

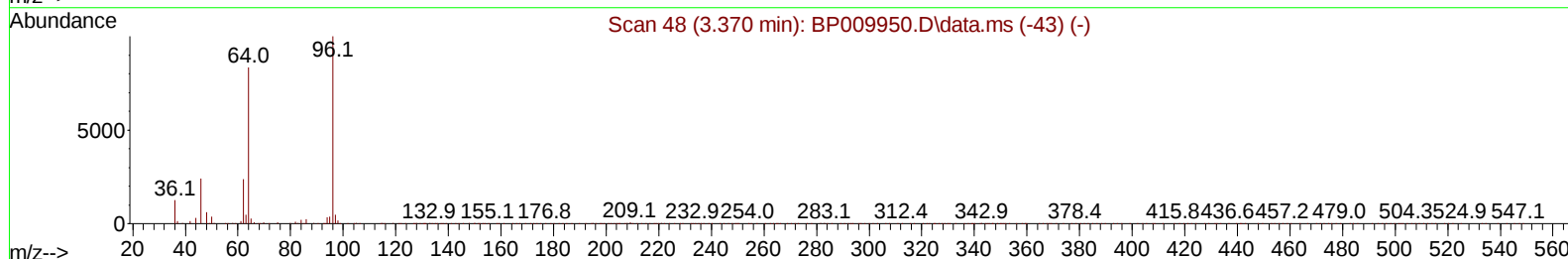
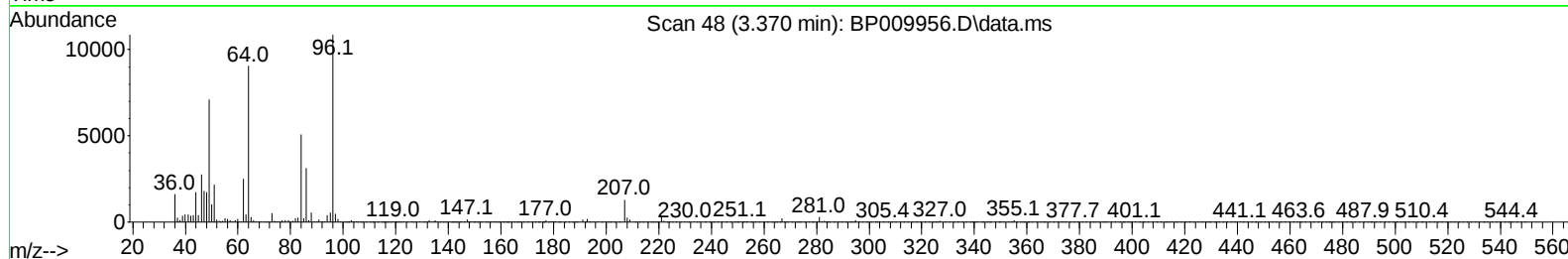
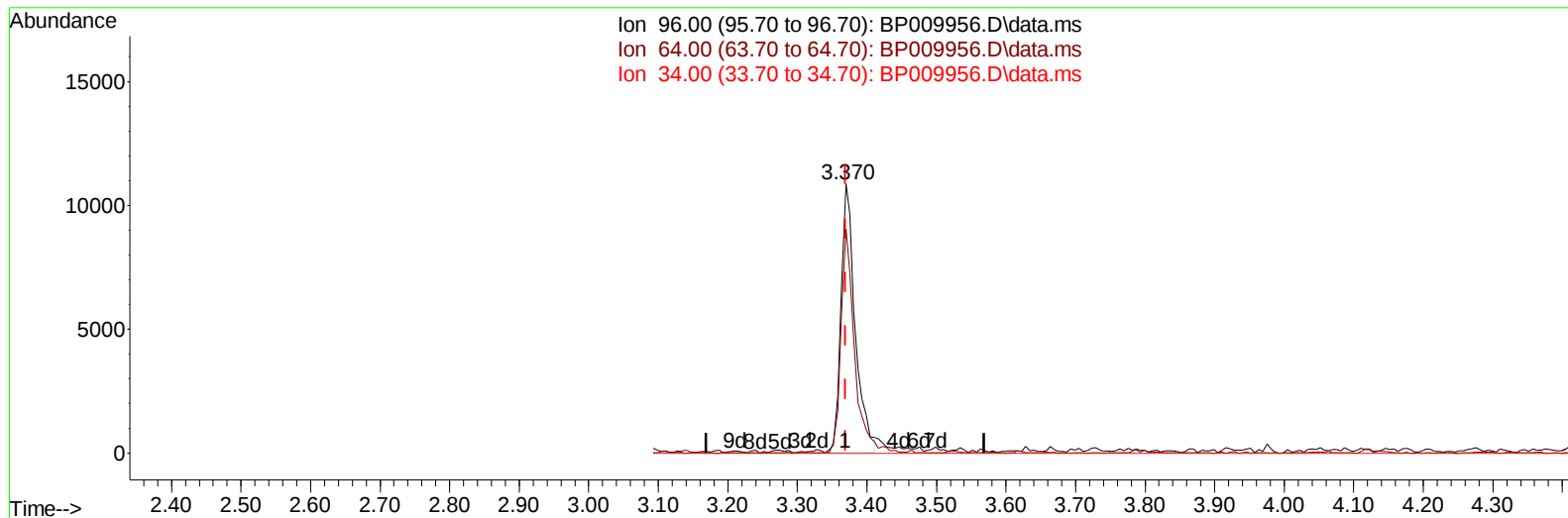
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
Data File : BP009956.D
Acq On : 19 Apr 2022 21:25
Operator : CG/JU
Sample : N2388-10
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF5

Manual IntegrationsAPPROVED

Quant Time: Apr 20 01:33:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 20 01:32:14 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/20/2022
Supervised By :Yogesh Patel 04/25/2022



TIC: BP009956.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 5.79 ng/uL

response 15502

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.32#
34.00	0.00	0.00
0.00	0.00	0.00

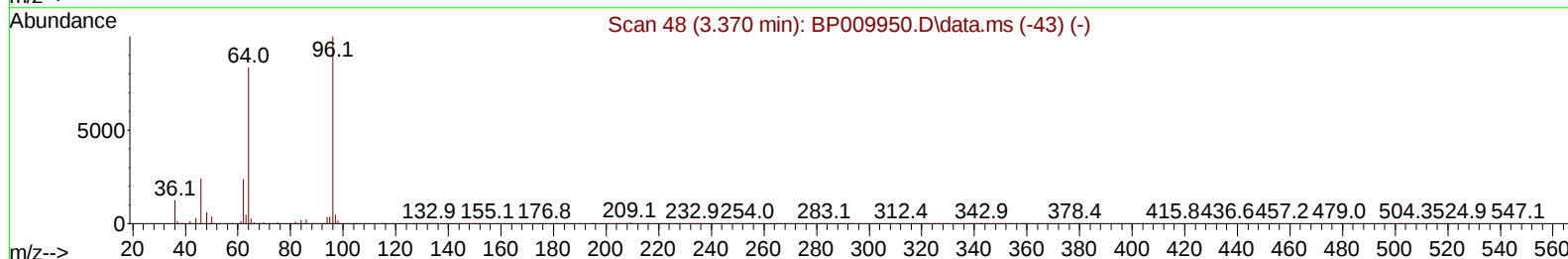
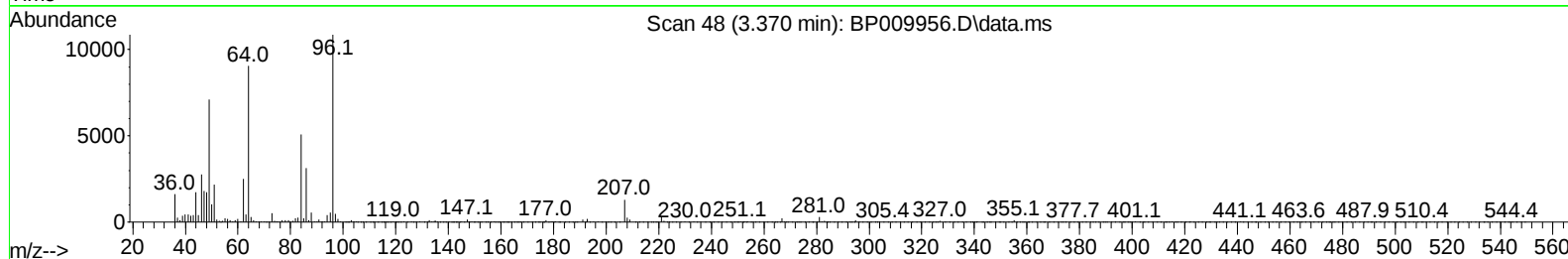
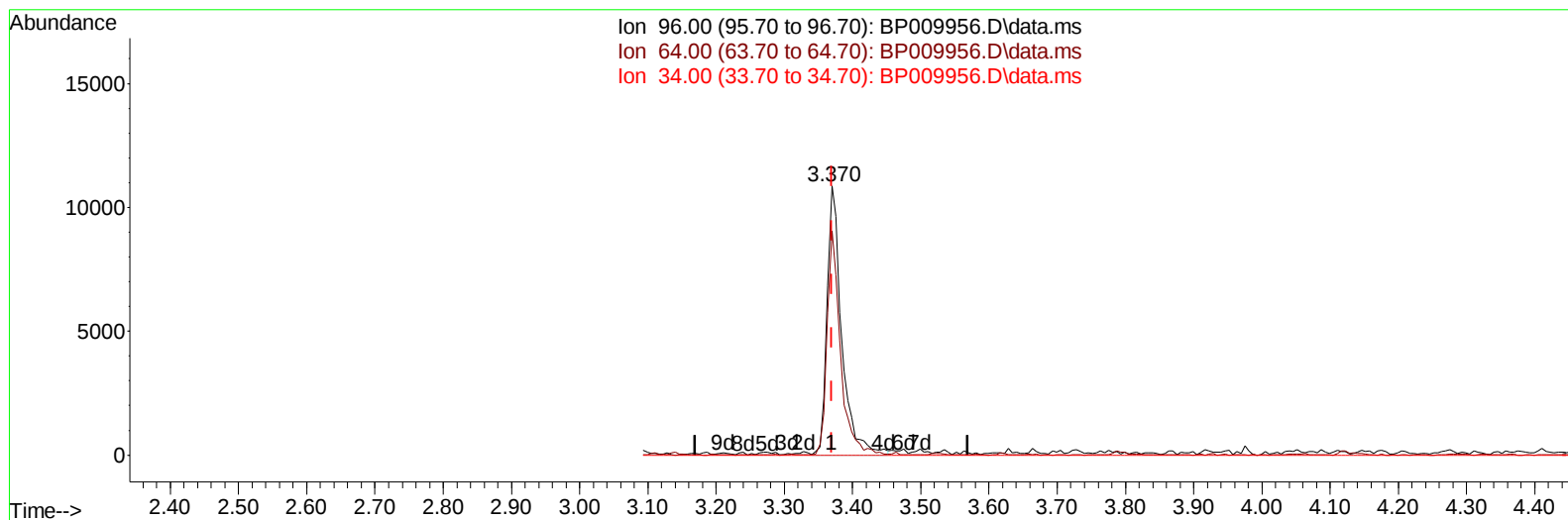
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TIC: BP009956.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 6.10 ng/uL m

response 16327

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.32#
34.00	0.00	0.00
0.00	0.00	0.00

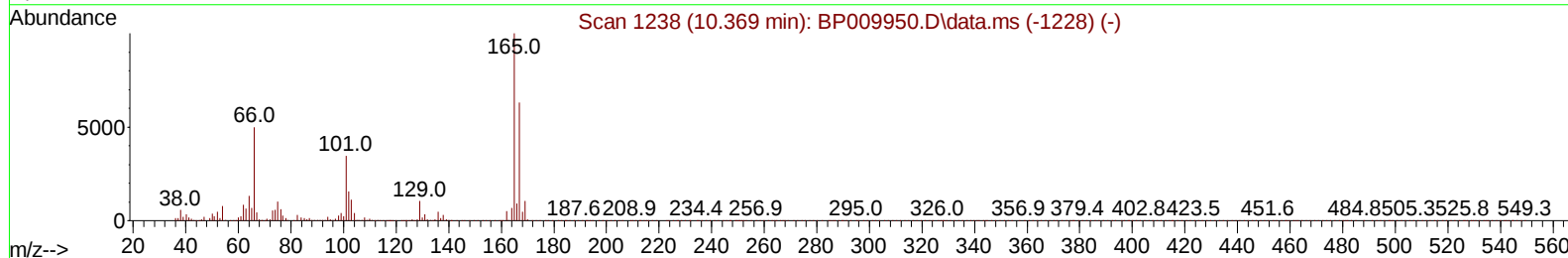
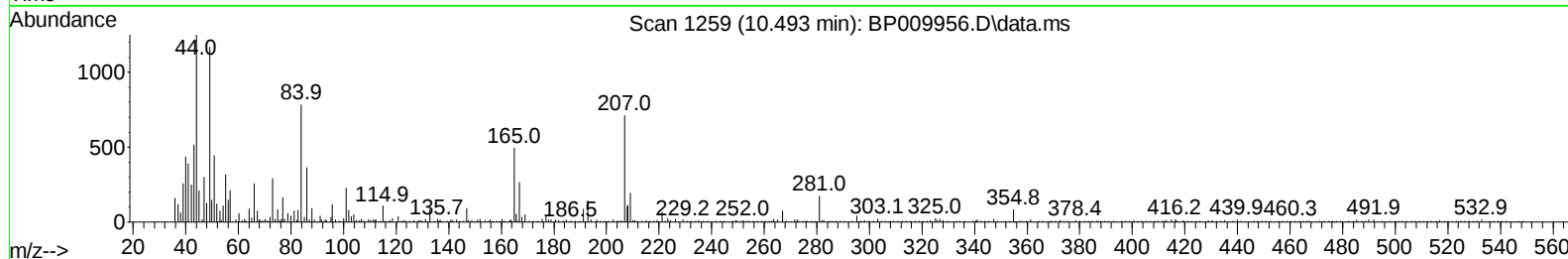
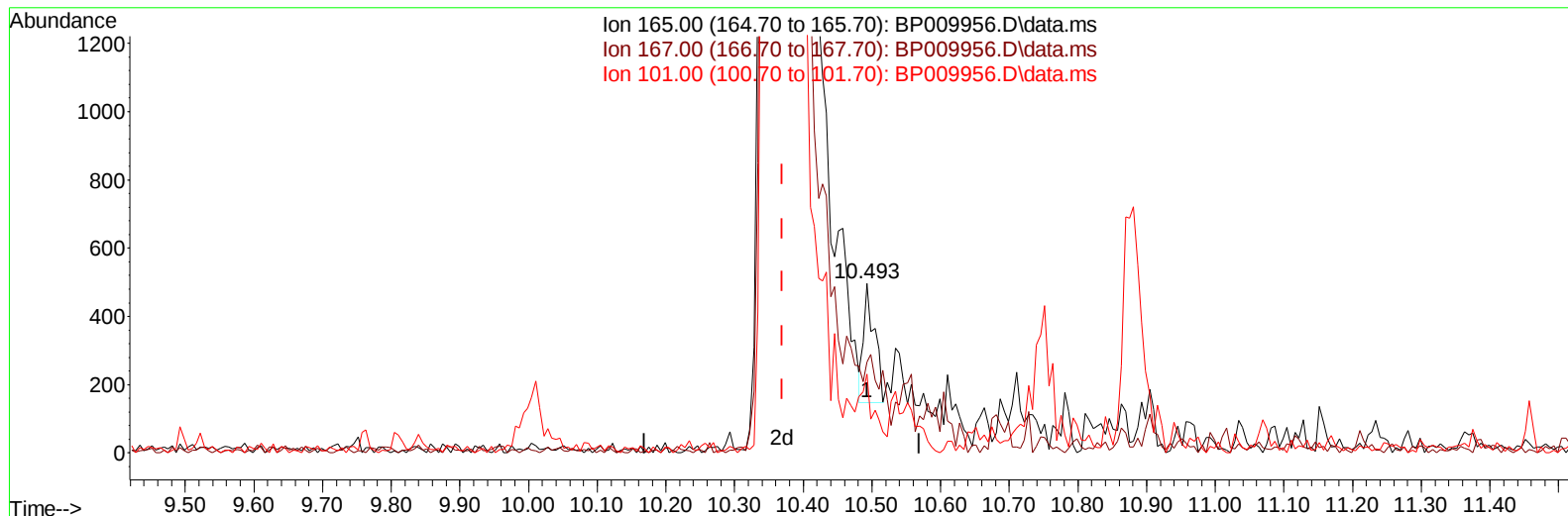
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Instrument :
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TIC: BP009956.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.493min (+ 0.123) 0.04 ng/u1

response 390

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	48.10
101.00	49.80	41.77
0.00	0.00	0.00

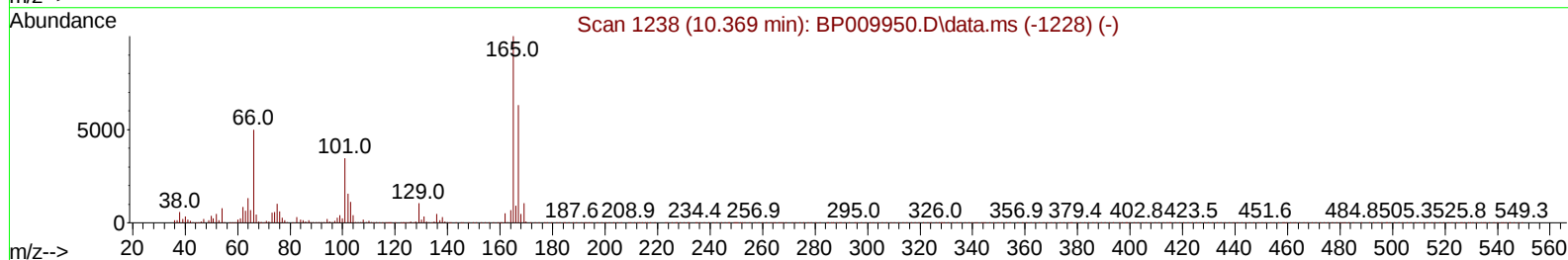
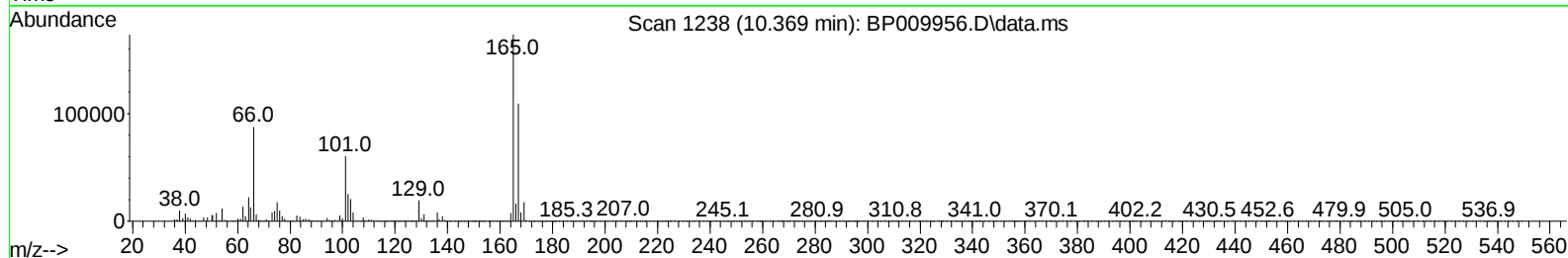
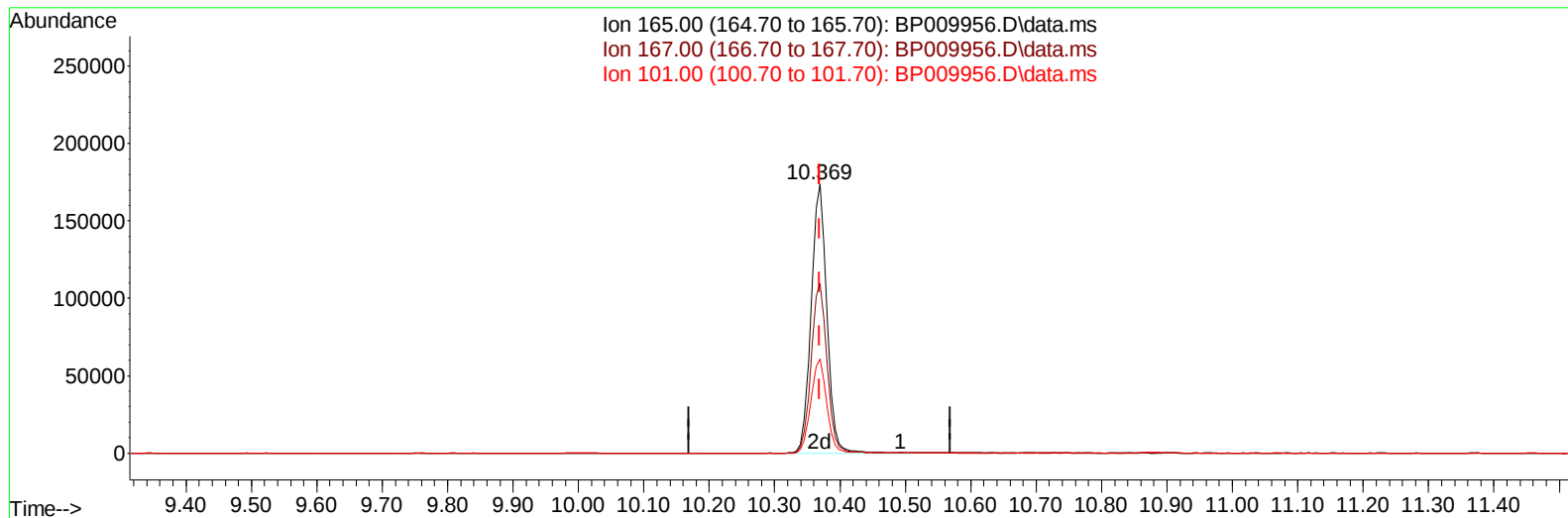
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Instrument :
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ClientSampleId :
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TIC: BP009956.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.369min (-0.000) 32.87 ng/ul m

response 291056

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	63.03
101.00	49.80	34.95#
0.00	0.00	0.00

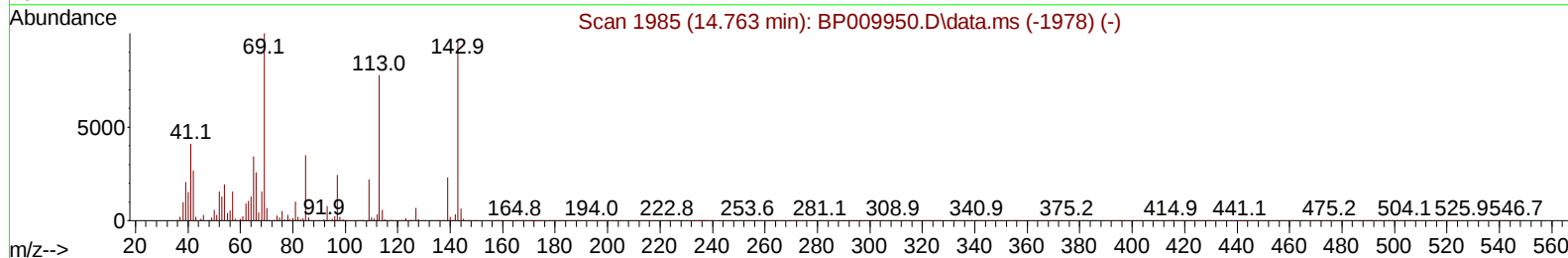
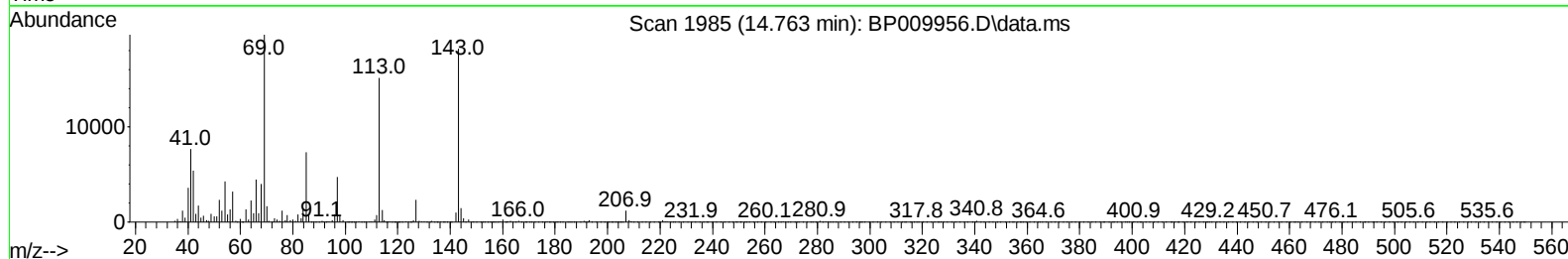
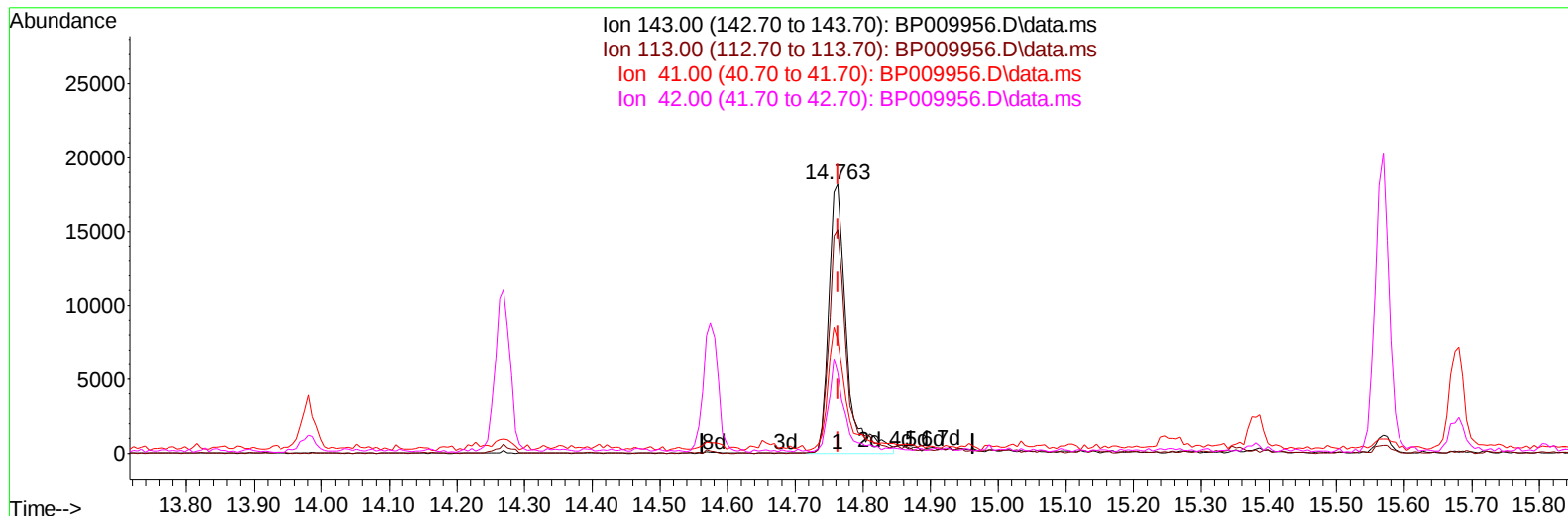
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Instrument :
 BNA_P
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TIC: BP009956.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 5.94 ng/ul m

response 33335

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.04#
41.00	23.20	42.13#
42.00	18.00	29.82#

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 Misc :
 ALS Vial : 8 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	119526	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	535147	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	393627	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	869206	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240	878696	20.000	ng/ul	# 0.00
88) Perylene-d12	23.863	264	841649	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	16327m	6.096	ng/uL	0.00
4) Pyridine-d5	3.793	84	64112	8.682	ng/ul	0.00
7) Phenol-d5	7.099	99	88024	8.388	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	250795	40.145	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	241904	30.561	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	174162	19.882	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	164908	42.732	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	162579	38.426	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	291056m	32.875	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	398432	31.597	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1261842	40.810	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1365944	37.289	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	33335m	5.945	ng/ul	0.00
60) Fluorene-d10	15.569	176	1078640	41.469	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	120629	23.427	ng/ul	0.00
73) Anthracene-d10	17.428	188	1761021	42.315	ng/ul	0.00
81) Pyrene-d10	19.657	212	2132558	44.437	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	1844191	42.076	ng/ul	0.00

Target Compounds Qvalue

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

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