

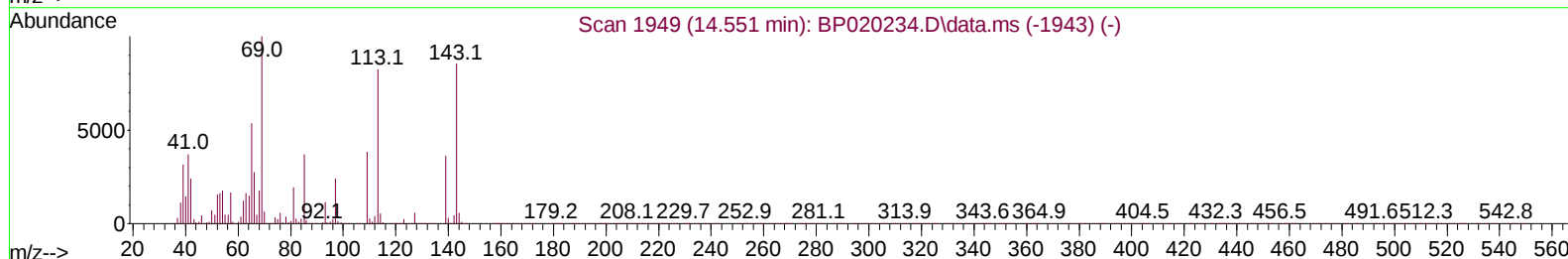
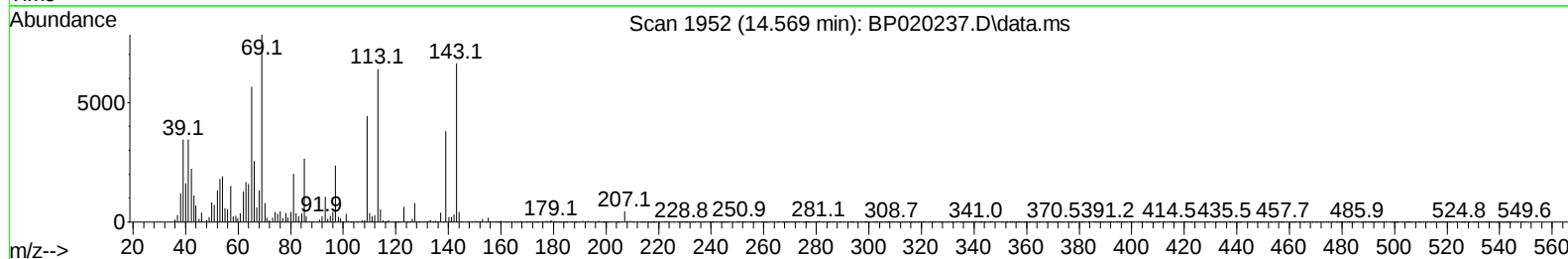
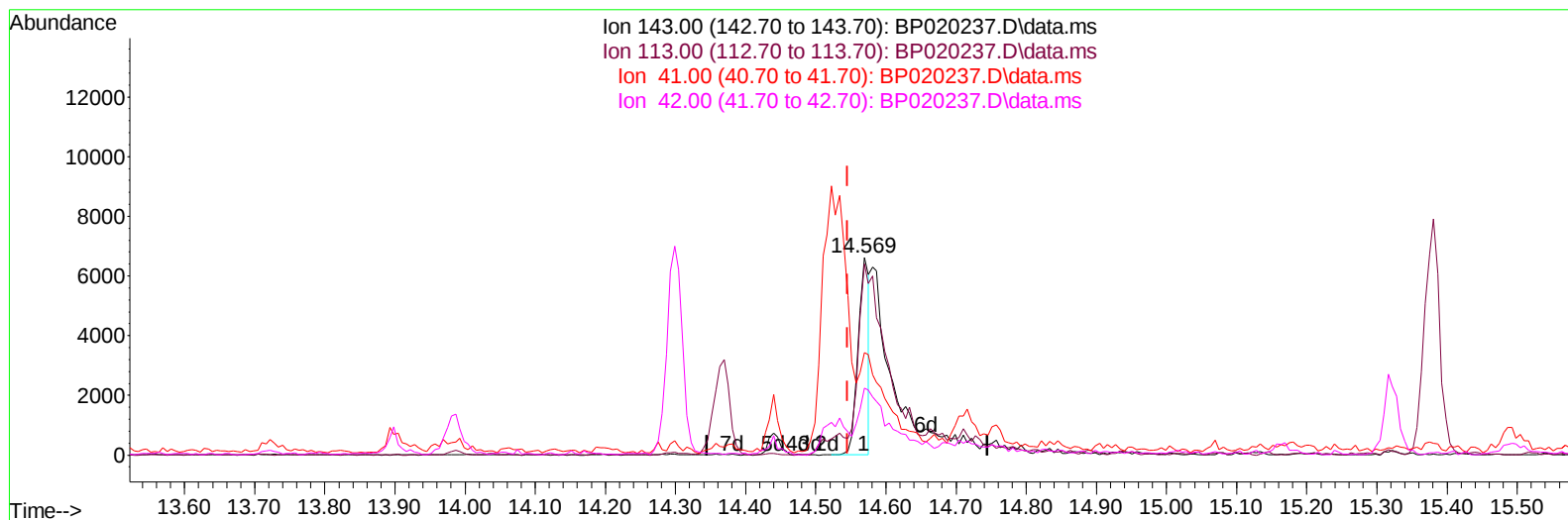
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
Data File : BP020237.D
Acq On : 08 May 2024 11:44
Operator : MA/JU
Sample : P2369-23DL 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Y5DL

Manual IntegrationsAPPROVED

Quant Time: May 08 13:26:13 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon May 06 12:26:12 2024
Response via : Initial Calibration

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



TIC: BP020237.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (+ 0.024) 2.10 ng/u1

response 7355

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	96.70
41.00	50.30	51.92
42.00	32.50	33.74

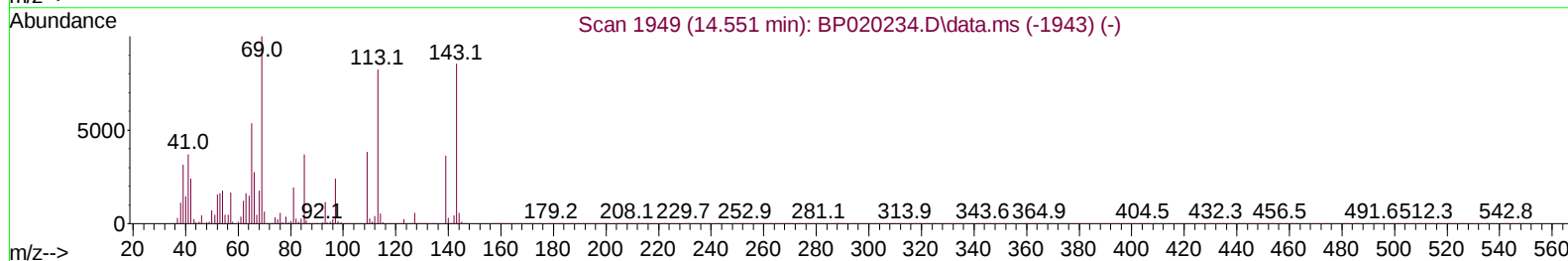
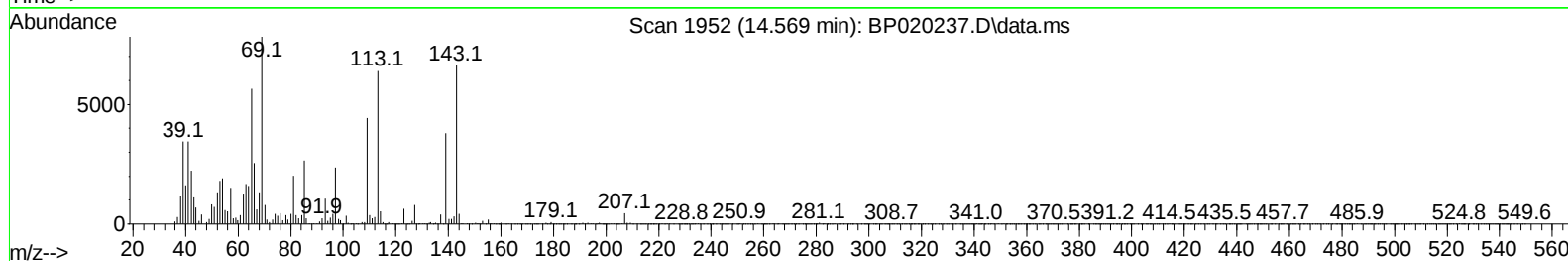
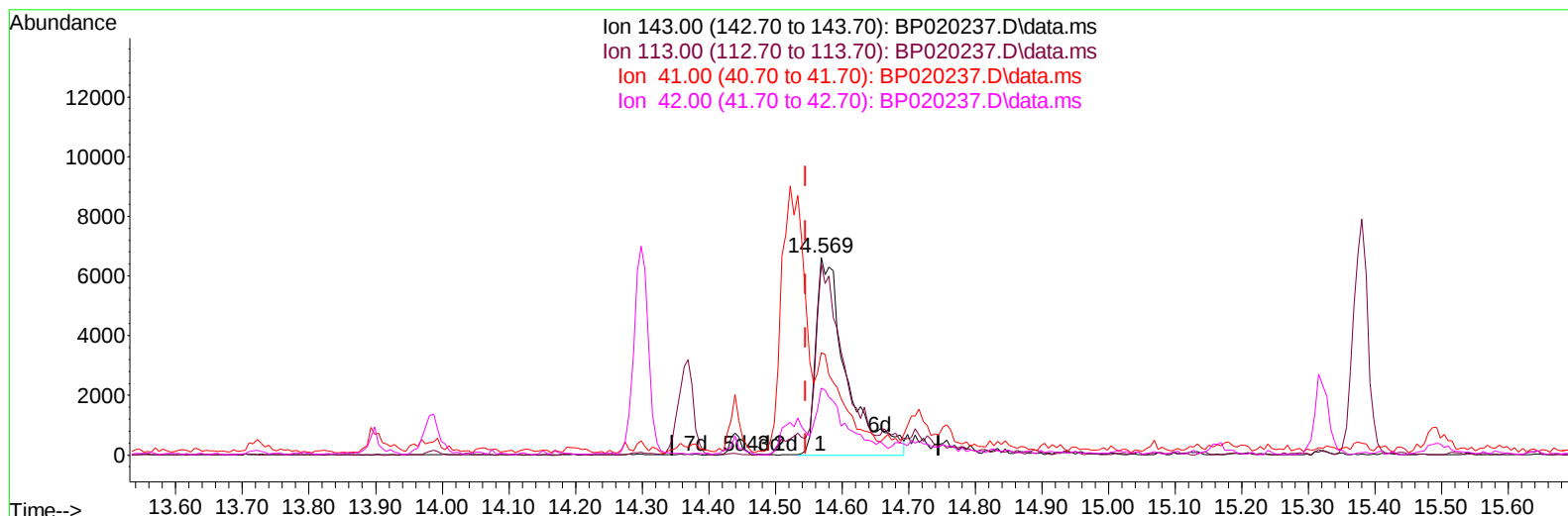
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TIC: BP020237.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (+ 0.024) 6.00 ng/ul m

response 21032

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	96.70
41.00	50.30	51.92
42.00	32.50	33.74

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Manual IntegrationsAPPROVED

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Quant Time: May 08 13:34:16 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.605	152		106410	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136		450520	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.298	164		312480	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.145	188		696219	20.000	ng/ul	0.00
79) Chrysene-d12	21.657	240		651320	20.000	ng/ul	0.02
88) Perylene-d12	25.033	264		653731	20.000	ng/ul	0.04
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.217	96		1940	0.770	ng/uL	0.00
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	6.793	99		40193	4.398	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.958	67		27079	4.878	ng/ul	0.00
11) 2-Chlorophenol-d4	7.146	132		32201	4.929	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113		33398	4.520	ng/ul	0.01
21) Nitrobenzene-d5	8.769	128		16863	4.998	ng/ul	0.01
24) 2-Nitrophenol-d4	9.481	143		18344	4.748	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.016	165		45077	6.394	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131		39023	3.955	ng/ul	0.02
46) Dimethylphthalate-d6	13.722	166		142577	6.248	ng/ul	0.02
49) Acenaphthylene-d8	13.987	160		144628	5.737	ng/ul	0.00
54) 4-Nitrophenol-d4	14.569	143		21032m	5.997	ng/ul	0.02
60) Fluorene-d10	15.322	176		122130	6.417	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.493	200		17421	3.943	ng/ul	0.02
73) Anthracene-d10	17.251	188		204464	6.789	ng/ul	0.01
81) Pyrene-d10	19.657	212		260010	6.829	ng/ul	0.02
92) Benzo(a)pyrene-d12	24.804	264		212483	6.722	ng/ul	0.04
Target Compounds							
						Qvalue	
71) Pentachlorophenol	16.787	266		172554	32.684	ng/ul	98

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

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