

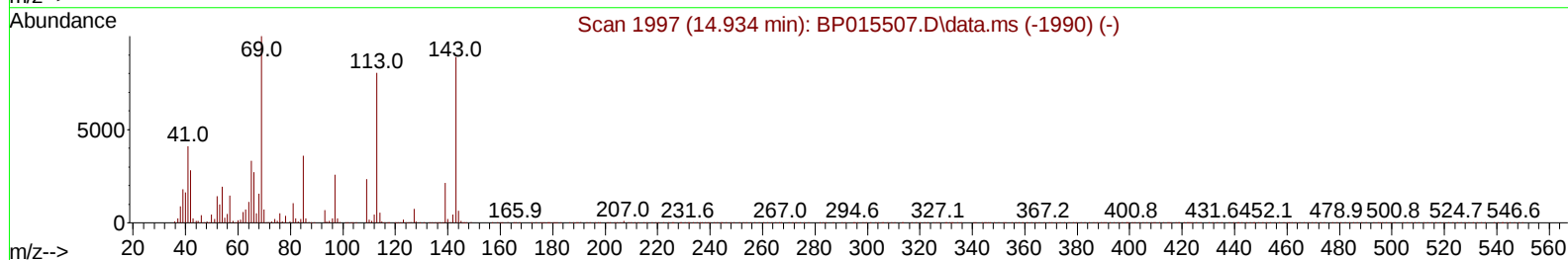
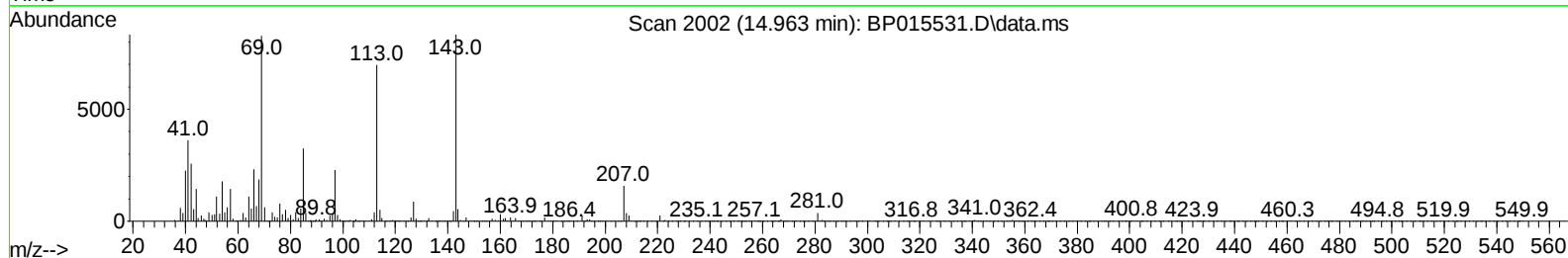
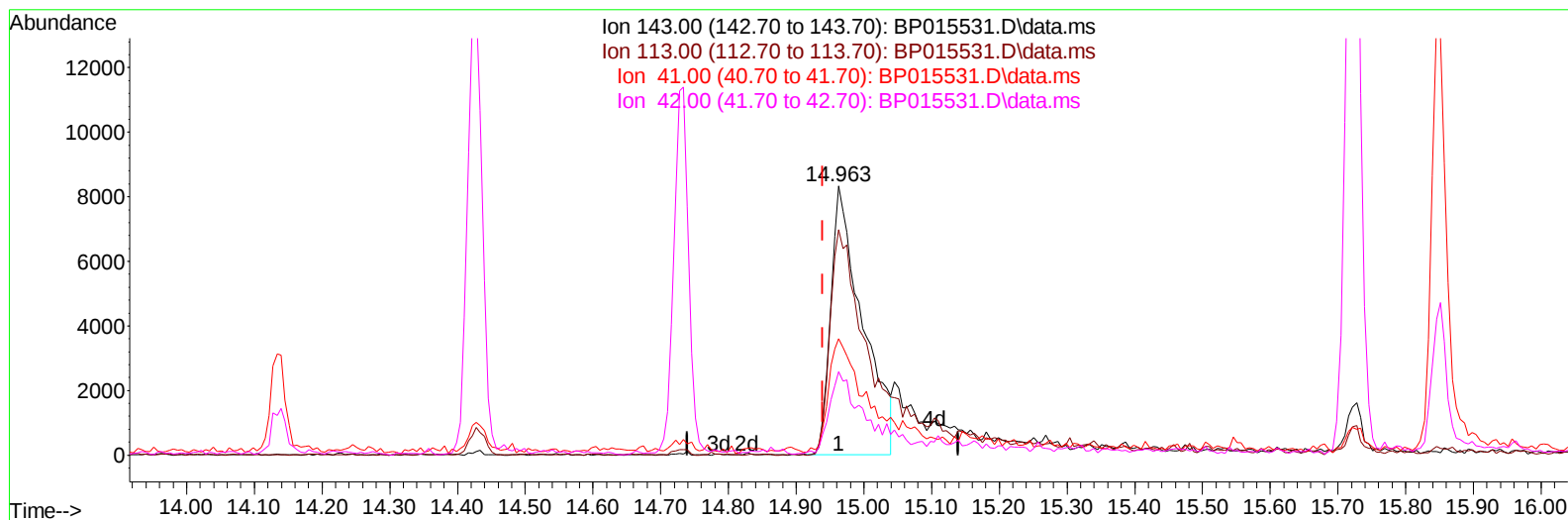
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015531.D
 Acq On : 14 Jun 2023 22:49
 Operator : MA/JU
 Sample : 03104-03
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONL9

Manual IntegrationsAPPROVED

Quant Time: Jun 15 00:05:33 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: BP015531.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.963min (+ 0.023) 4.13 ng/u1

response 26859

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	83.88
41.00	42.20	43.34
42.00	29.40	30.91

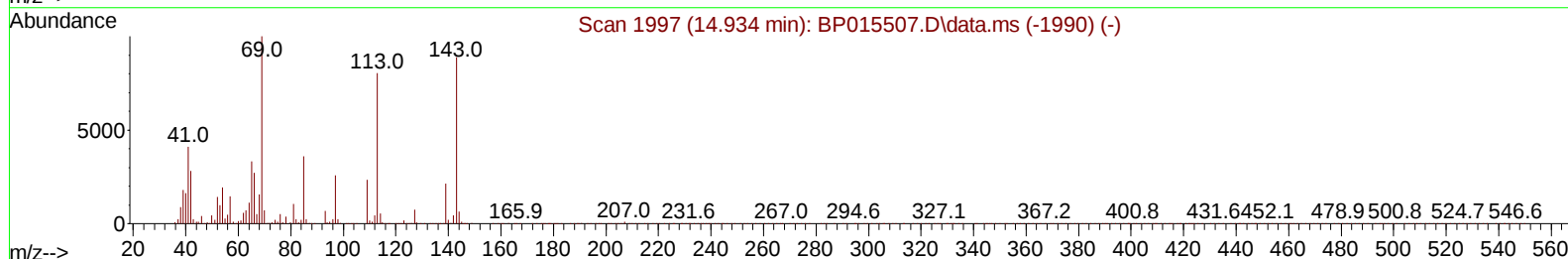
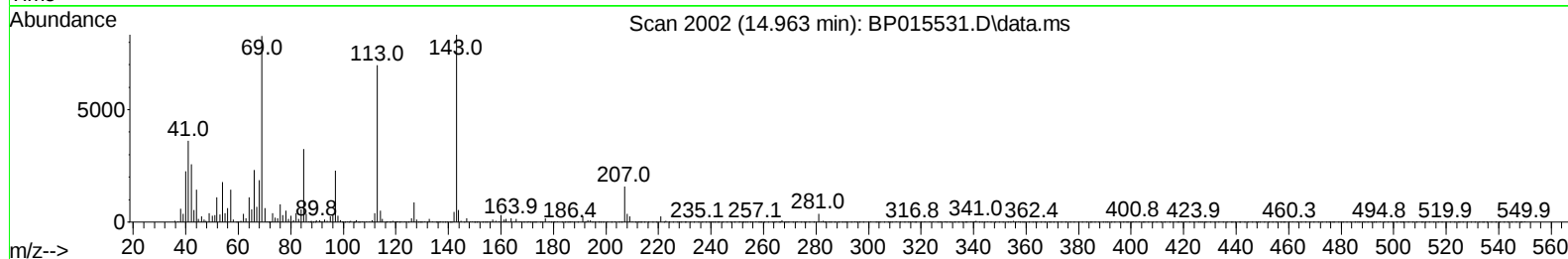
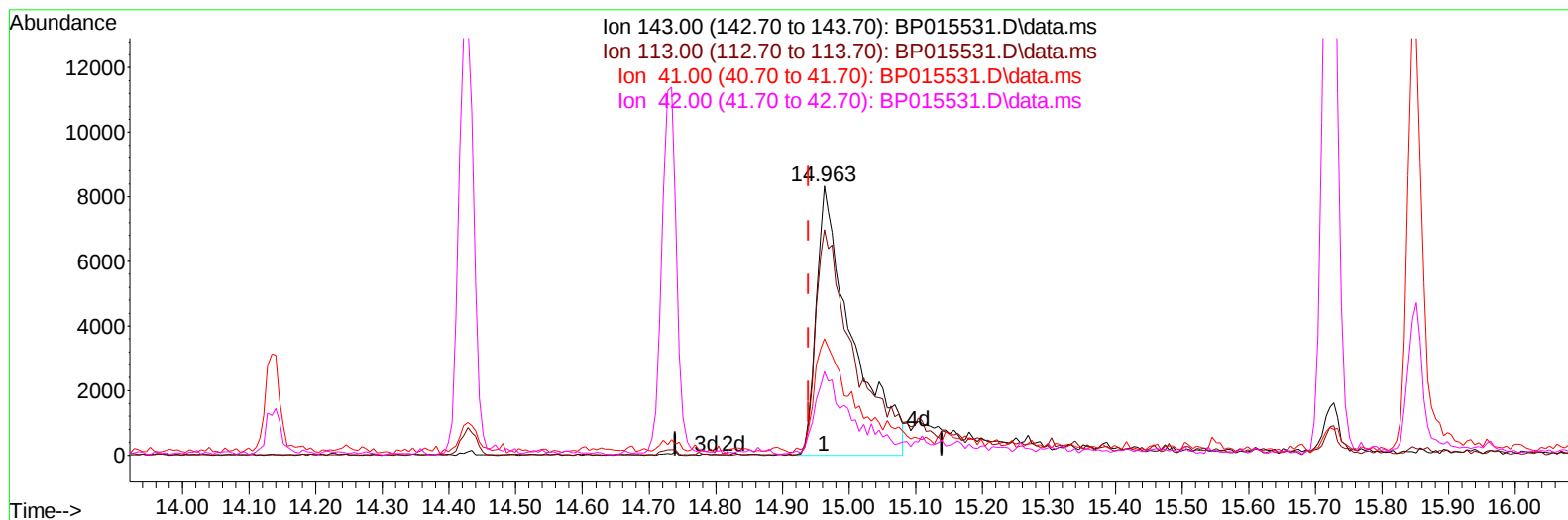
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TIC: BP015531.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.963min (+ 0.023) 4.74 ng/ul m

response 30817

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	83.88
41.00	42.20	43.34
42.00	29.40	30.91

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
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Quant Time: Jun 15 00:06:27 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	139421	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	652323	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	460438	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	1094248	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	1152948	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	1245847	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	15241	4.048	ng/uL	0.00
4) Pyridine-d5	3.864	84	54261	5.548	ng/ul	0.02
7) Phenol-d5	7.246	99	95387	7.426	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	263335	34.329	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	257395	27.640	ng/ul	0.00
15) 4-Methylphenol-d8	8.799	113	191293	18.147	ng/ul	-0.01
21) Nitrobenzene-d5	9.263	128	179893	35.126	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.993	143	198579	33.953	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	324738	30.438	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	352732	23.034	ng/ul	0.00
46) Dimethylphthalate-d6	14.139	166	1318625	36.317	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1412353	35.574	ng/ul	0.00
54) 4-Nitrophenol-d4	14.963	143	30817m	4.740	ng/ul	0.02
60) Fluorene-d10	15.722	176	1137650	37.157	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	237846	34.311	ng/ul	0.00
73) Anthracene-d10	17.586	188	2012533	40.423	ng/ul	0.00
81) Pyrene-d10	19.810	212	2527062	40.952	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2546602	40.738	ng/ul	0.00

Target Compounds Qvalue

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

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