

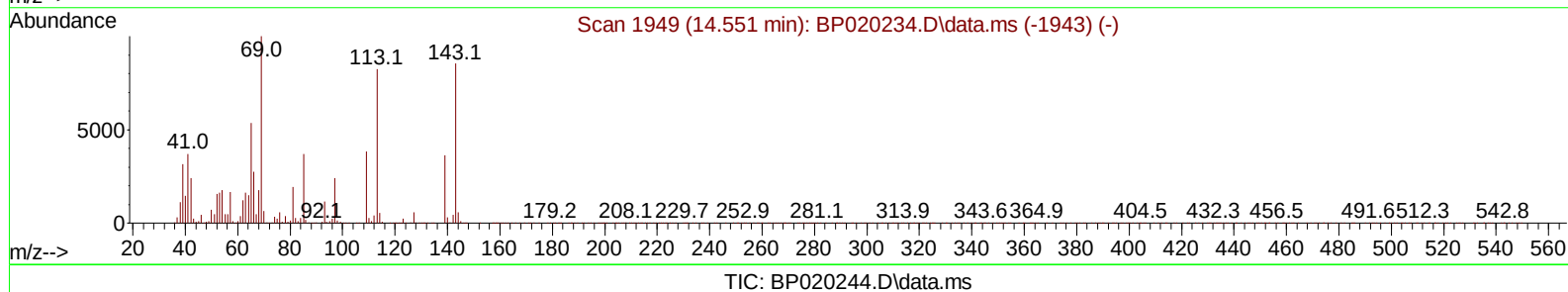
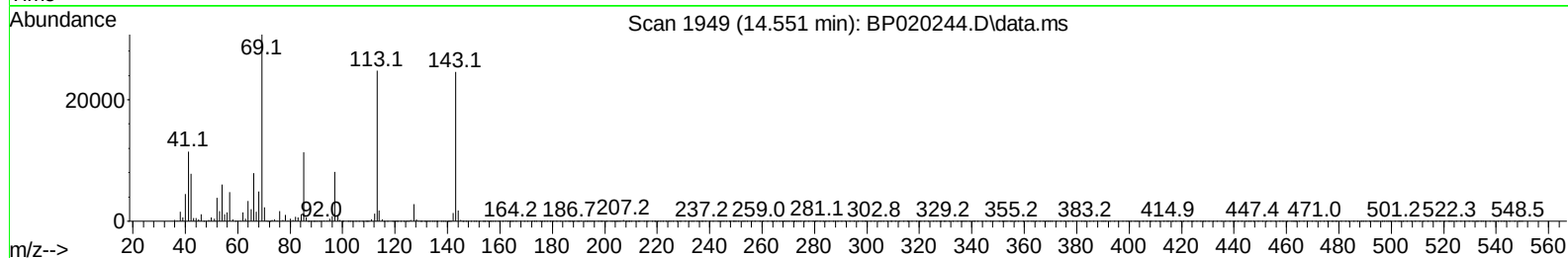
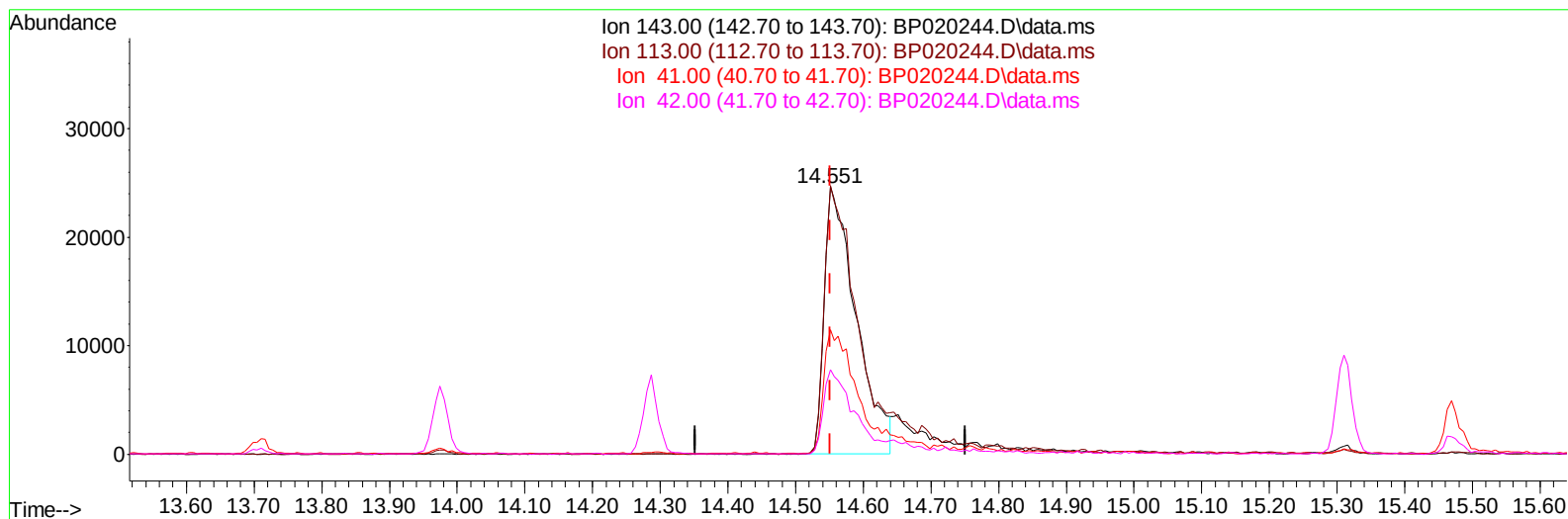
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020244.D
Acq On : 08 May 2024 19:09
Operator : MA/JU
Sample : PB160727BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK727

Manual IntegrationsAPPROVED

Quant Time: May 08 22:59:21 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 08 22:57:55 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/09/2024
Supervised By :mohammad ahmed 05/11/2024



(54) 4-Nitrophenol-d4 (S)

14.551min (-0.000) 23.97 ng/ul

response 79993

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	100.45
41.00	50.30	46.61
42.00	32.50	31.62

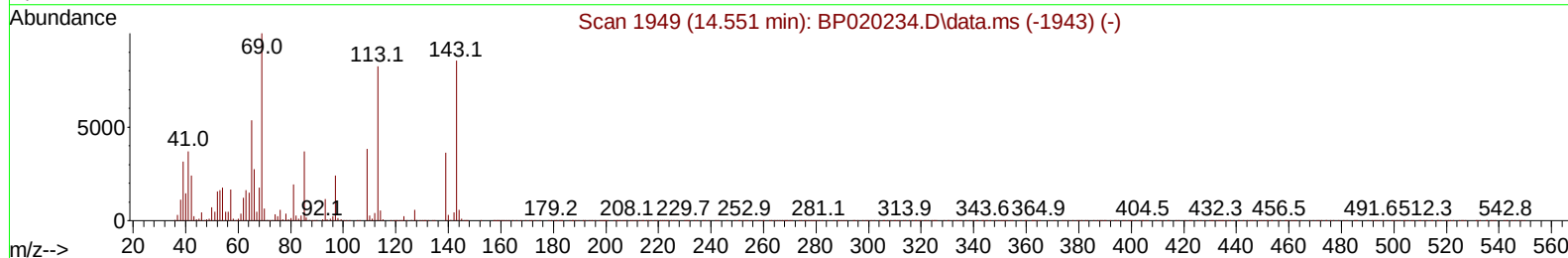
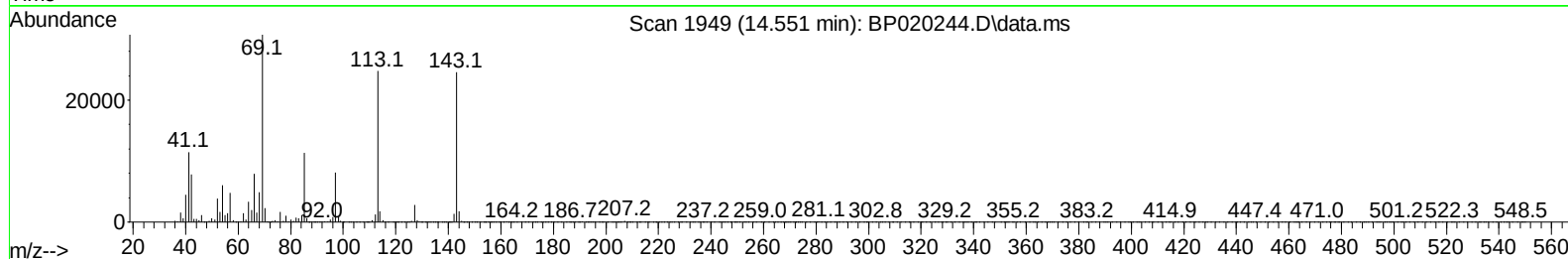
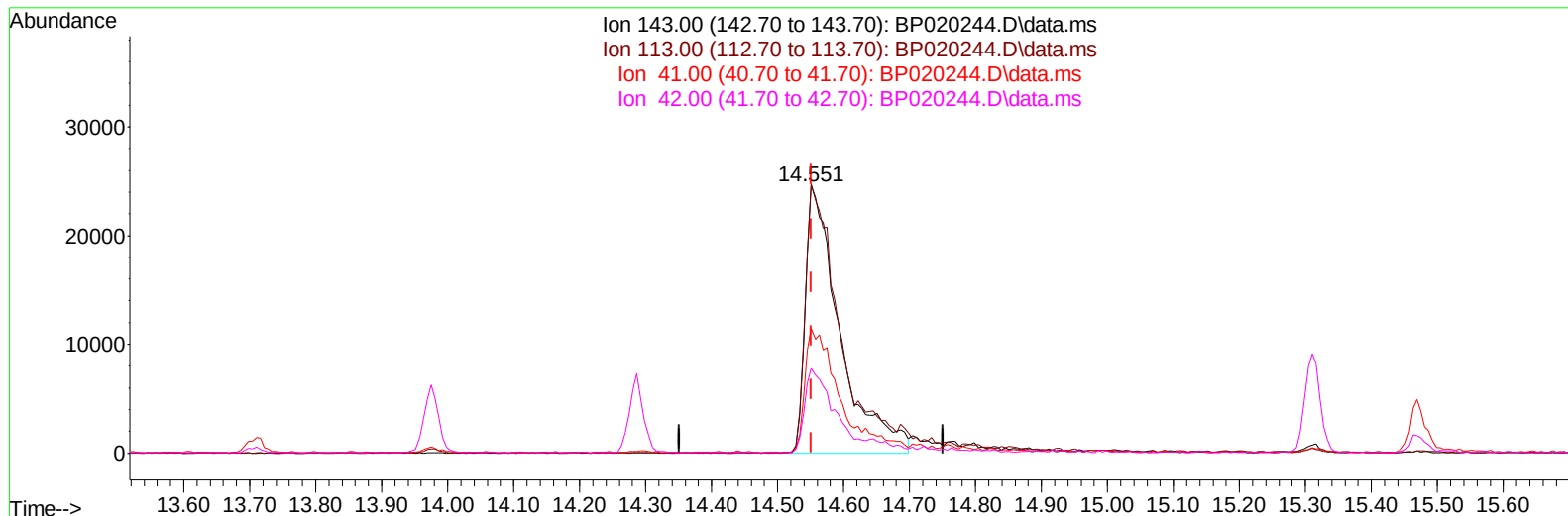
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TIC: BP020244.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.551min (-0.000) 26.56 ng/ul m

response 88647

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	100.45
41.00	50.30	46.61
42.00	32.50	31.62

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/09/2024
 Supervised By :mohammad ahmed 05/11/2024

Quant Time: May 08 23:02:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 08 22:57:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	100983	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	424416	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.287	164	297341	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.133	188	670586	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	610151	20.000	ng/ul	0.00
88) Perylene-d12	24.992	264	588548	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	11738	4.909	ng/uL	0.00
4) Pyridine-d5	3.617	84	166586	23.977	ng/ul	0.00
7) Phenol-d5	6.805	99	212620	24.517	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	135724	25.763	ng/ul	0.00
11) 2-Chlorophenol-d4	7.158	132	166286	26.822	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	169612	24.191	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	84269	26.515	ng/ul	0.00
24) 2-Nitrophenol-d4	9.493	143	99376	27.305	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	155525	23.417	ng/ul	0.00
31) 4-Chloroaniline-d4	10.546	131	253024	27.221	ng/ul	0.00
46) Dimethylphthalate-d6	13.710	166	617886	28.456	ng/ul	0.00
49) Acenaphthylene-d8	13.975	160	661502	27.578	ng/ul	0.00
54) 4-Nitrophenol-d4	14.551	143	88647m	26.563	ng/ul	0.00
60) Fluorene-d10	15.310	176	519900	28.706	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.469	200	94914	22.305	ng/ul	-0.01
73) Anthracene-d10	17.239	188	852376	29.382	ng/ul	0.00
81) Pyrene-d10	19.639	212	1054022	29.552	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.768	264	846402	29.743	ng/ul	0.00

Target Compounds Qvalue

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

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