

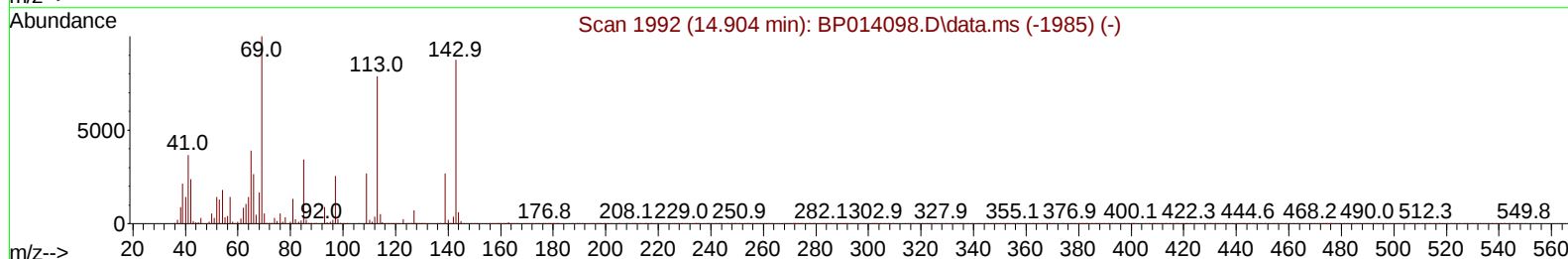
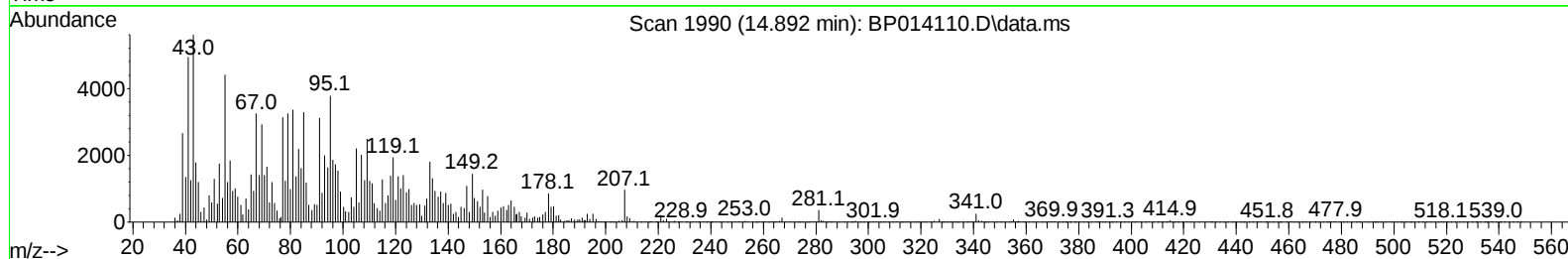
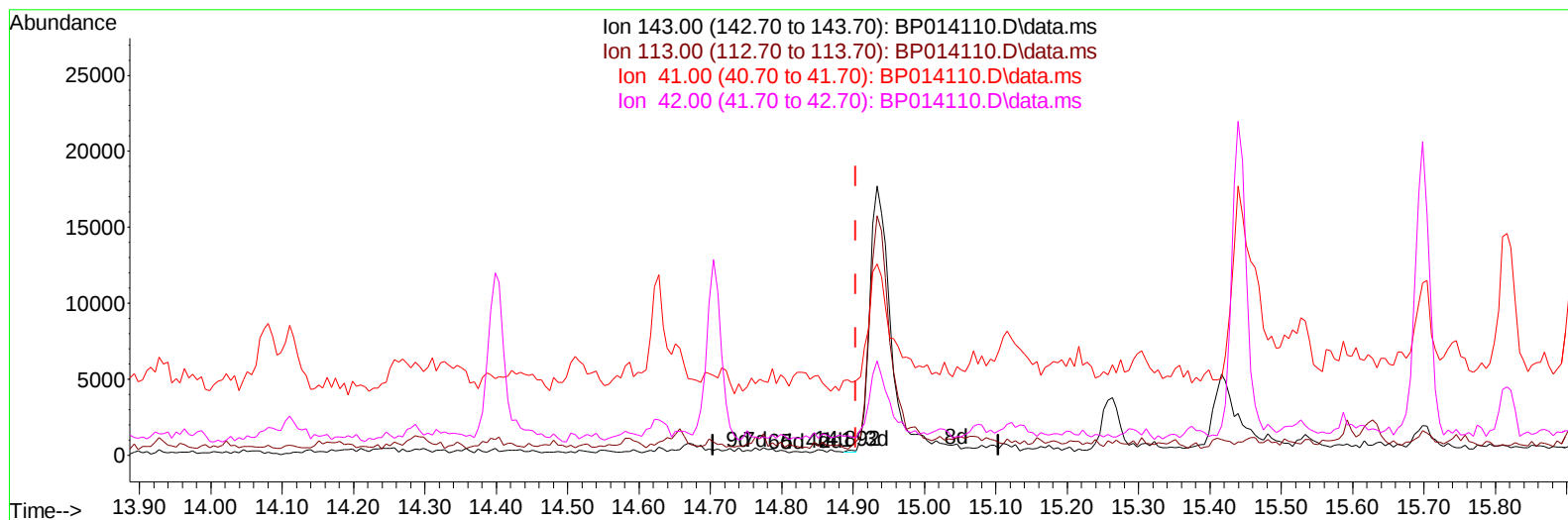
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
 Data File : BP014110.D
 Acq On : 16 Mar 2023 19:22
 Operator : CG/JU
 Sample : 01884-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB927

Manual IntegrationsAPPROVED

Quant Time: Mar 17 01:01:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 17 00:57:25 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/17/2023
 Supervised By :Jagrut Upadhyay 03/17/2023



TIC: BP014110.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.892min (-0.012) 0.01 ng/u1

response 98

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	139.13#
41.00	40.20	1650.17#
42.00	28.50	422.74#

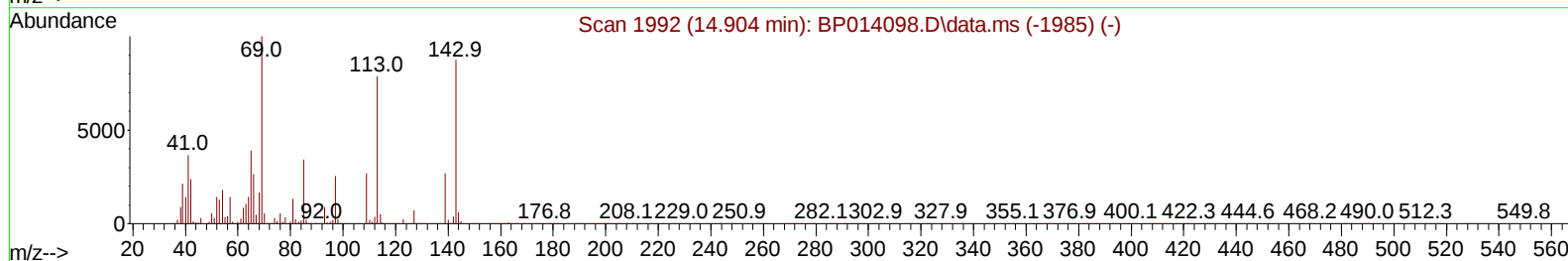
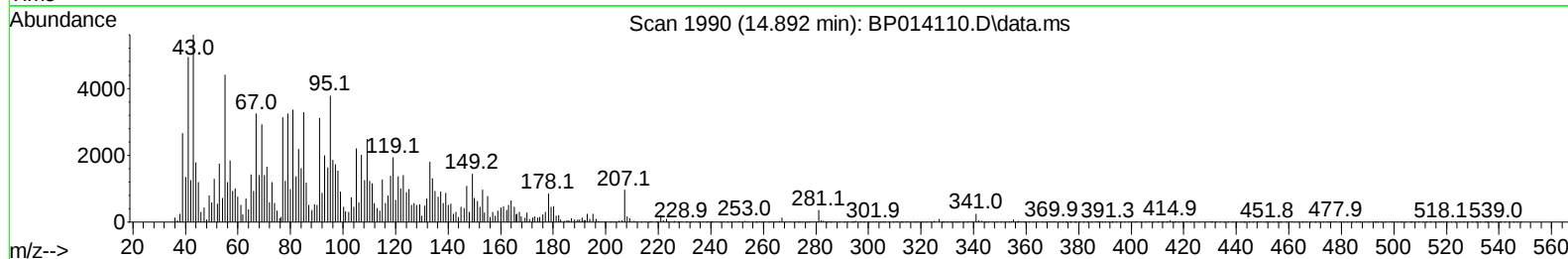
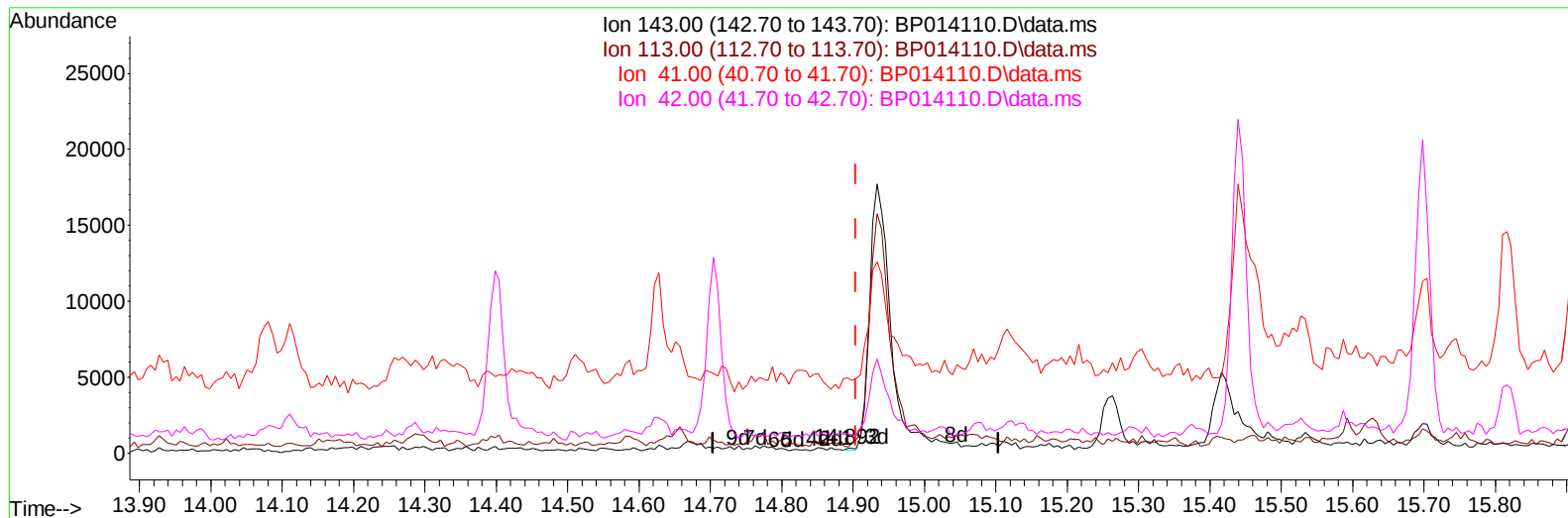
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(54) 4-Nitrophenol-d4 (S)

14.892min (-0.012) 0.01 ng/u1

response 98

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	139.13#
41.00	40.20	1650.17#
42.00	28.50	422.74#

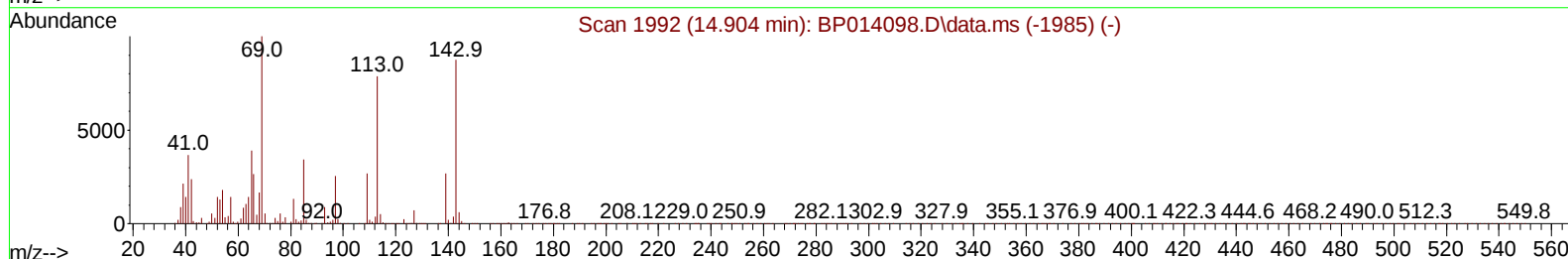
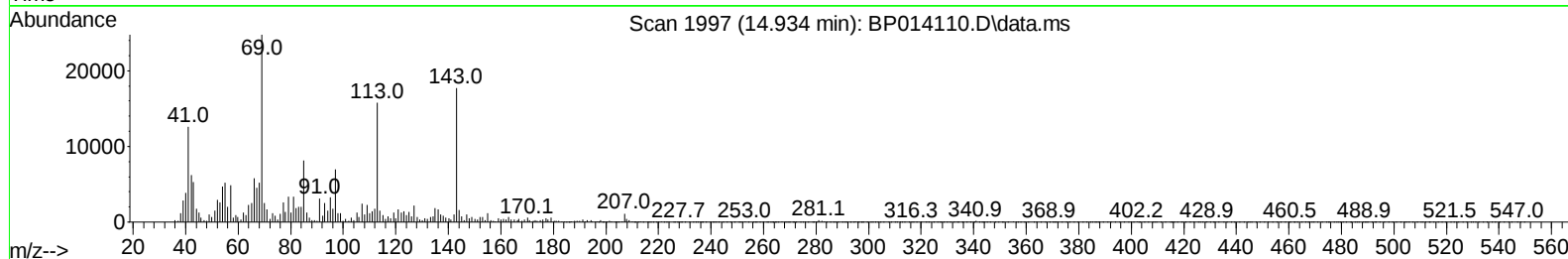
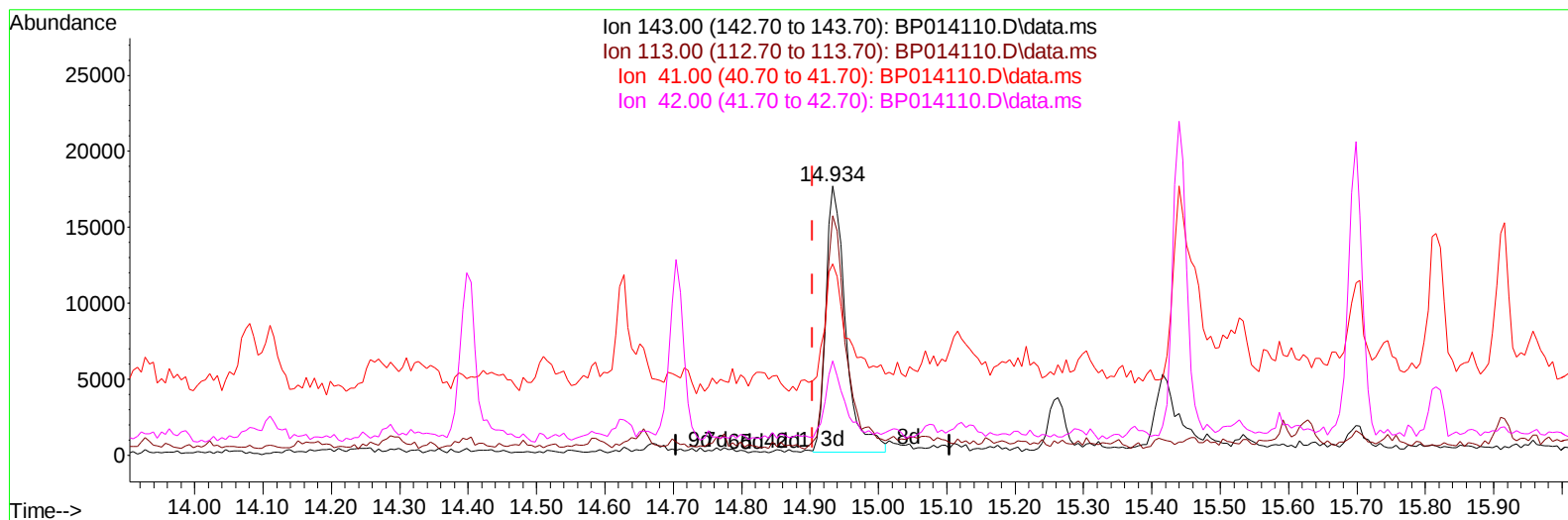
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TIC: BP014110.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.029) 4.39 ng/ul m

response 35405

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	88.96
41.00	40.20	71.02#
42.00	28.50	35.00#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
 Data File : BP014110.D
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 Operator : CG/JU
 Sample : 01884-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.063	152	189003	20.000	ng/ul	0.00
20) Naphthalene-d8	10.887	136	795493	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	539300	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	1202426	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.545	240	1202720	20.000	ng/ul	# 0.00
88) Perylene-d12	24.080	264	1221577	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.428	96	20504	4.099	ng/uL	0.00
4) Pyridine-d5	3.864	84	175853	12.697	ng/ul	0.00
7) Phenol-d5	7.222	99	128277	7.691	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.387	67	300025	30.844	ng/ul	0.00
11) 2-Chlorophenol-d4	7.593	132	268983	21.167	ng/ul	0.00
15) 4-Methylphenol-d8	8.775	113	260084	19.090	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	182841	28.807	ng/ul	0.00
24) 2-Nitrophenol-d4	9.963	143	190493	25.072	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.504	165	350235	24.498	ng/ul	0.00
31) 4-Chloroaniline-d4	11.028	131	423927	22.407	ng/ul	0.00
46) Dimethylphthalate-d6	14.110	166	1320313	28.847	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160	1385425	28.574	ng/ul	0.00
54) 4-Nitrophenol-d4	14.934	143	35405m	4.390	ng/ul	0.03
60) Fluorene-d10	15.698	176	1043625	26.905	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.816	200	166141	18.281	ng/ul	0.00
73) Anthracene-d10	17.563	188	1717791	29.428	ng/ul	0.00
81) Pyrene-d10	19.786	212	2036651	30.703	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.915	264	1913875	29.454	ng/ul	0.00
Target Compounds						
52) Acenaphthene	14.769	153	78865	2.217	ng/ul	Qvalue 98

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

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