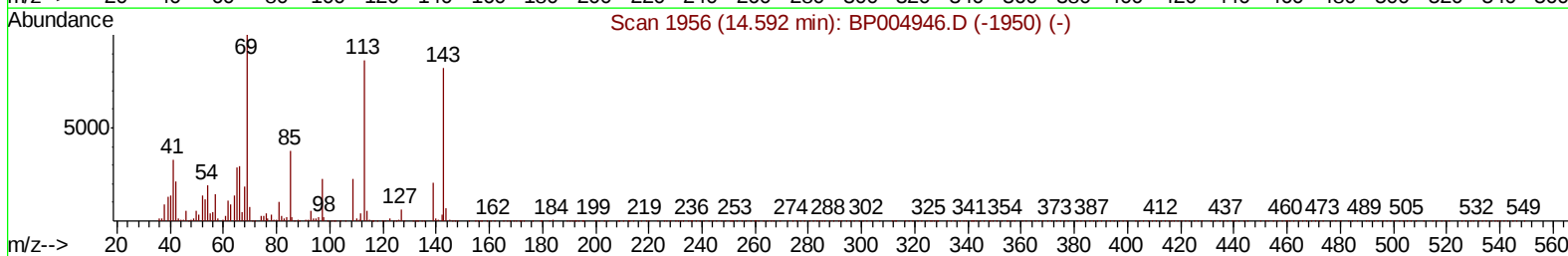
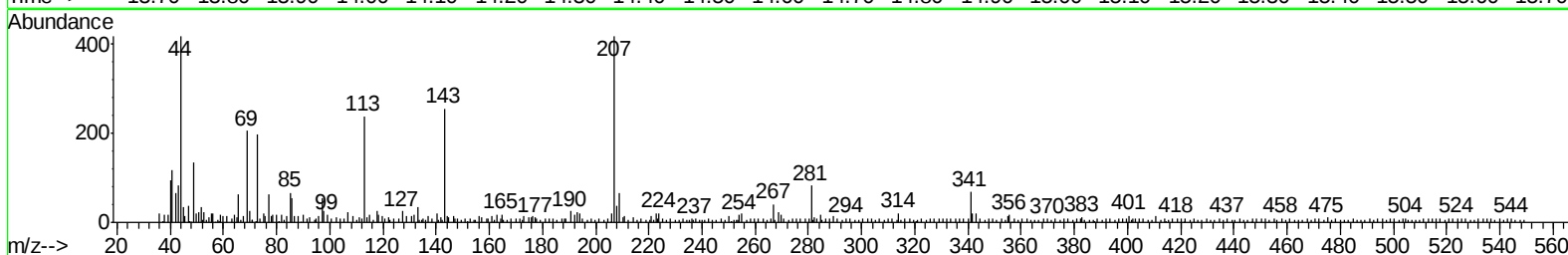
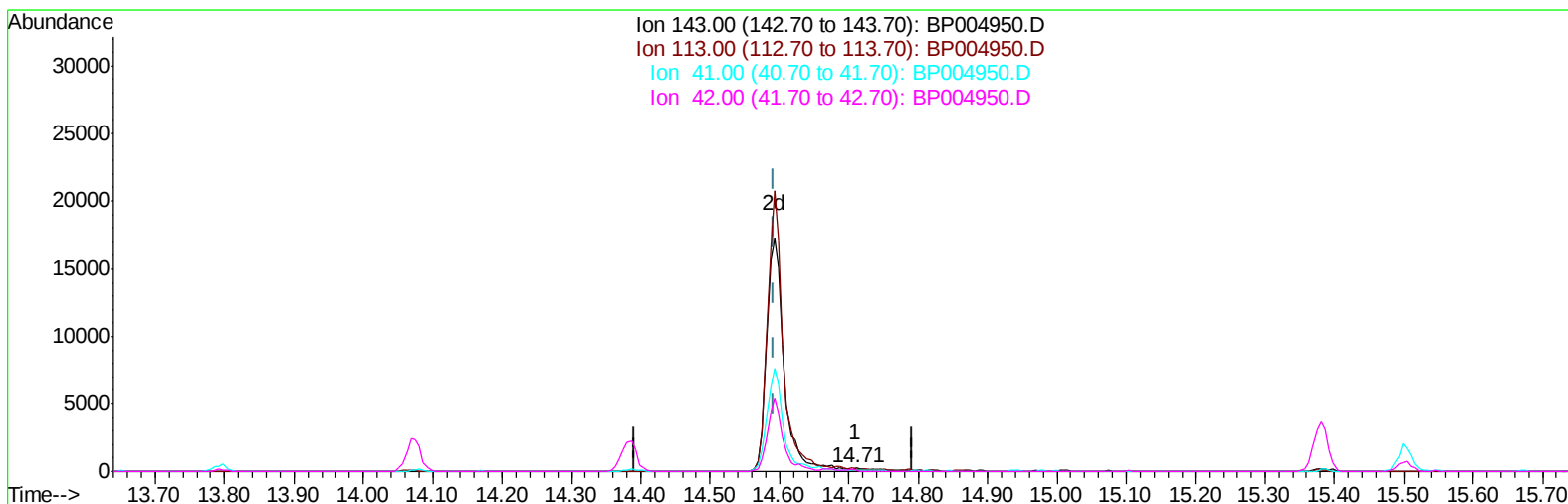
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP030921\
Data File : BP004950.D
Acq On : 09 Mar 2021 19:17
Operator : CG/JU
Sample : PB134968BL
Misc : SOM-MDL-S-PB-06
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK68

Manual Integrations
APPROVED

mohammad
3/10/2021 11:54:31 AM

Quant Time: Mar 10 00:48:32 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: BP004950.D

(52) 4-Nitrophenol-d4 (S)

14.710min (+0.118) 0.17ng/ul

response 166

Ion	Exp%	Act%
143.00	100	100
113.00	98.40	93.33
41.00	40.30	45.49
42.00	28.70	26.27

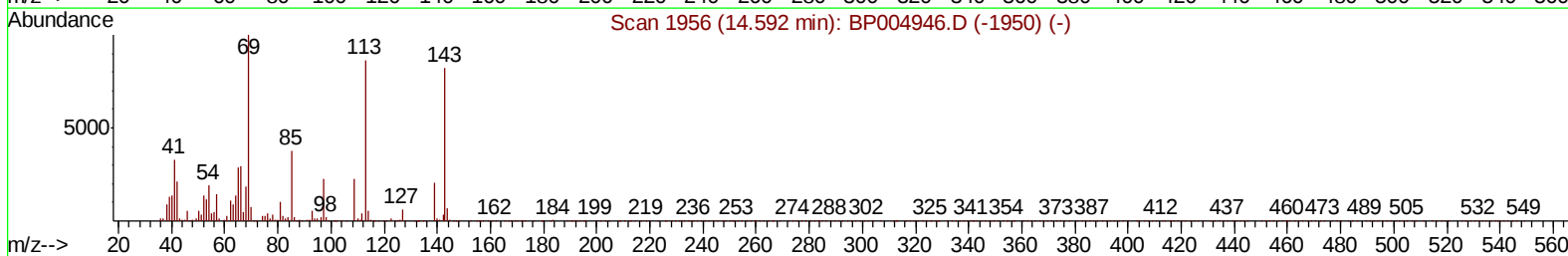
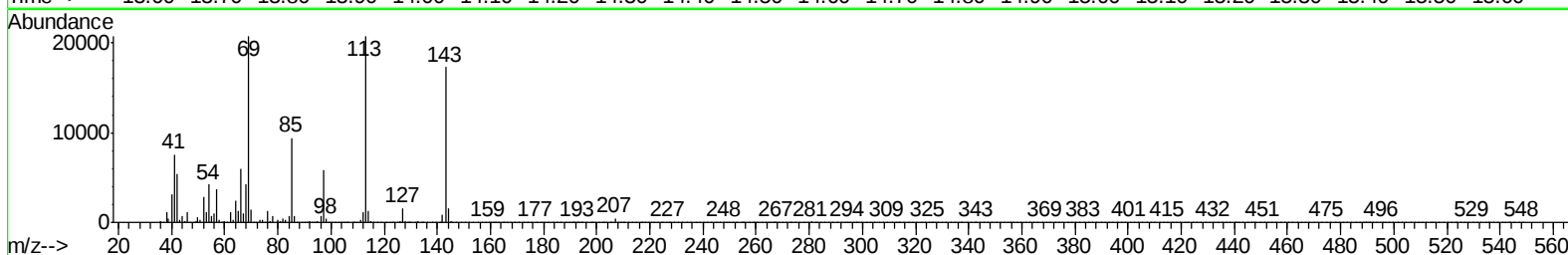
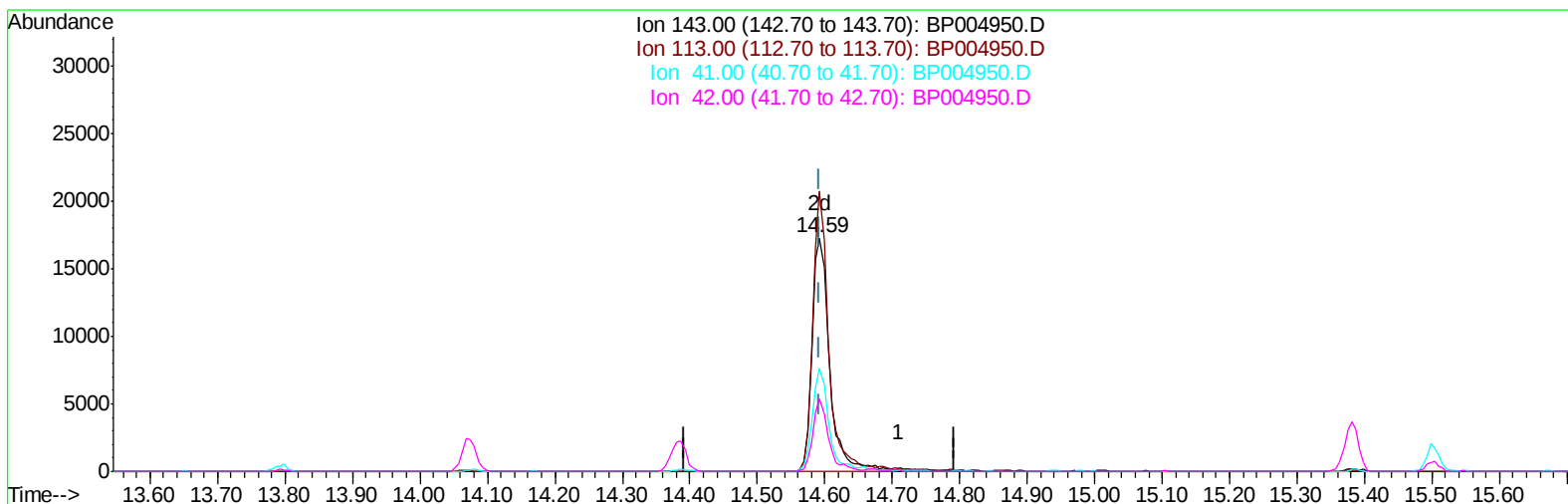
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TIC: BP004950.D

(52) 4-Nitrophenol-d4 (S)

14.592min (+0.000) 30.65ng/ul m

response 30790

Ion	Exp%	Act%
143.00	100	100
113.00	98.40	119.99#
41.00	40.30	43.94
42.00	28.70	31.30

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	33144	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	117675	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.39	164	79441	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.14	188	179706	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	204913	20.00	ng/ul	0.00
86) Perylene-d12	23.52	264	211868	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	5253	5.92	ng/uL	0.00
5) Phenol-d5	6.90	99	64767	24.88	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.07	67	36500	26.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	54489	26.49	ng/ul	0.00
13) 4-Methylphenol-d8	8.44	113	48259	22.01	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	24240	27.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	27007	27.77	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	47189	25.22	ng/ul	0.00
29) 4-Chloroaniline-d4	10.66	131	68201	33.30	ng/ul	0.00
44) Dimethylphthalate-d6	13.80	166	195656	33.47	ng/ul	0.00
47) Acenaphthylene-d8	14.07	160	206664	28.96	ng/ul	0.00
52) 4-Nitrophenol-d4	14.59	143	30790m	30.65	ng/ul	0.00
58) Fluorene-d10	15.38	176	159224	31.68	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.50	200	34366	29.82	ng/ul	0.00
71) Anthracene-d10	17.24	188	274900	34.13	ng/ul	0.00
79) Pyrene-d10	19.48	212	334036	35.15	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	380740	35.14	ng/ul	0.00

Target Compounds Ovalue

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

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