

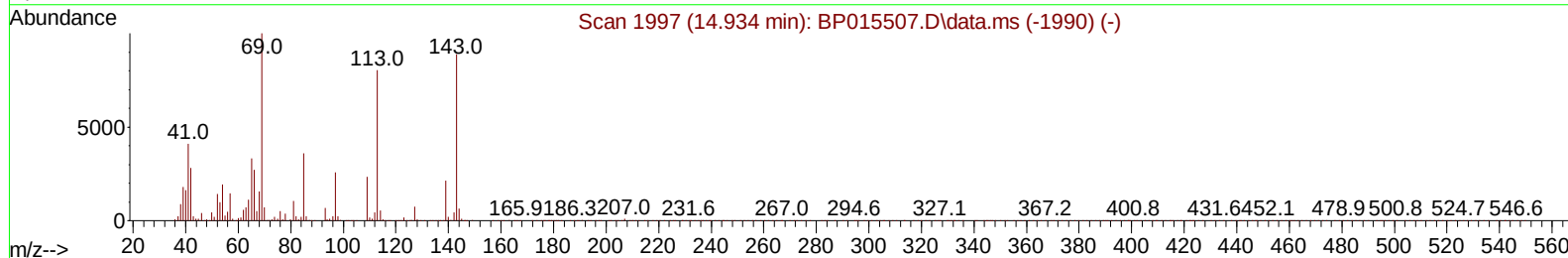
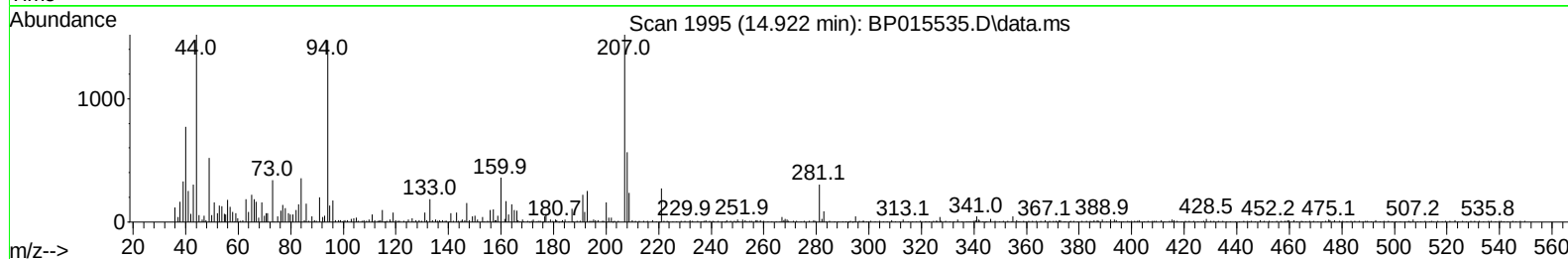
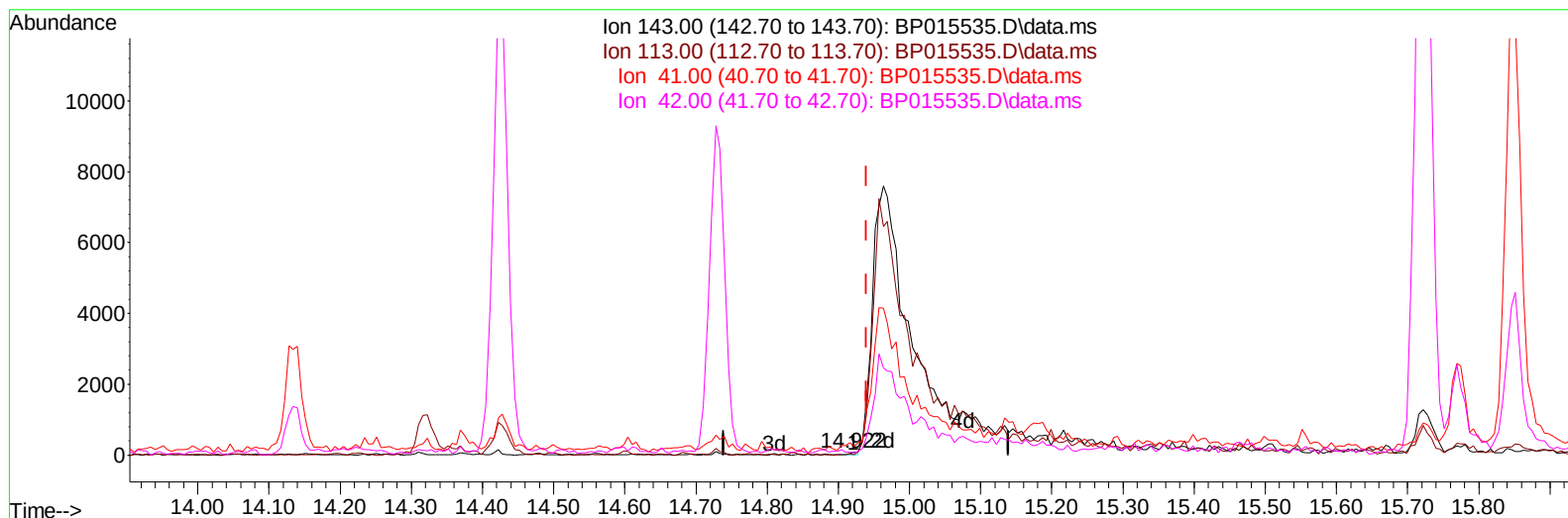
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015535.D
 Acq On : 15 Jun 2023 01:03
 Operator : MA/JU
 Sample : 03104-09
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONLO

Manual IntegrationsAPPROVED

Quant Time: Jun 15 02:04:59 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: BP015535.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (-0.018) 0.01 ng/u1

response 58

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	21.52#
41.00	42.20	316.46#
42.00	29.40	83.54#

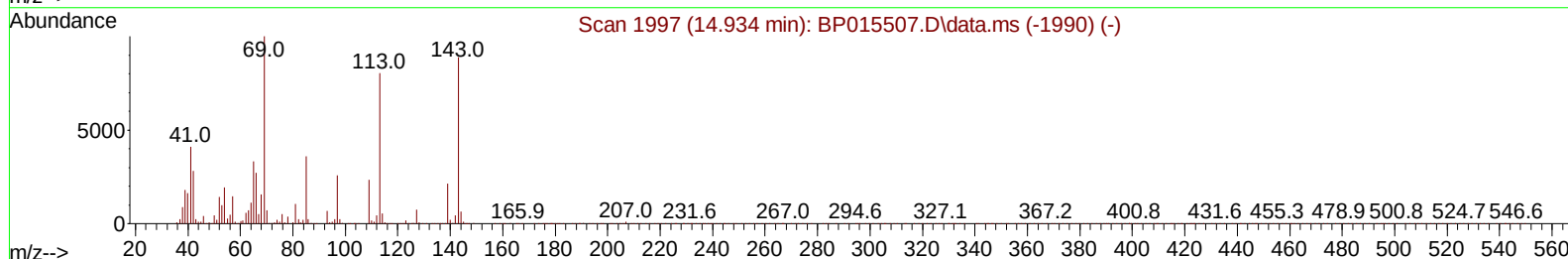
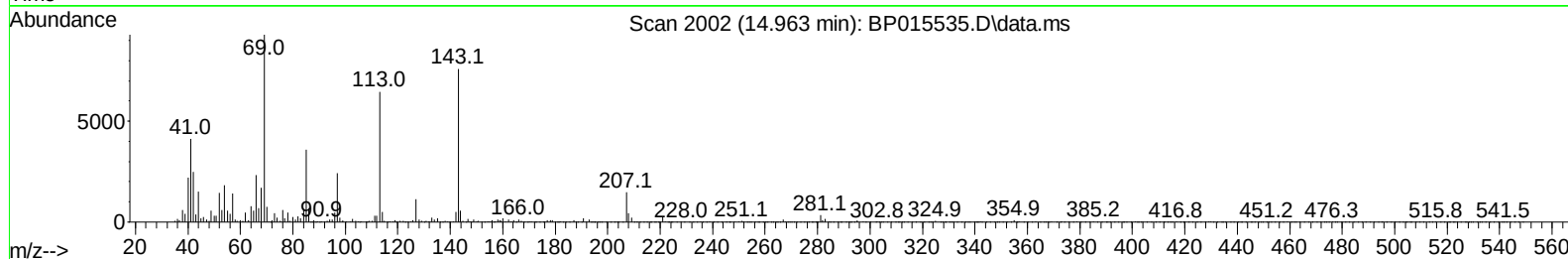
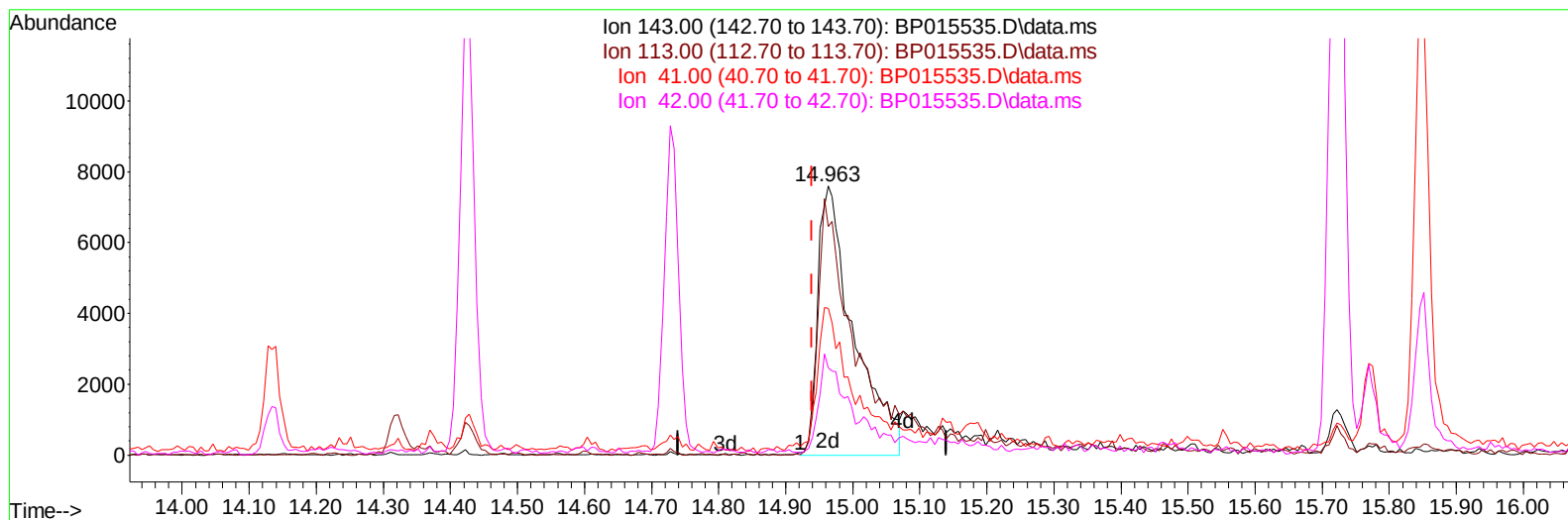
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TIC: BP015535.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.963min (+ 0.024) 5.38 ng/ul m

response 28066

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	85.06
41.00	42.20	54.50#
42.00	29.40	32.55

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
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Quant Time: Jun 15 02:25:05 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	112897	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	531835	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.728	164	369529	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.487	188	891602	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	916082	20.000	ng/ul	0.00
88) Perylene-d12	24.116	264	986097	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	15868	5.205	ng/uL	0.00
4) Pyridine-d5	3.870	84	21007	2.653	ng/ul	0.02
7) Phenol-d5	7.246	99	81995	7.884	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.411	67	242373	39.020	ng/ul	0.00
11) 2-Chlorophenol-d4	7.611	132	234669	31.120	ng/ul	-0.01
15) 4-Methylphenol-d8	8.799	113	146251	17.133	ng/ul	-0.01
21) Nitrobenzene-d5	9.263	128	159833	38.279	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.993	143	186733	39.161	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	294276	33.832	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	159652	12.787	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	1248164	42.834	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1322558	41.508	ng/ul	0.00
54) 4-Nitrophenol-d4	14.963	143	28066m	5.378	ng/ul	0.02
60) Fluorene-d10	15.722	176	1061857	43.213	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	223740	39.611	ng/ul	0.00
73) Anthracene-d10	17.587	188	1780525	43.891	ng/ul	0.00
81) Pyrene-d10	19.810	212	2299570	46.901	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2240205	45.277	ng/ul	0.00

Target Compounds Qvalue

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

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