

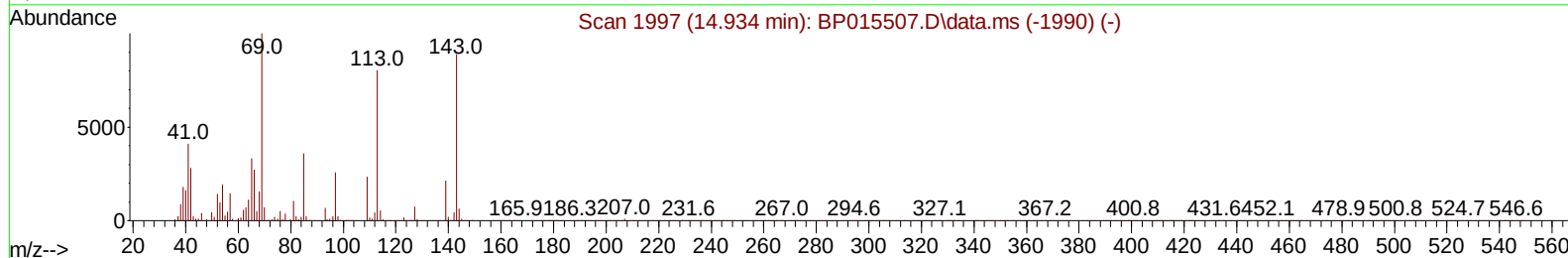
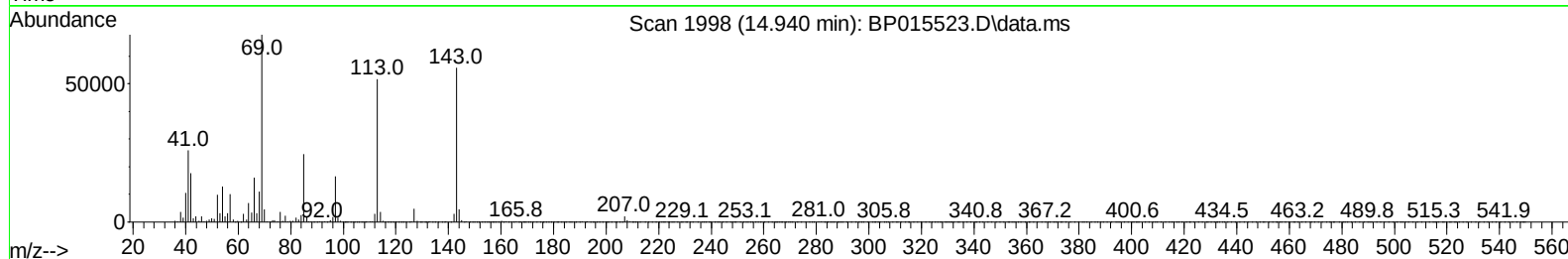
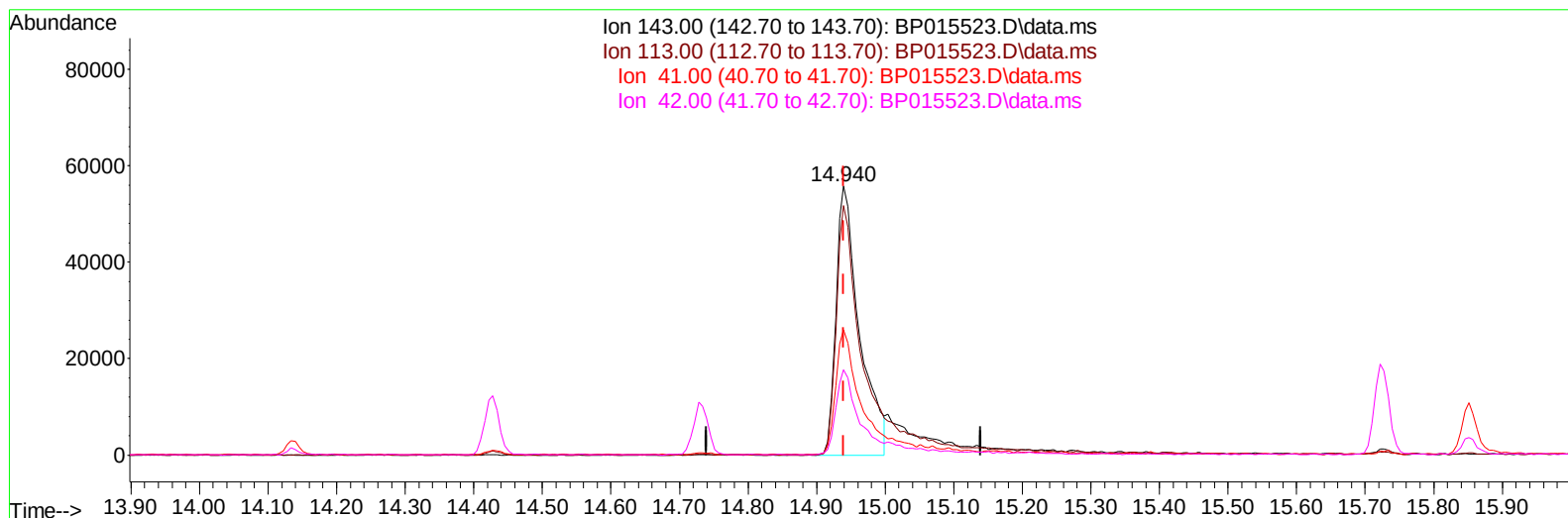
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015523.D
 Acq On : 14 Jun 2023 18:18
 Operator : MA/JU
 Sample : PB153379BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK379

Manual IntegrationsAPPROVED

Quant Time: Jun 14 23:50:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 09 23:33:38 2023
 Response via : Initial Calibration

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



TIC: BP015523.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.940min (-0.000) 21.46 ng/ul

response 131955

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	92.81
41.00	42.20	46.53
42.00	29.40	31.72

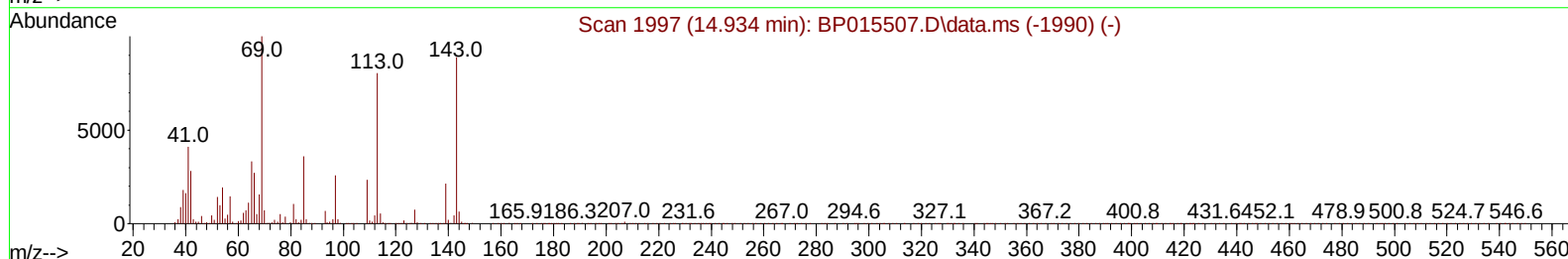
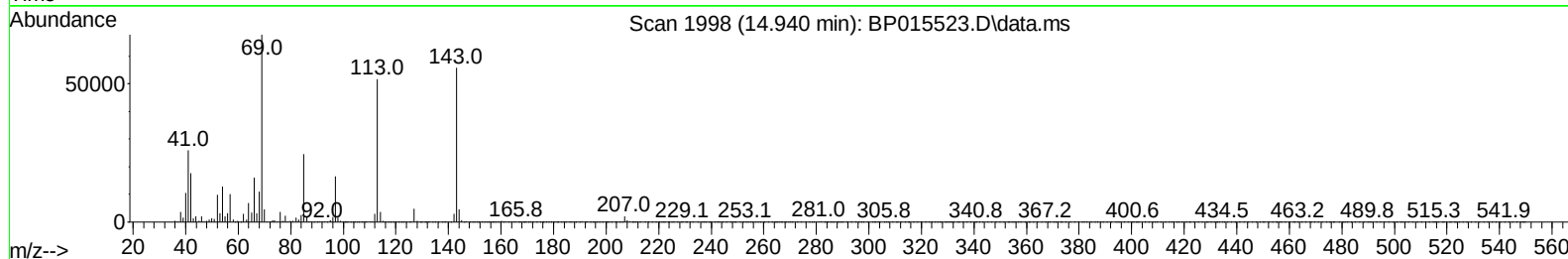
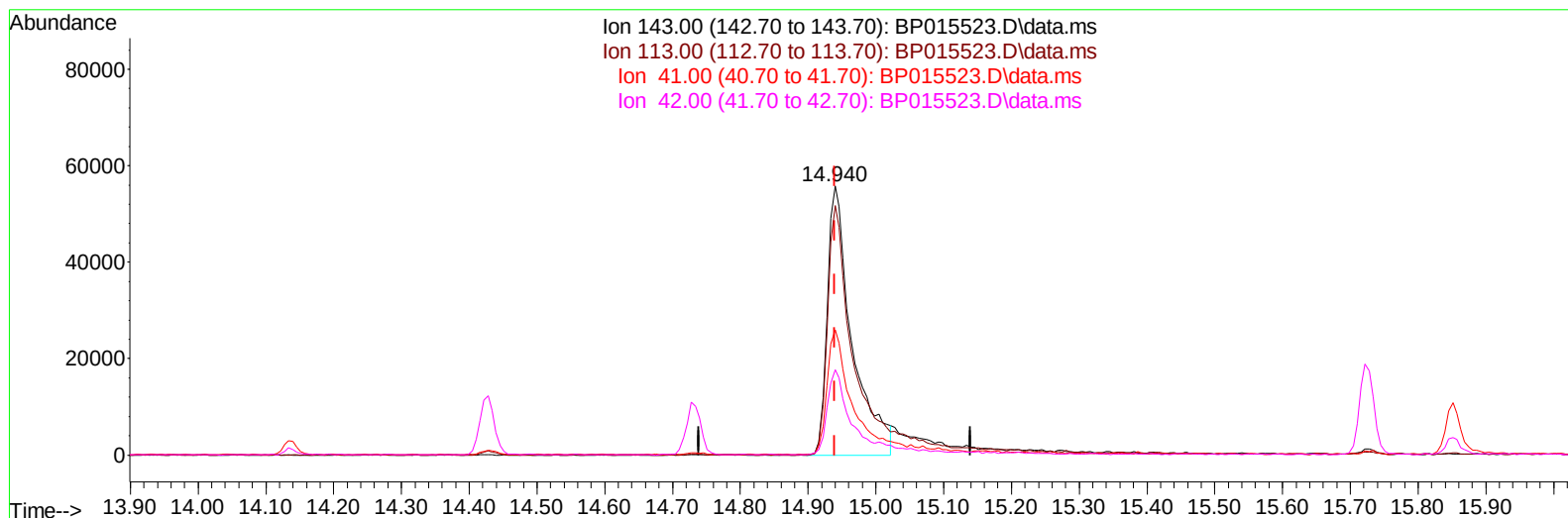
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TIC: BP015523.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.940min (-0.000) 23.03 ng/ul m

response 141646

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	92.81
41.00	42.20	46.53
42.00	29.40	31.72

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
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Quant Time: Jun 14 23:51:14 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	128249	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	607920	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	435506	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	1047416	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	1066793	20.000	ng/ul	0.00
88) Perylene-d12	24.121	264	1134069	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	23663	6.832	ng/uL	0.00
4) Pyridine-d5	3.852	84	130035	14.455	ng/ul	0.00
7) Phenol-d5	7.240	99	358067	30.306	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.411	67	230010	32.597	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	285563	33.336	ng/ul	0.00
15) 4-Methylphenol-d8	8.799	113	310039	31.973	ng/ul	-0.01
21) Nitrobenzene-d5	9.269	128	152164	31.882	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	177654	32.594	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	306343	30.811	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	218204	15.290	ng/ul	0.00
46) Dimethylphthalate-d6	14.140	166	1136877	33.104	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1243472	33.113	ng/ul	0.00
54) 4-Nitrophenol-d4	14.940	143	141646m	23.032	ng/ul	0.00
60) Fluorene-d10	15.728	176	967319	33.402	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	184179	27.757	ng/ul	0.00
73) Anthracene-d10	17.586	188	1613794	33.863	ng/ul	0.00
81) Pyrene-d10	19.810	212	2005983	35.133	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	1989390	34.961	ng/ul	0.00

Target Compounds Qvalue

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

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