

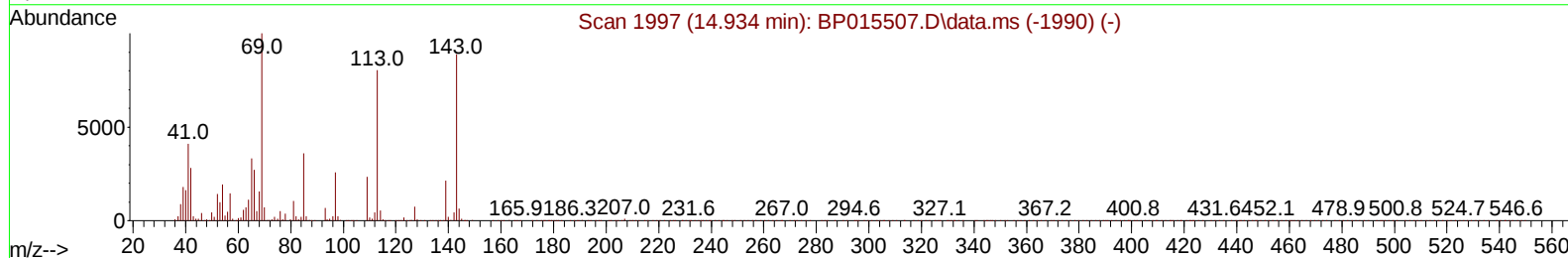
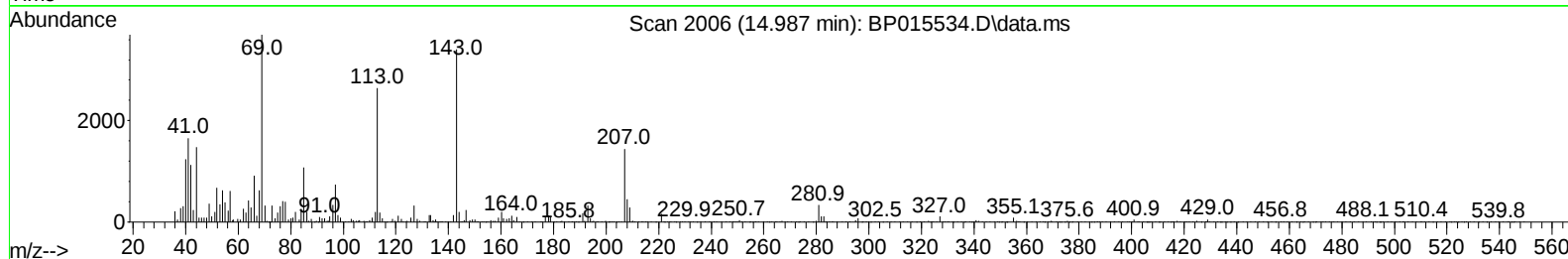
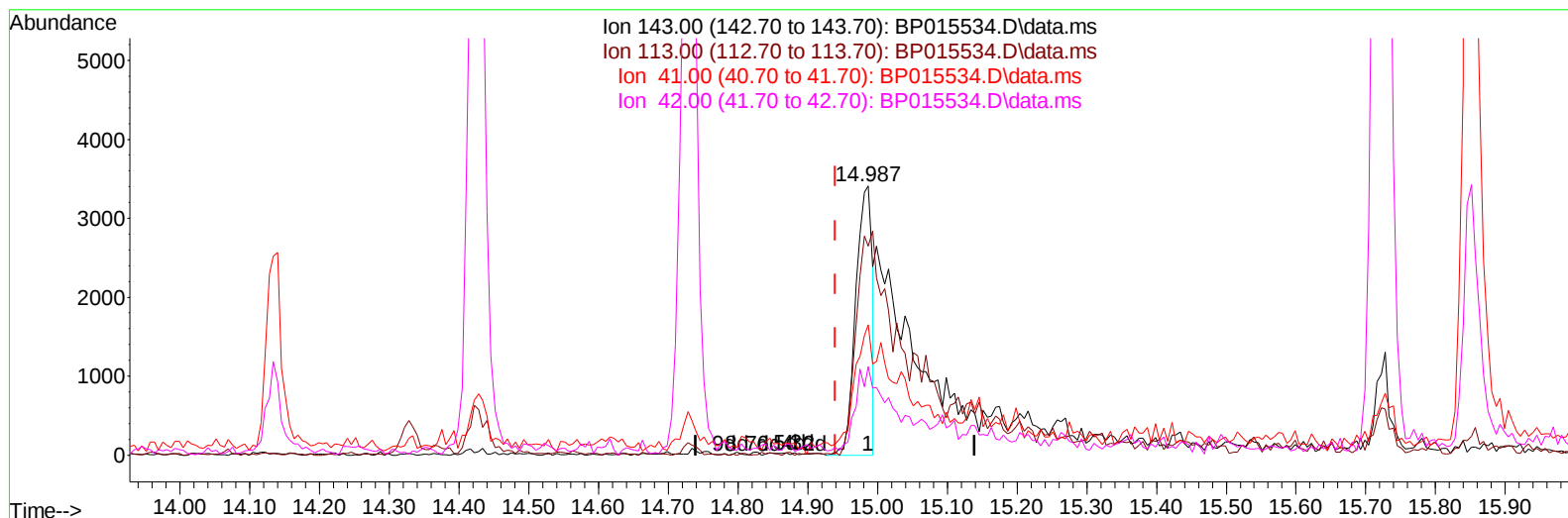
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015534.D
 Acq On : 15 Jun 2023 00:29
 Operator : MA/JU
 Sample : 03104-08
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONK9

Manual IntegrationsAPPROVED

Quant Time: Jun 15 02:04:46 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: BP015534.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.987min (+ 0.047) 1.05 ng/u1

response 5535

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	77.72
41.00	42.20	48.55
42.00	29.40	32.99

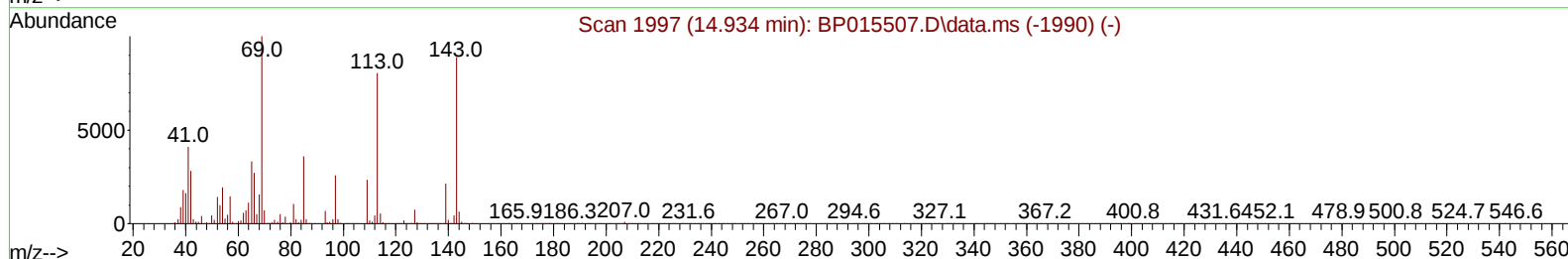
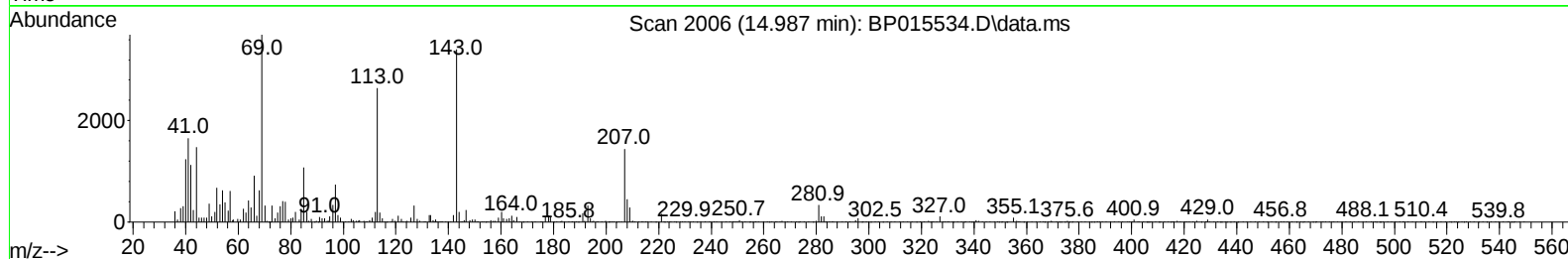
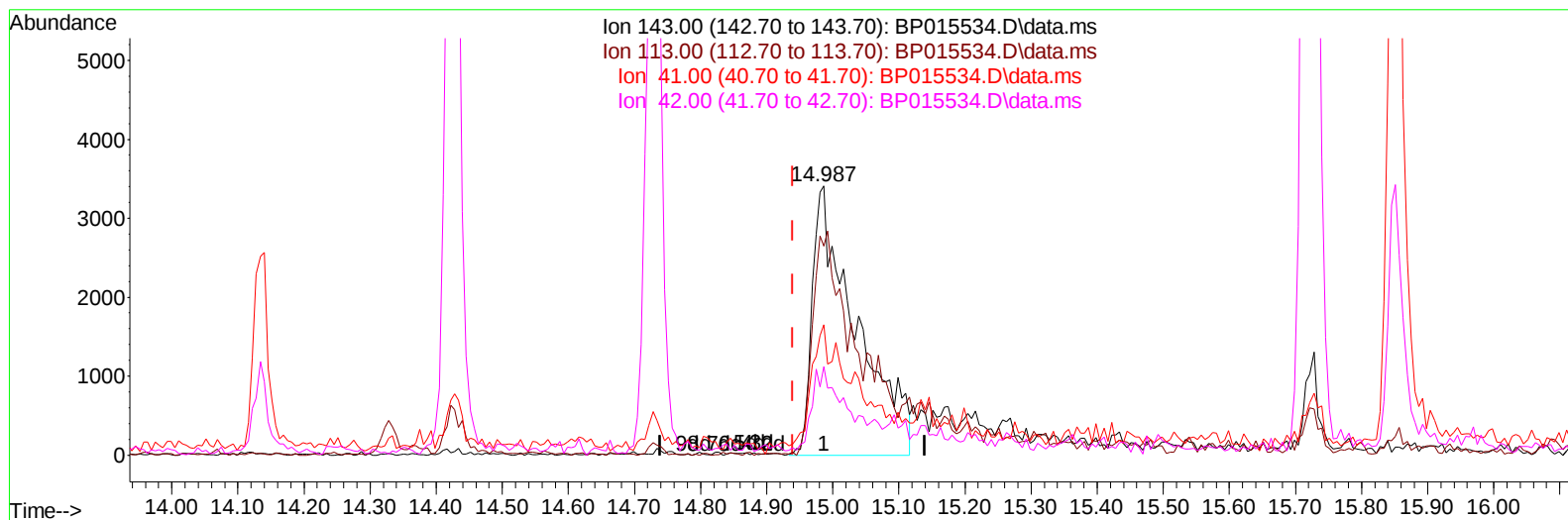
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(54) 4-Nitrophenol-d4 (S)

14.987min (+ 0.047) 2.99 ng/ul m

response 15673

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	77.72
41.00	42.20	48.55
42.00	29.40	32.99

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	111816	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	528125	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.728	164	371691	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.487	188	884463	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	910158	20.000	ng/ul	0.00
88) Perylene-d12	24.116	264	993958	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	12783	4.233	ng/uL	0.00
4) Pyridine-d5	3.858	84	47523	6.059	ng/ul	0.01
7) Phenol-d5	7.246	99	53304	5.175	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.411	67	193799	31.501	ng/ul	0.00
11) 2-Chlorophenol-d4	7.611	132	175935	23.557	ng/ul	-0.01
15) 4-Methylphenol-d8	8.805	113	117247	13.868	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	128198	30.919	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.993	143	140037	29.575	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	225535	26.111	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	286717	23.126	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	983554	33.557	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1067494	33.308	ng/ul	0.00
54) 4-Nitrophenol-d4	14.987	143	15673m	2.986	ng/ul	0.05
60) Fluorene-d10	15.722	176	876051	35.444	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	177372	31.656	ng/ul	0.00
73) Anthracene-d10	17.587	188	1589491	39.498	ng/ul	0.00
81) Pyrene-d10	19.810	212	2006437	41.189	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2002199	40.146	ng/ul	0.00

Target Compounds Qvalue

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

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