

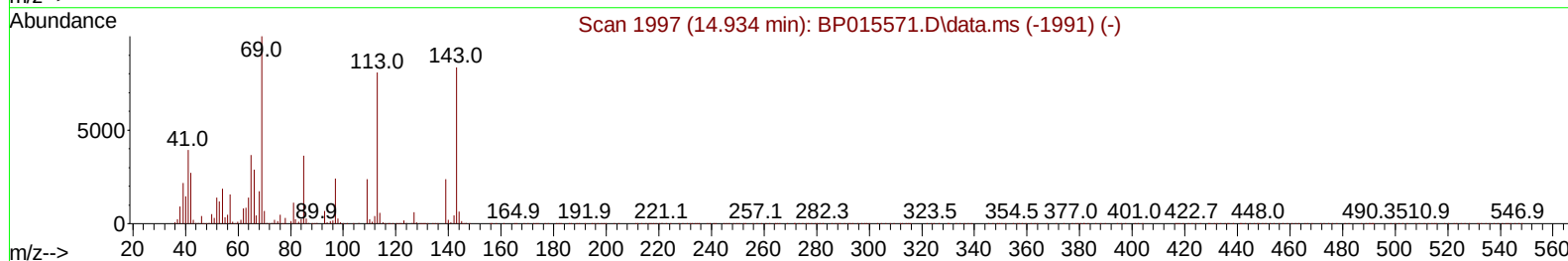
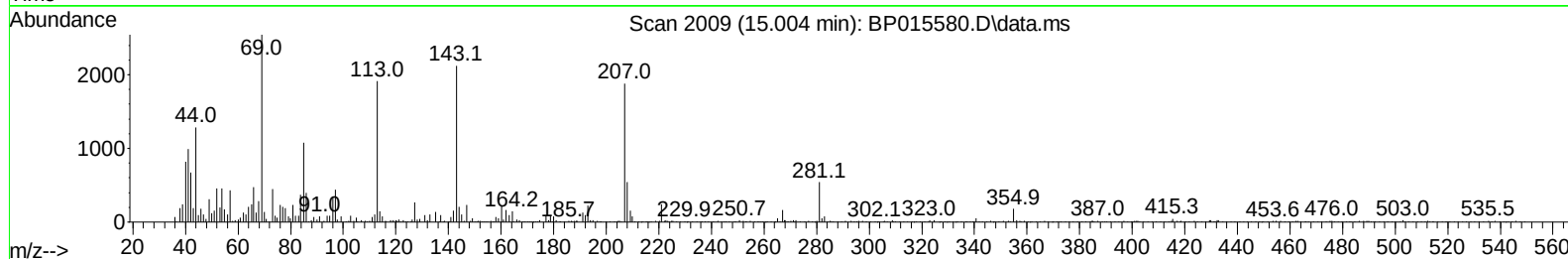
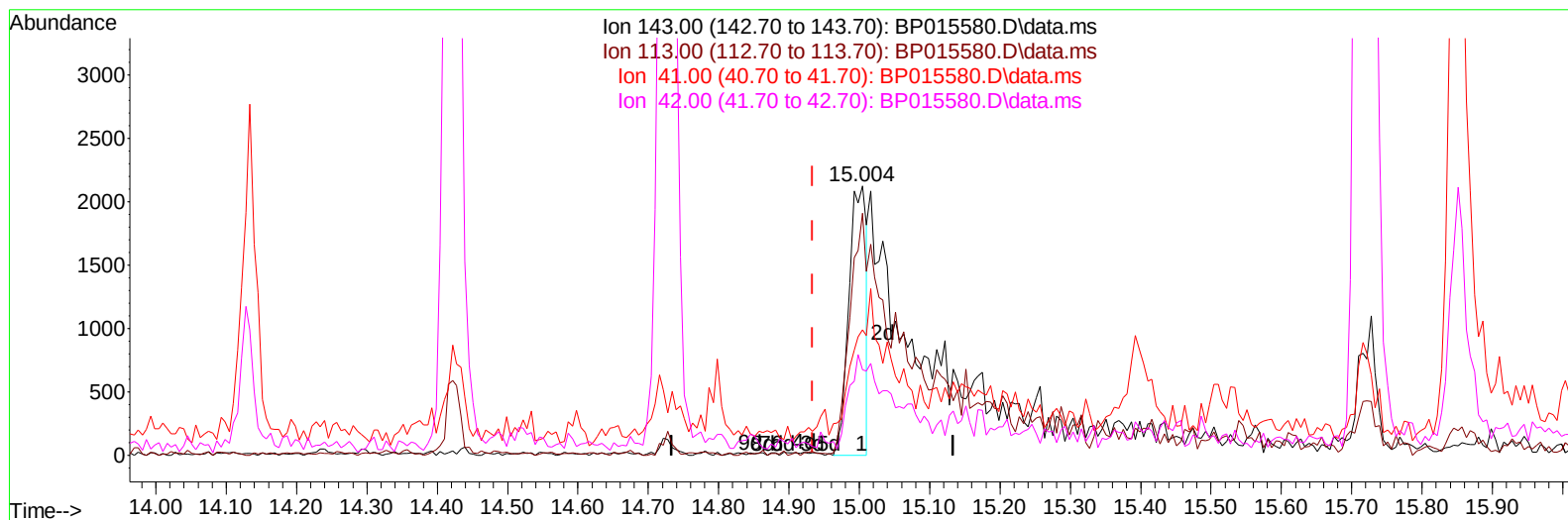
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
 Data File : BP015580.D
 Acq On : 16 Jun 2023 14:40
 Operator : MA/JU
 Sample : 03104-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONQ7

Manual IntegrationsAPPROVED

Quant Time: Jun 16 22:06:41 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 06/19/2023
 Supervised By :mohammad ahmed 06/20/2023



TIC: BP015580.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.004min (+ 0.070) 0.68 ng/ul

response 3696

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	90.00
41.00	42.20	46.63
42.00	29.40	31.68

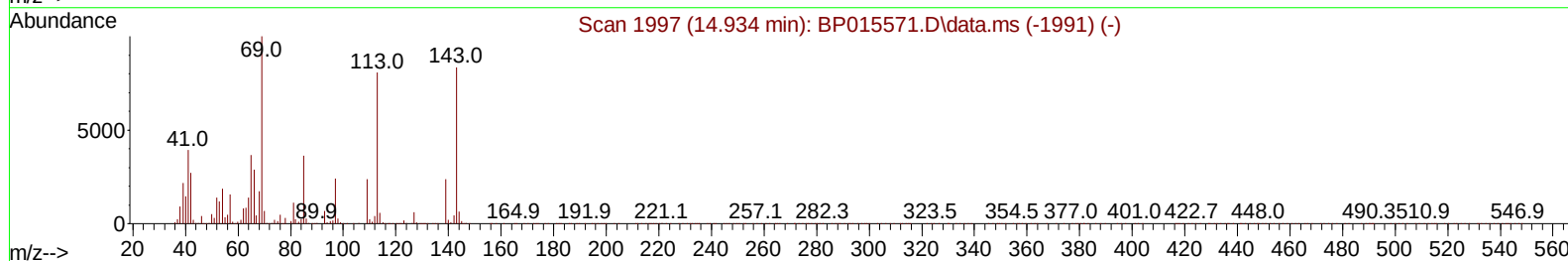
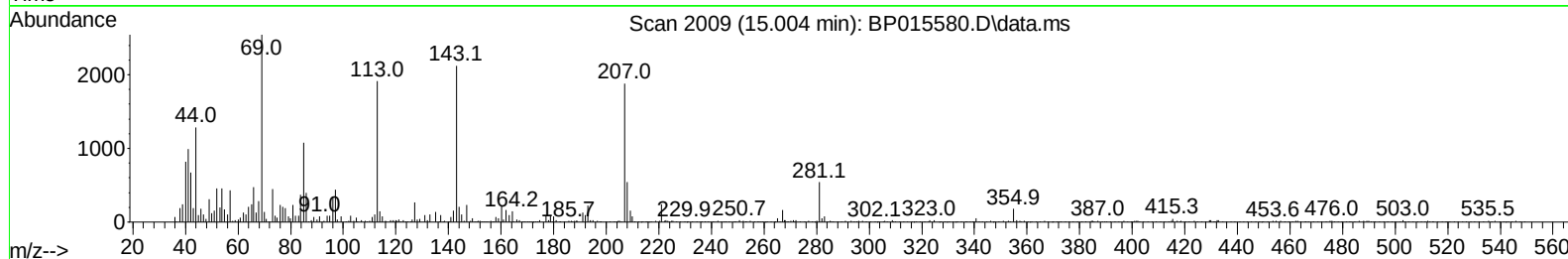
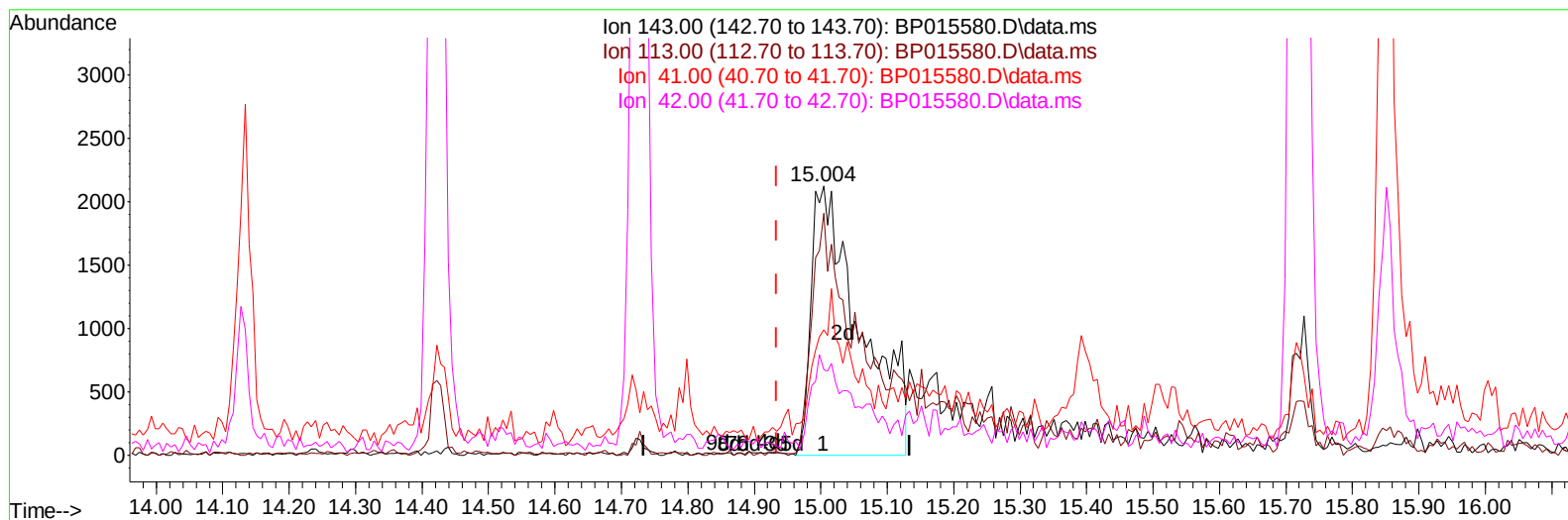
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(54) 4-Nitrophenol-d4 (S)

15.004min (+ 0.070) 2.00 ng/ul m

response 10843

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	90.00
41.00	42.20	46.63
42.00	29.40	31.68

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	108881	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	507468	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.728	164	383455	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	919965	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	925800	20.000	ng/ul	0.00
88) Perylene-d12	24.110	264	1011593	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	11948	4.063	ng/uL	0.00
4) Pyridine-d5	3.864	84	30880	4.043	ng/ul	0.02
7) Phenol-d5	7.246	99	42103	4.197	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.405	67	162916	27.195	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	142415	19.582	ng/ul	0.00
15) 4-Methylphenol-d8	8.804	113	96329	11.701	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	107181	26.902	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	111858	24.585	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	180598	21.760	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	215587	18.096	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	902963	29.862	ng/ul	0.00
49) Acenaphthylene-d8	14.422	160	936110	28.312	ng/ul	0.00
54) 4-Nitrophenol-d4	15.004	143	10843m	2.002	ng/ul	0.07
60) Fluorene-d10	15.722	176	783644	30.733	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	118890	20.400	ng/ul	0.00
73) Anthracene-d10	17.586	188	1411986	33.733	ng/ul	0.00
81) Pyrene-d10	19.810	212	1761531	35.550	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.945	264	1721618	33.918	ng/ul	0.00

Target Compounds Qvalue

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

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