

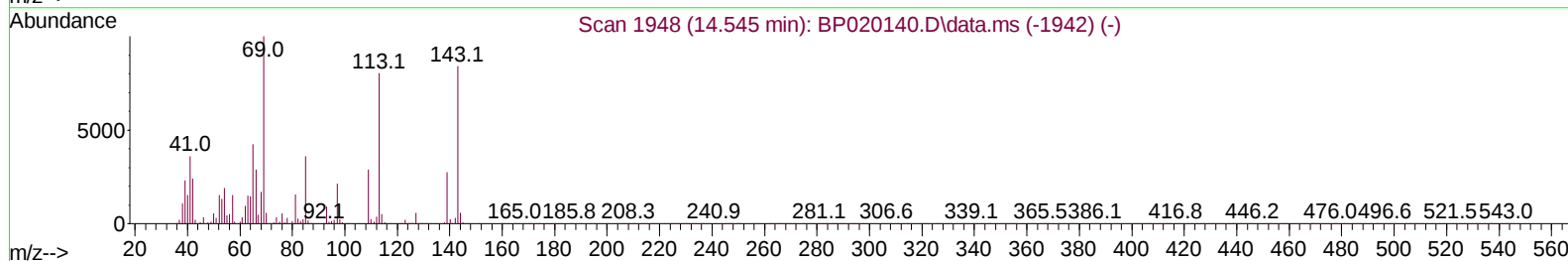
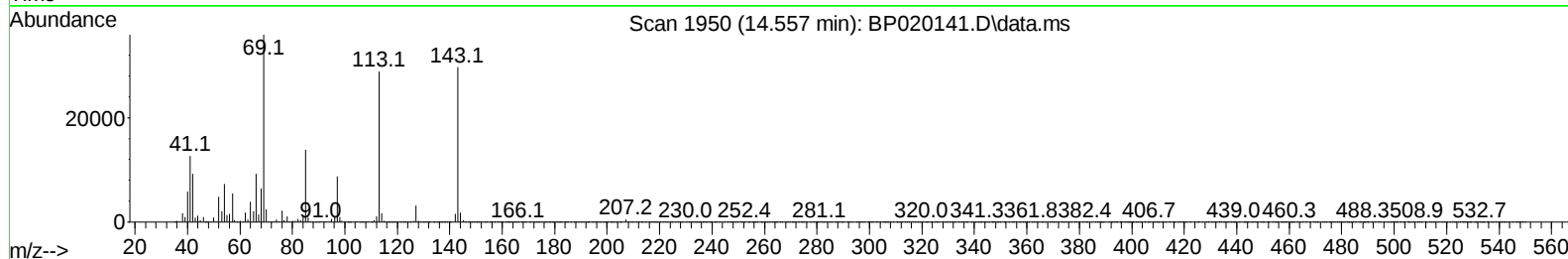
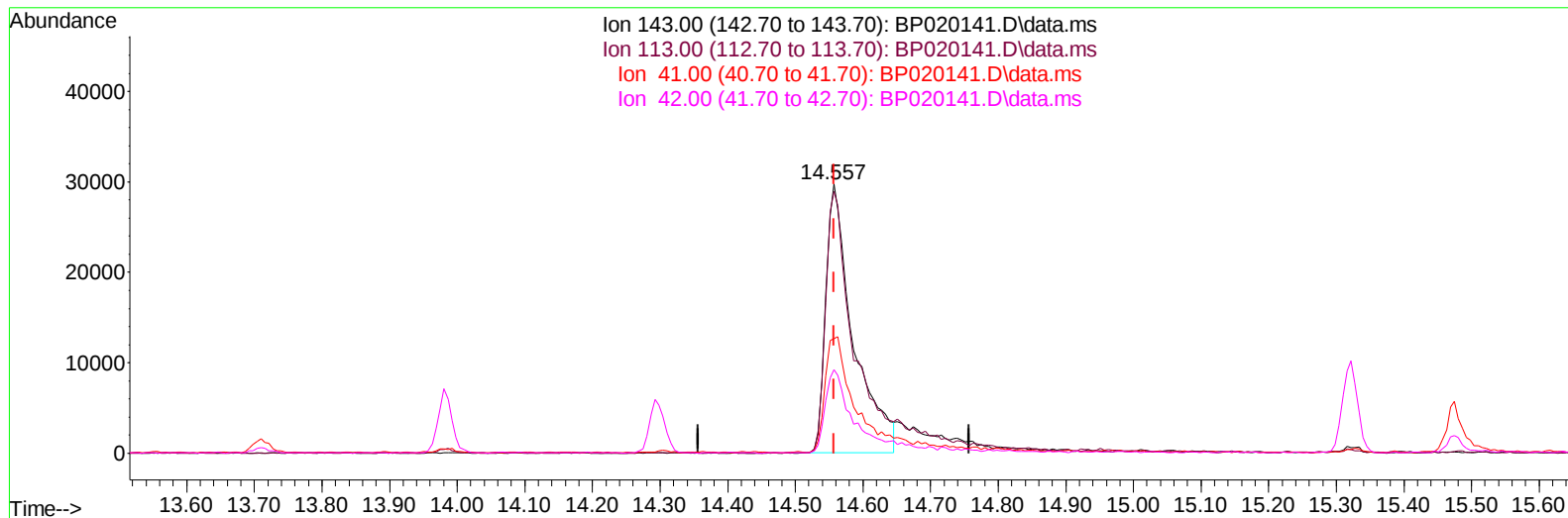
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020141.D
Acq On : 02 May 2024 10:02
Operator : MA/JU
Sample : PB160361BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK361

Manual IntegrationsAPPROVED

Quant Time: May 02 11:21:10 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 01 17:34:36 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/02/2024
Supervised By :mohammad ahmed 05/06/2024



TIC: BP020141.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (-0.000) 28.92 ng/u1

response 84171

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	97.15
41.00	50.30	42.55
42.00	32.50	31.09

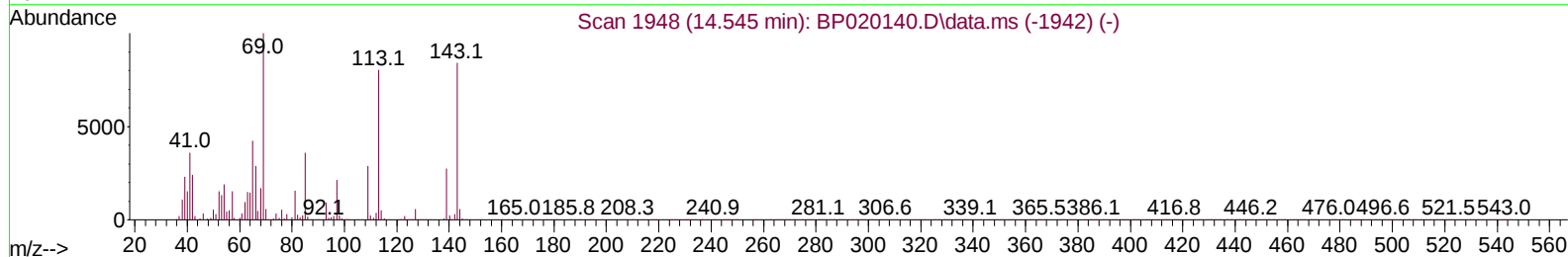
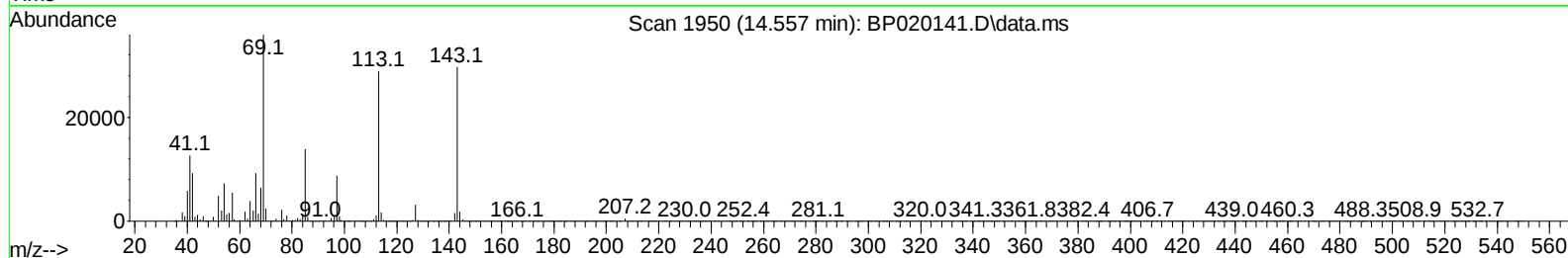
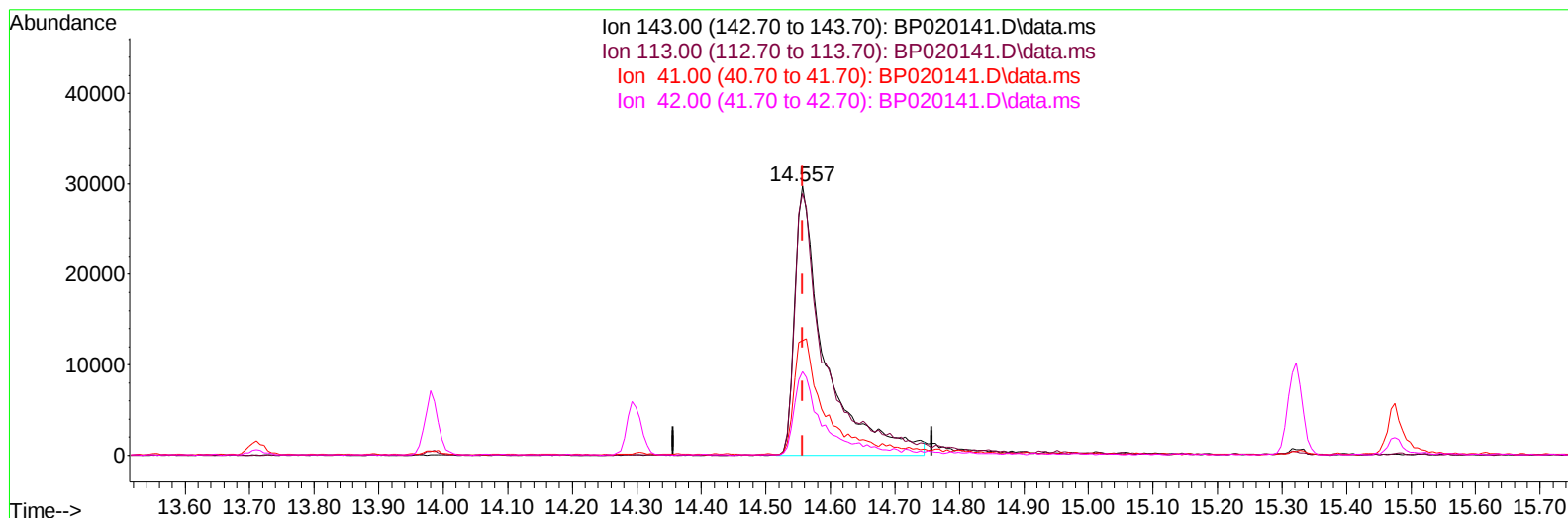
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TIC: BP020141.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (-0.000) 33.45 ng/ul m

response 97355

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	97.15
41.00	50.30	42.55
42.00	32.50	31.09

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/02/2024
 Supervised By :mohammad ahmed 05/06/2024

Quant Time: May 02 13:00:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 01 17:34:36 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	84082	20.000	ng/ul	0.00
20) Naphthalene-d8	10.398	136	355818	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.298	164	259304	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.133	188	604361	20.000	ng/ul	0.00
79) Chrysene-d12	21.639	240	596812	20.000	ng/ul	0.00
88) Perylene-d12	25.015	264	614272	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	11736	5.895	ng/uL	0.00
4) Pyridine-d5	3.628	84	131010	22.647	ng/ul	0.00
7) Phenol-d5	6.816	99	209983	29.080	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	137964	31.452	ng/ul	0.00
11) 2-Chlorophenol-d4	7.163	132	159685	30.935	ng/ul	0.00
15) 4-Methylphenol-d8	8.334	113	174321	29.860	ng/ul	0.00
21) Nitrobenzene-d5	8.787	128	85432	32.063	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	99472	32.601	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.034	165	163917	29.439	ng/ul	0.00
31) 4-Chloroaniline-d4	10.563	131	139689	17.925	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	666528	35.199	ng/ul	-0.01
49) Acenaphthylene-d8	13.981	160	679545	32.486	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.557	143	97355m	33.452	ng/ul	0.00
60) Fluorene-d10	15.322	176	537043	34.003	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.475	200	108884	28.391	ng/ul	0.00
73) Anthracene-d10	17.239	188	909634	34.792	ng/ul	0.00
81) Pyrene-d10	19.639	212	1121582	32.149	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.774	264	1029999	34.679	ng/ul	-0.01

Target Compounds Qvalue

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

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