

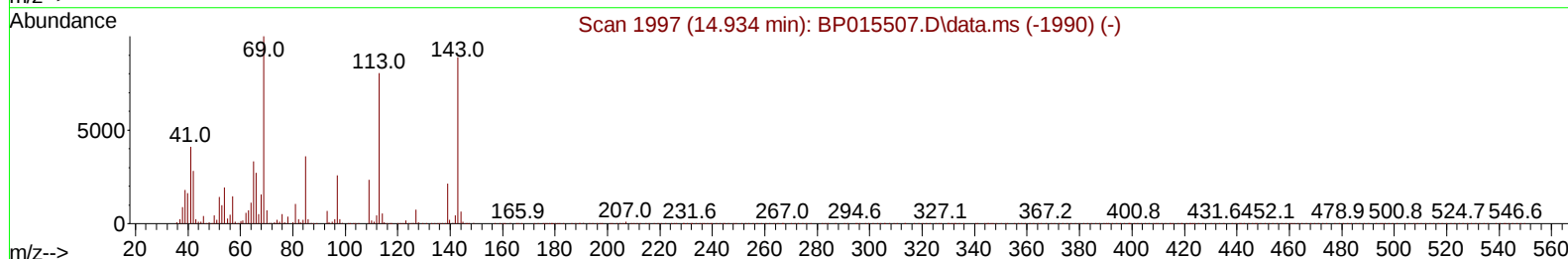
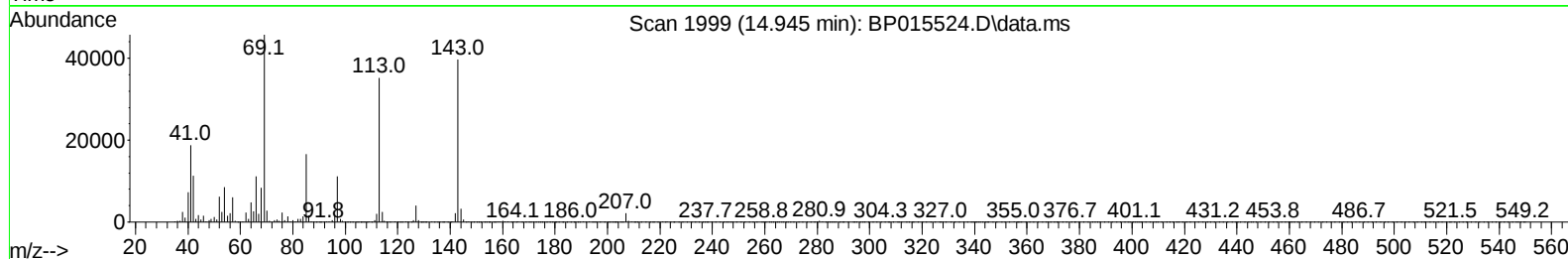
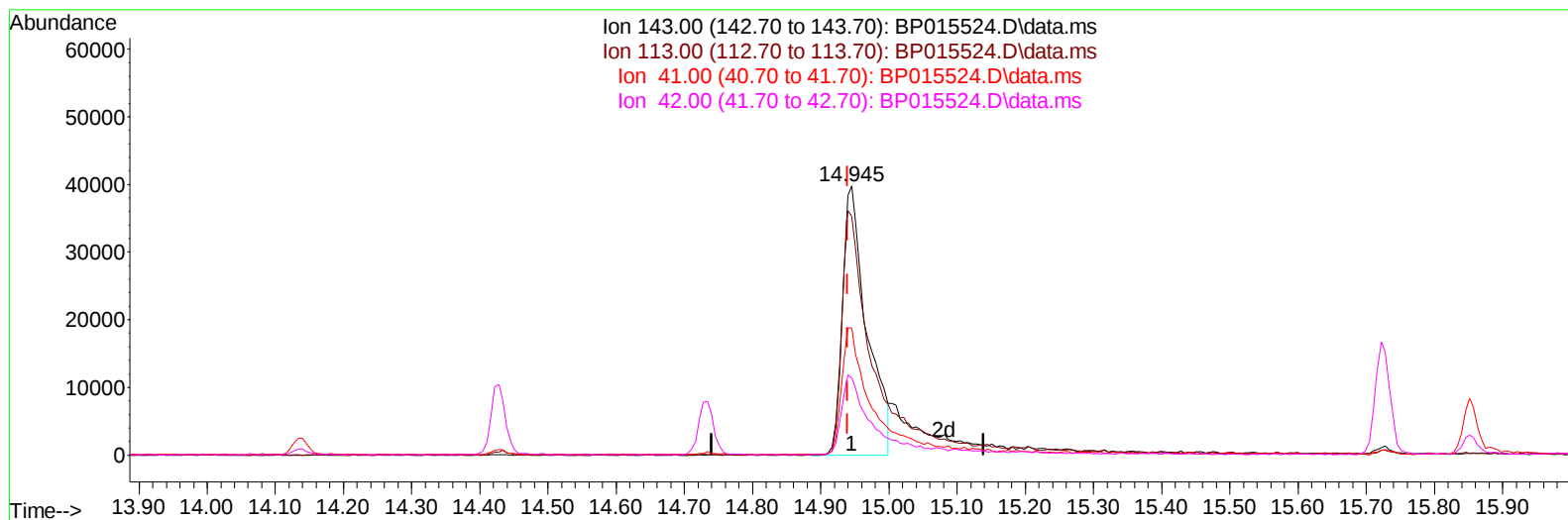
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015524.D
 Acq On : 14 Jun 2023 18:52
 Operator : MA/JU
 Sample : PB153412BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK412

Manual IntegrationsAPPROVED

Quant Time: Jun 14 23:53:14 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: BP015524.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.945min (+ 0.006) 21.46 ng/u1

response 99274

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	88.75
41.00	42.20	47.28
42.00	29.40	28.62

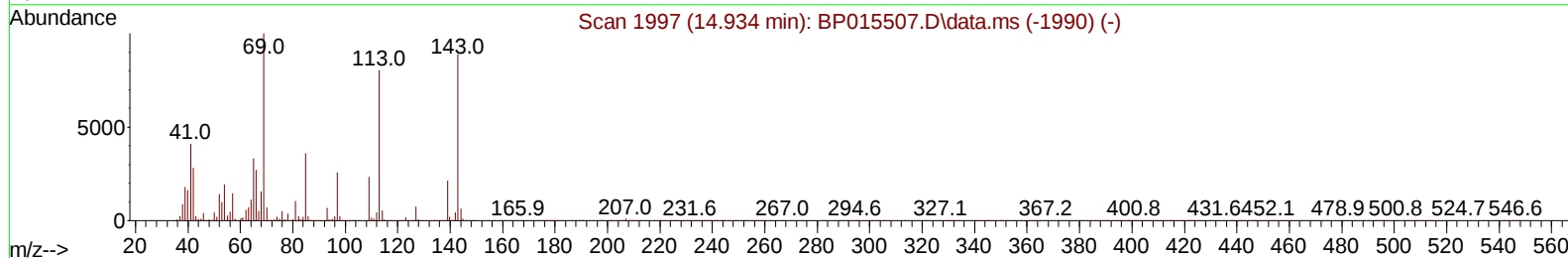
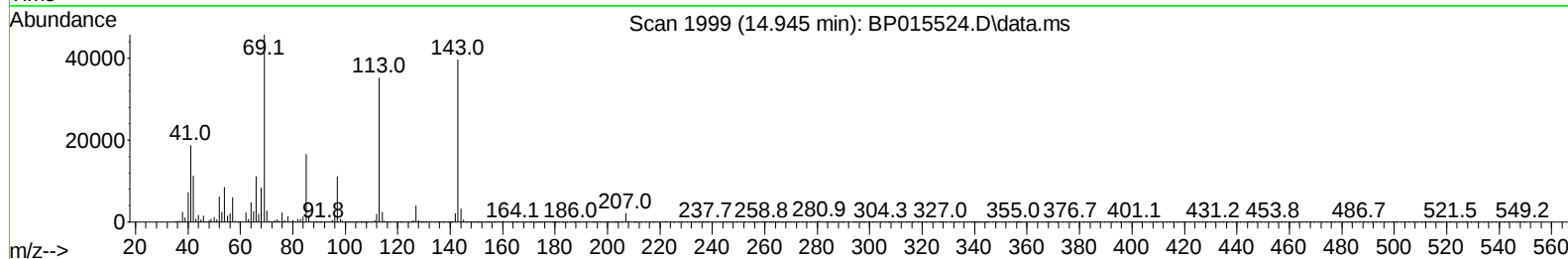
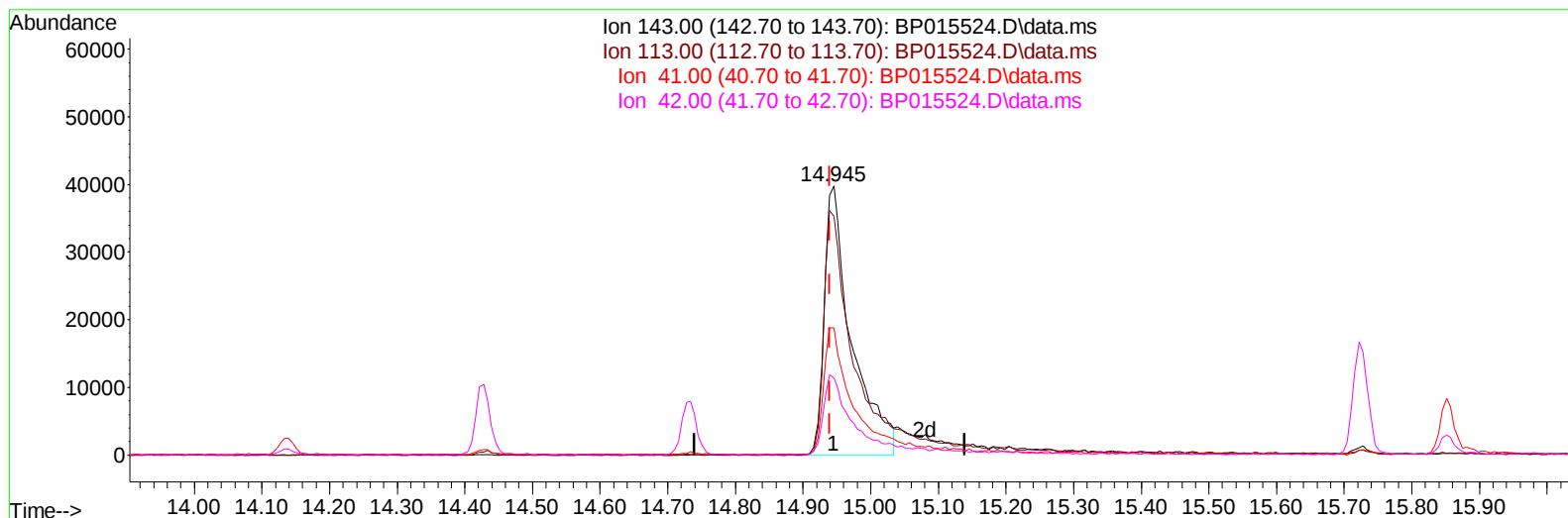
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TIC: BP015524.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.945min (+ 0.006) 24.05 ng/ul m

response 111215

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	88.75
41.00	42.20	47.28
42.00	29.40	28.62

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	95925	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	457241	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	327522	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	778231	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	769530	20.000	ng/ul	0.00
88) Perylene-d12	24.121	264	822252	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	16538	6.384	ng/uL	0.00
4) Pyridine-d5	3.846	84	236351	35.126	ng/ul	0.00
7) Phenol-d5	7.246	99	294082	33.278	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	193183	36.603	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	234081	36.534	ng/ul	0.00
15) 4-Methylphenol-d8	8.805	113	259487	35.777	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	124747	34.750	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.993	143	145885	35.586	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	253263	33.867	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	367078	34.197	ng/ul	0.00
46) Dimethylphthalate-d6	14.140	166	989154	38.299	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1038250	36.764	ng/ul	0.00
54) 4-Nitrophenol-d4	14.945	143	111215m	24.046	ng/ul	0.00
60) Fluorene-d10	15.728	176	806858	37.047	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	147620	29.942	ng/ul	0.00
73) Anthracene-d10	17.586	188	1371518	38.734	ng/ul	0.00
81) Pyrene-d10	19.810	212	1689378	41.018	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	1604499	38.890	ng/ul	0.00

Target Compounds Qvalue

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

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