

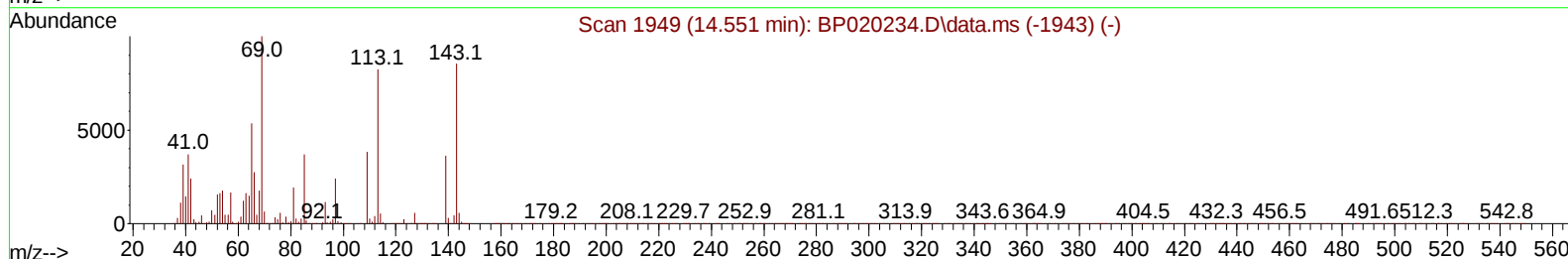
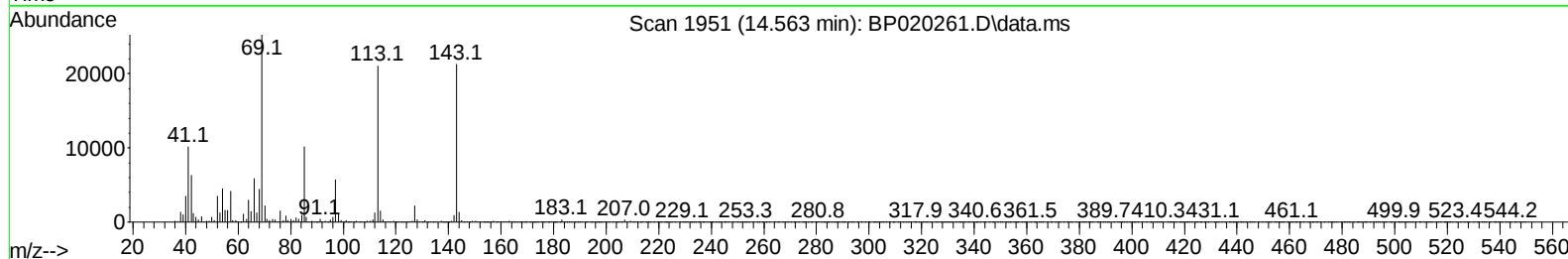
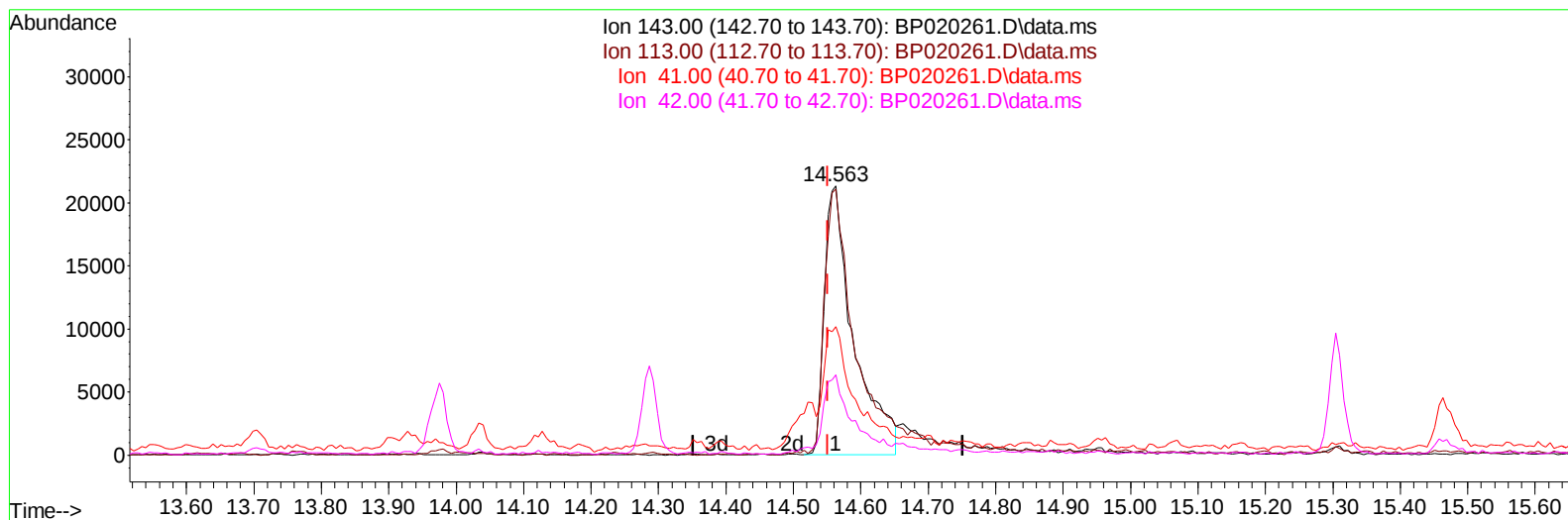
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
 Data File : BP020261.D
 Acq On : 09 May 2024 08:32
 Operator : MA/JU
 Sample : P2369-10
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43N4

Manual IntegrationsAPPROVED

Quant Time: May 09 08:59:40 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



TIC: BP020261.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.012) 19.21 ng/u1

response 64315

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	98.95
41.00	50.30	47.82
42.00	32.50	29.73

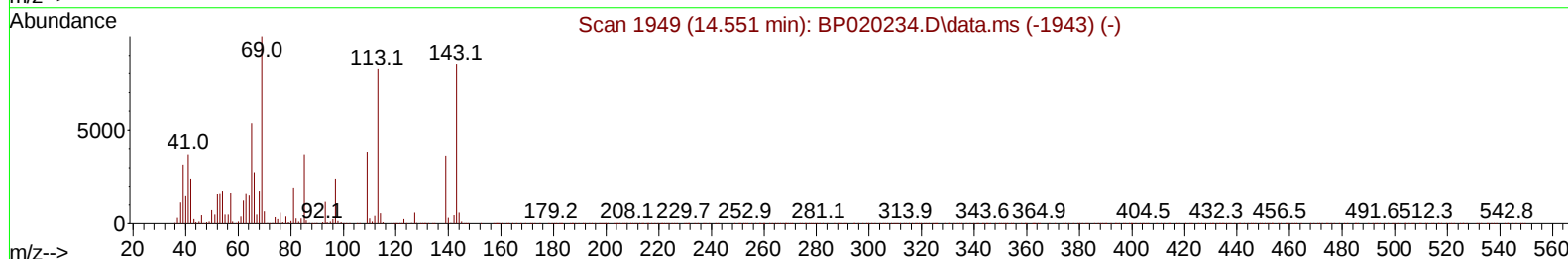
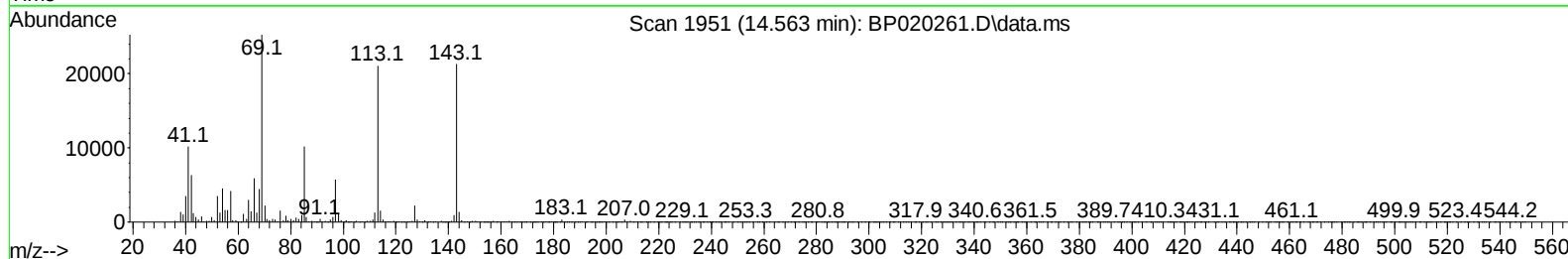
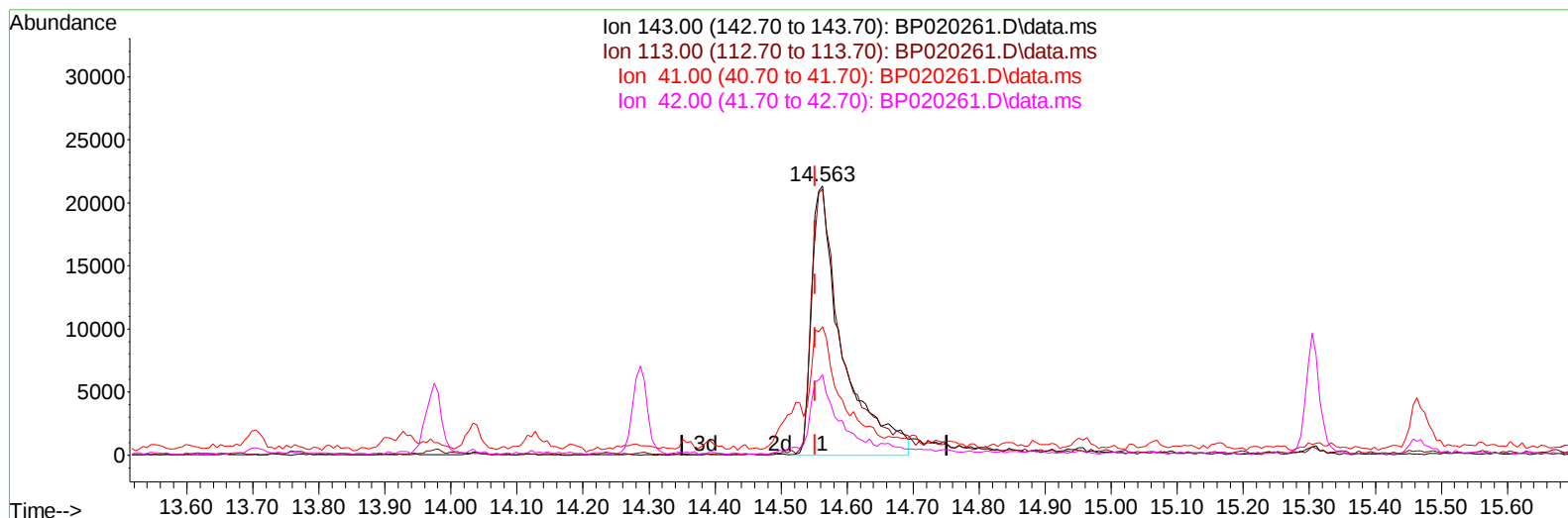
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(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.012) 20.69 ng/ul m

response 69277

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	98.95
41.00	50.30	47.82
42.00	32.50	29.73

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.616	152		94589	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136		409542	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.287	164		298332	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.122	188		643721	20.000	ng/ul	-0.02
79) Chrysene-d12	21.633	240		511264	20.000	ng/ul	0.00
88) Perylene-d12	24.998	264		522311	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.217	96		8617	3.847	ng/uL	0.00
4) Pyridine-d5	3.622	84		104943	16.126	ng/ul	0.00
7) Phenol-d5	6.811	99		180180	22.181	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67		107592	21.804	ng/ul	0.00
11) 2-Chlorophenol-d4	7.158	132		137198	23.626	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113		148602	22.627	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128		71764	23.401	ng/ul	0.00
24) 2-Nitrophenol-d4	9.493	143		83665	23.823	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165		152019	23.721	ng/ul	0.00
31) 4-Chloroaniline-d4	10.552	131		187564	20.911	ng/ul	0.00
46) Dimethylphthalate-d6	13.704	166		589433	27.056	ng/ul	0.00
49) Acenaphthylene-d8	13.975	160		626478	26.031	ng/ul	0.00
54) 4-Nitrophenol-d4	14.563	143		69277m	20.690	ng/ul	0.01
60) Fluorene-d10	15.304	176		499030	27.462	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.463	200		78203	19.145	ng/ul	-0.02
73) Anthracene-d10	17.222	188		771541	27.706	ng/ul	-0.02
81) Pyrene-d10	19.639	212		884988	29.612	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.763	264		699032	27.680	ng/ul	-0.01
Target Compounds							
						Qvalue	
71) Pentachlorophenol	16.763	266		452197	92.637	ng/ul	99

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

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