

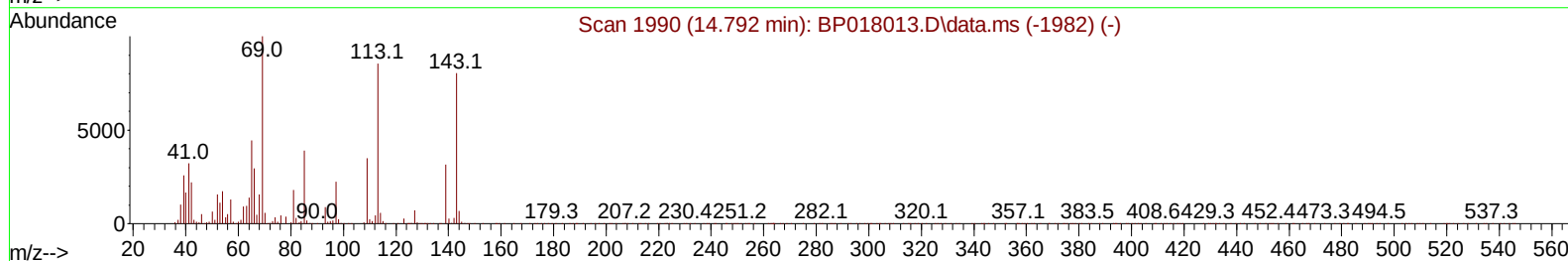
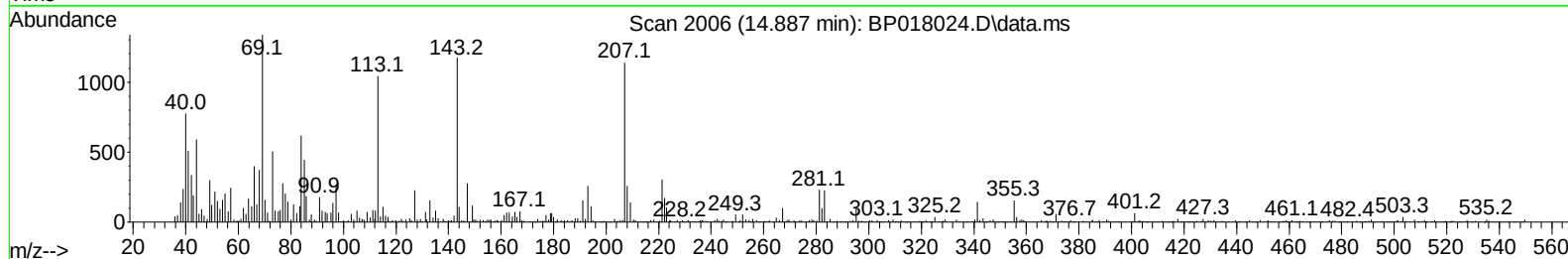
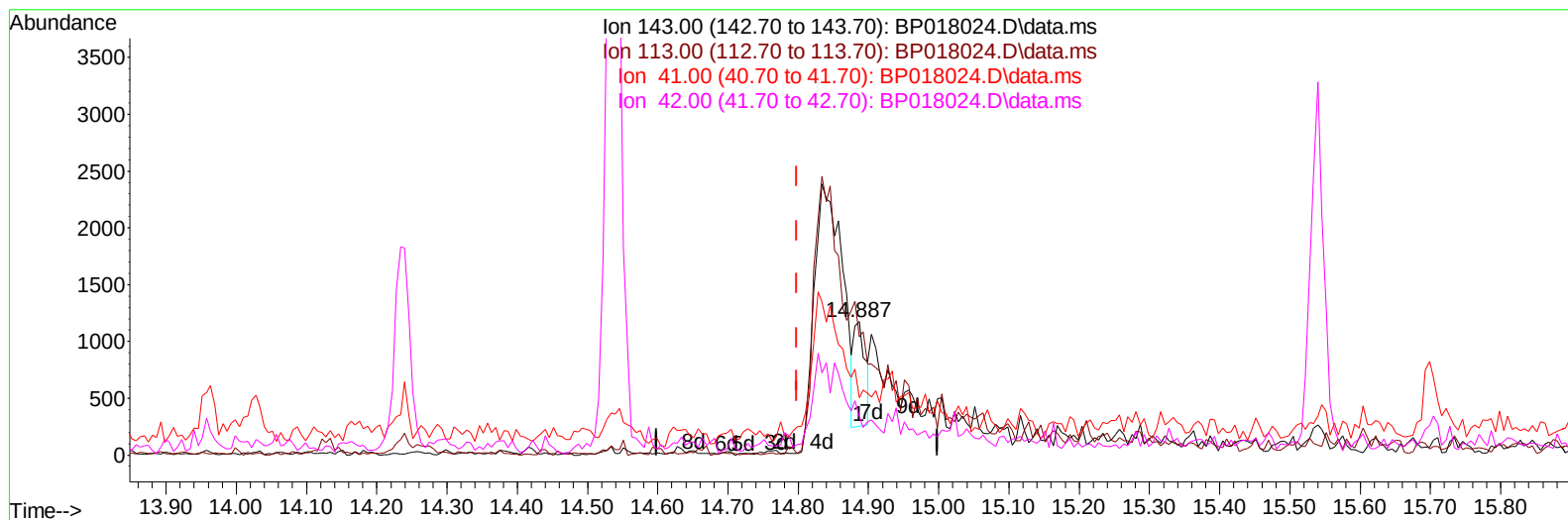
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
 Data File : BP018024.D
 Acq On : 10 Nov 2023 11:57
 Operator : MA/JU
 Sample : 05091-11
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCK18

Manual IntegrationsAPPROVED

Quant Time: Nov 10 22:17:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 09 22:34:38 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/11/2023
 Supervised By :mohammad ahmed 11/15/2023



TIC: BP018024.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.088) 0.34 ng/ul

response 1048

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	88.53
41.00	46.90	43.16
42.00	29.70	28.55

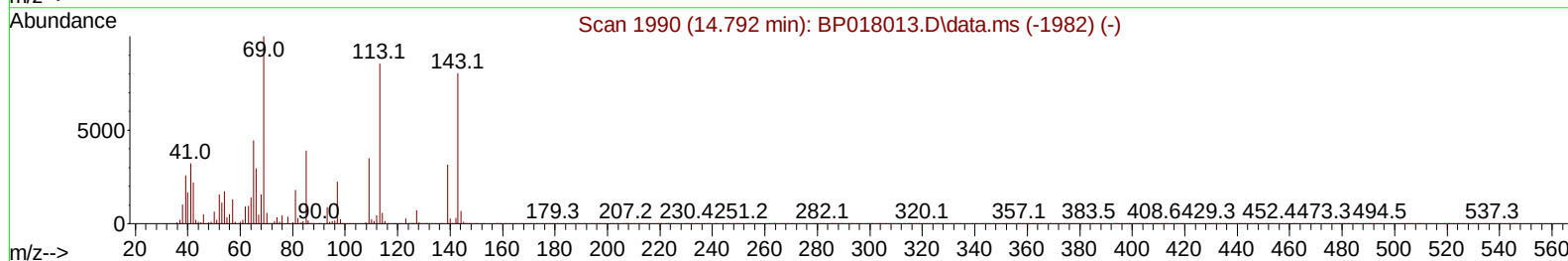
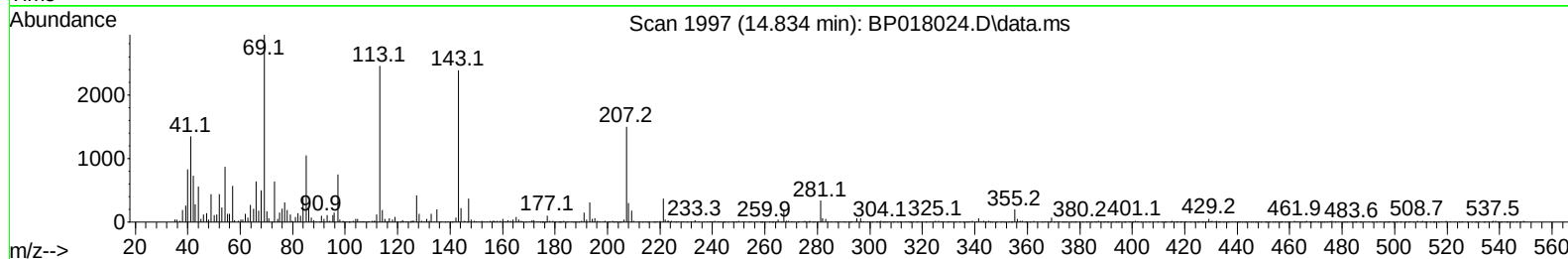
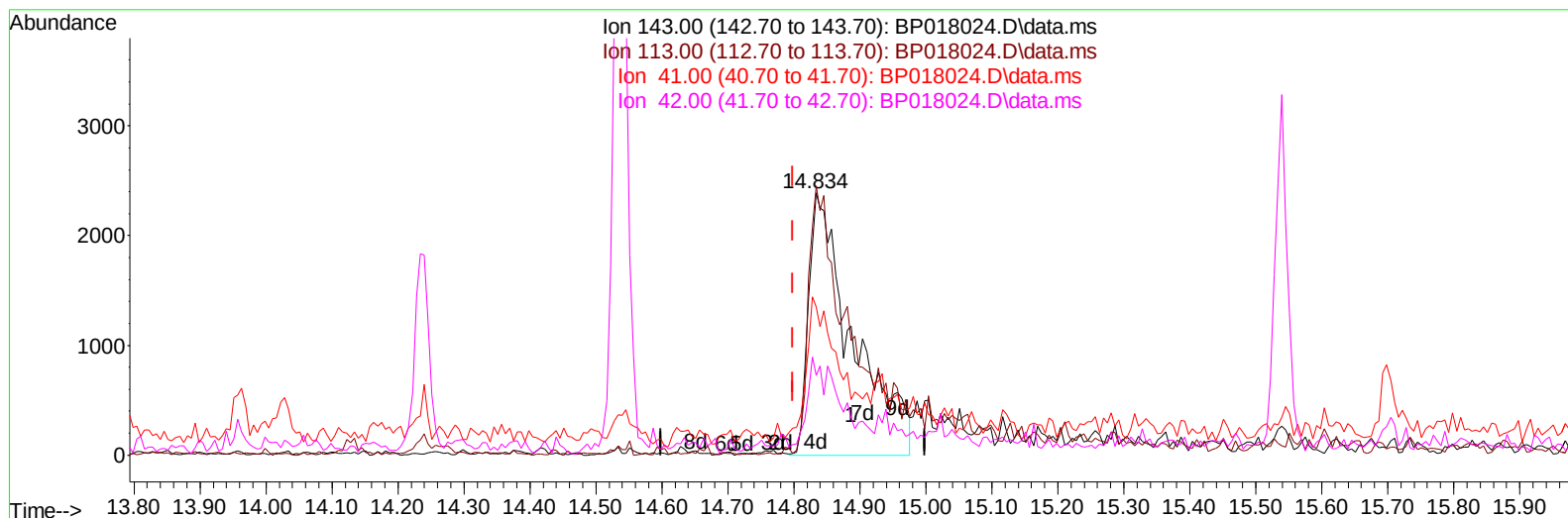
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TIC: BP018024.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (+ 0.035) 3.62 ng/ul m

response 11034

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	102.60
41.00	46.90	56.38#
42.00	29.70	30.35

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	87520	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	359188	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.540	164	226623	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.316	188	513029	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240	498033	20.000	ng/ul	0.00
88) Perylene-d12	23.957	264	548559	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	2490	0.974	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.034	99	49755	6.049	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	35191	7.063	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	44911	6.825	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	19958	3.025	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	22269	6.923	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	24346	6.801	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	36309	5.662	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	33582	3.821	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	151567	7.275	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	163558	7.323	ng/ul	0.00
54) 4-Nitrophenol-d4	14.834	143	11034m	3.616	ng/ul	0.04
60) Fluorene-d10	15.540	176	130264	7.544	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	9339	2.607	ng/ul	0.00
73) Anthracene-d10	17.416	188	211064	7.678	ng/ul	0.00
81) Pyrene-d10	19.669	212	285862	8.918	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.786	264	250475	7.856	ng/ul	-0.01

Target Compounds Qvalue

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

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