

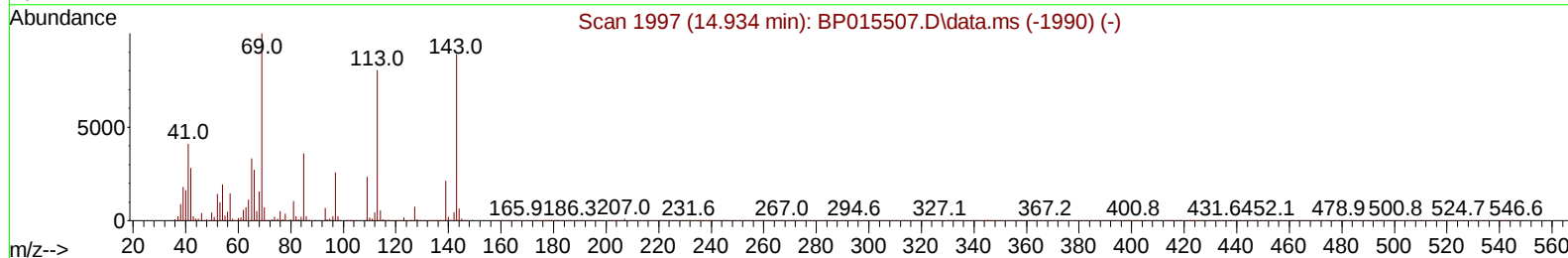
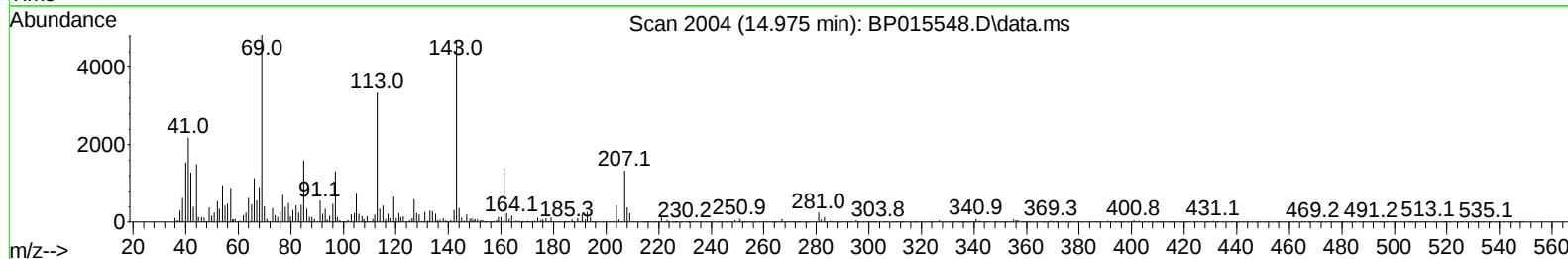
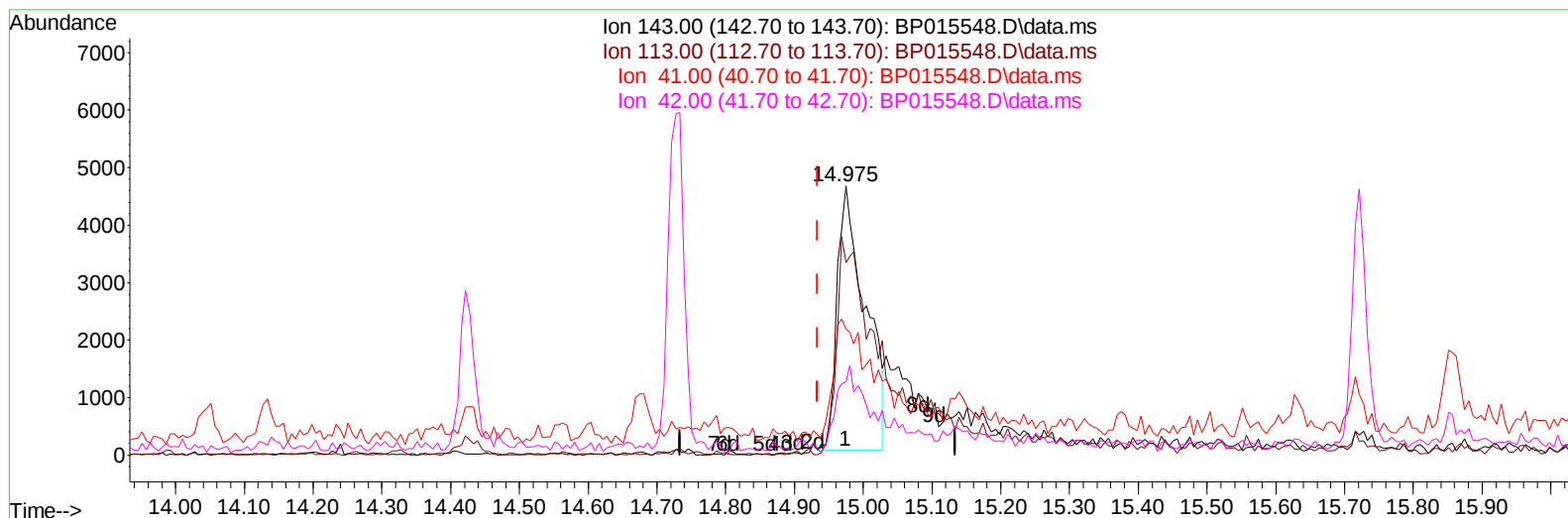
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015548.D
 Acq On : 15 Jun 2023 08:53
 Operator : MA/JU
 Sample : 03088-07
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFR3

Manual IntegrationsAPPROVED

Quant Time: Jun 15 09:22:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 15 05:36:46 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/15/2023
 Supervised By :mohammad ahmed 06/16/2023



TIC: BP015548.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.041) 3.44 ng/ul

response 13059

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	71.48
41.00	42.20	46.86
42.00	29.40	27.26

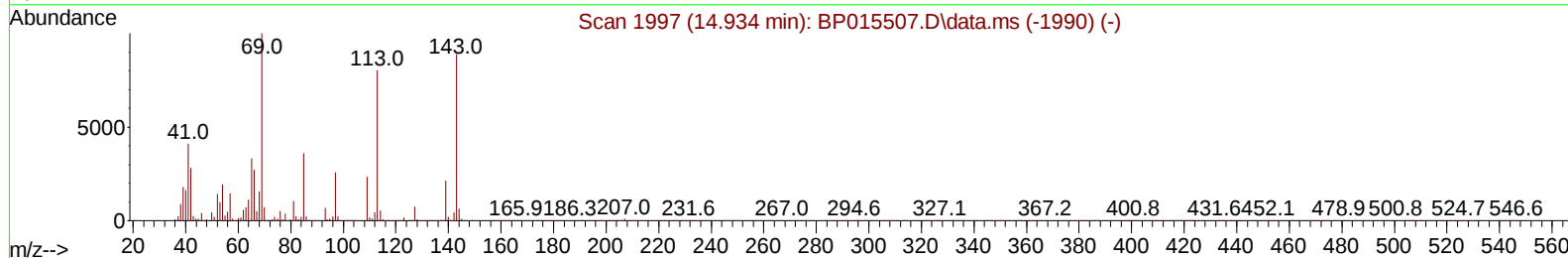
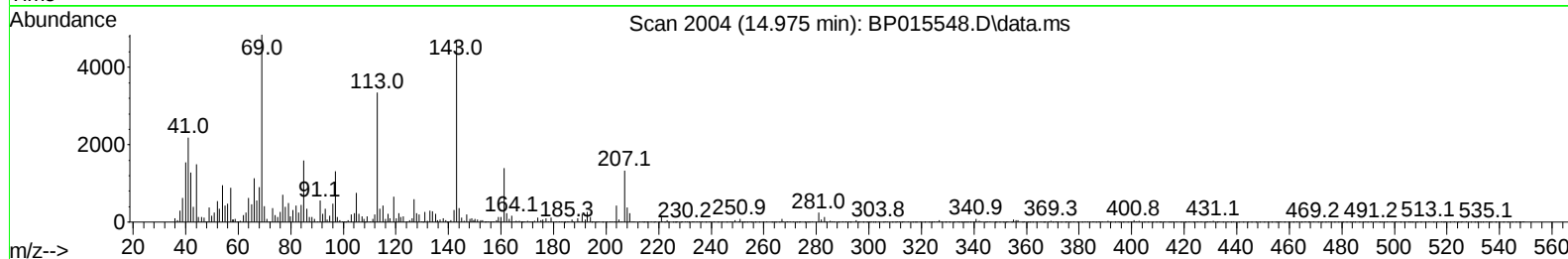
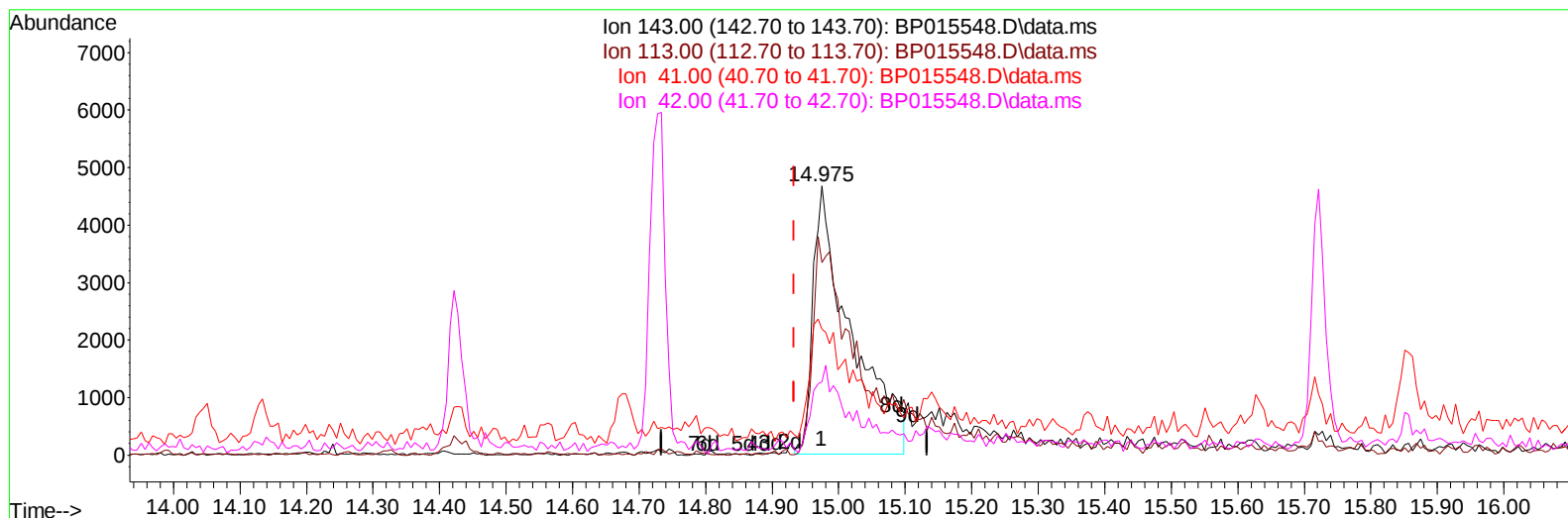
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(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.041) 4.86 ng/ul m

response 18448

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	71.48
41.00	42.20	46.86
42.00	29.40	27.26

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	82502	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	393367	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.727	164	269033	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	625294	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	594647	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	622141	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	4097	1.839	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.246	99	85202	11.210	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	55298	12.182	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	70555	12.803	ng/ul	0.00
15) 4-Methylphenol-d8	8.804	113	66749	10.700	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	35170	11.388	ng/ul	0.00
24) 2-Nitrophenol-d4	9.992	143	41221	11.688	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.539	165	70833	11.010	ng/ul	0.00
31) 4-Chloroaniline-d4	11.069	131	57878	6.268	ng/ul	0.01
46) Dimethylphthalate-d6	14.133	166	267528	12.610	ng/ul	0.00
49) Acenaphthylene-d8	14.422	160	289429	12.477	ng/ul	0.00
54) 4-Nitrophenol-d4	14.975	143	18448m	4.856	ng/ul	0.04
60) Fluorene-d10	15.722	176	232351	12.988	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	26436	6.674	ng/ul	0.00
73) Anthracene-d10	17.586	188	350177	12.309	ng/ul	0.00
81) Pyrene-d10	19.810	212	361574	11.361	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	228541	7.321	ng/ul	0.00

Target Compounds Qvalue

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

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