

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009958.D
 Acq On : 19 Apr 2022 22:33
 Operator : CG/JU
 Sample : N2421-19
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J92

Manual IntegrationsAPPROVED

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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
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Reviewed By :Jagrut

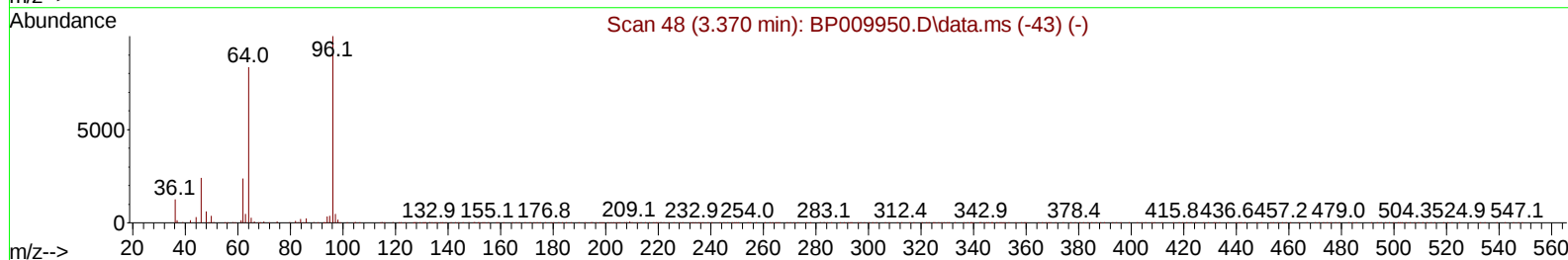
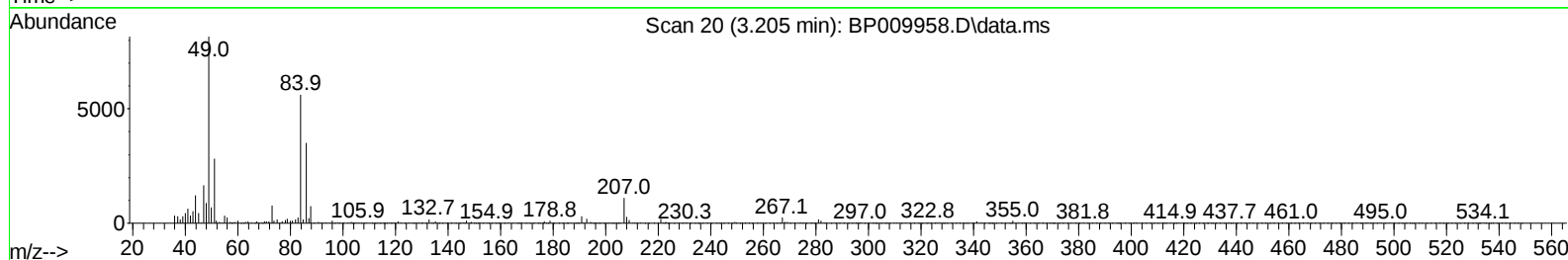
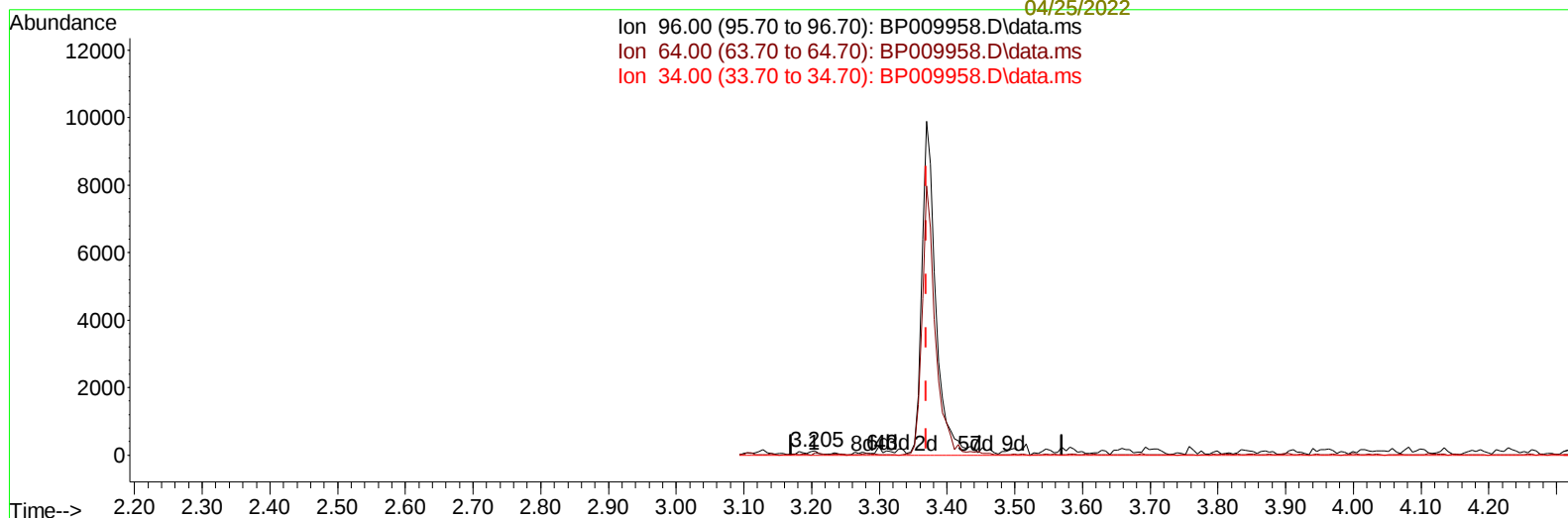
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Supervised By :Yogesh

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04/25/2022



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

3.205min (-0.165) 0.03 ng/uL

response 70

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	63.79
34.00	0.00	0.00
0.00	0.00	0.00

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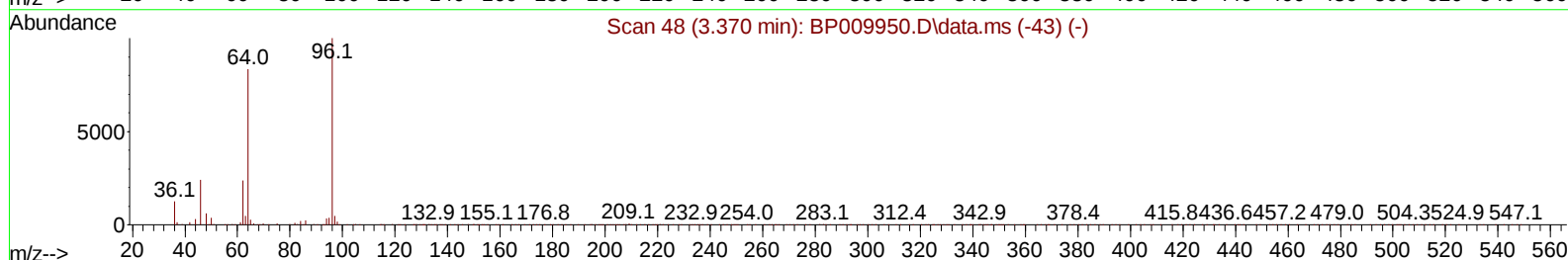
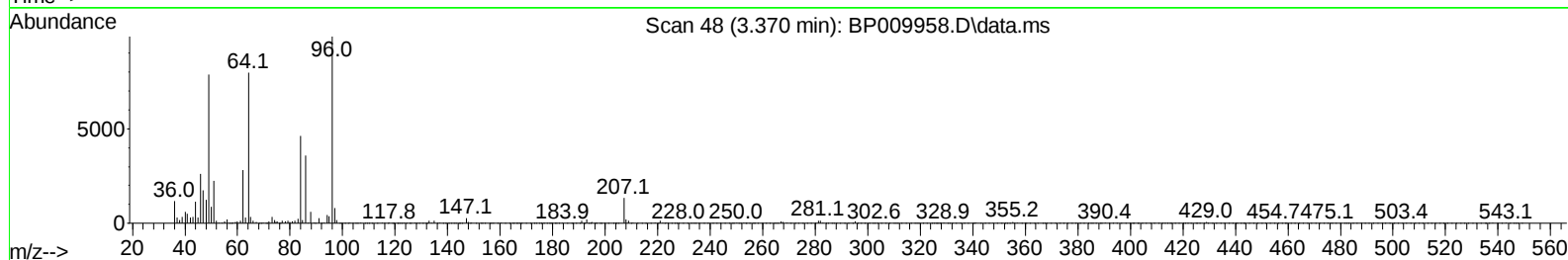
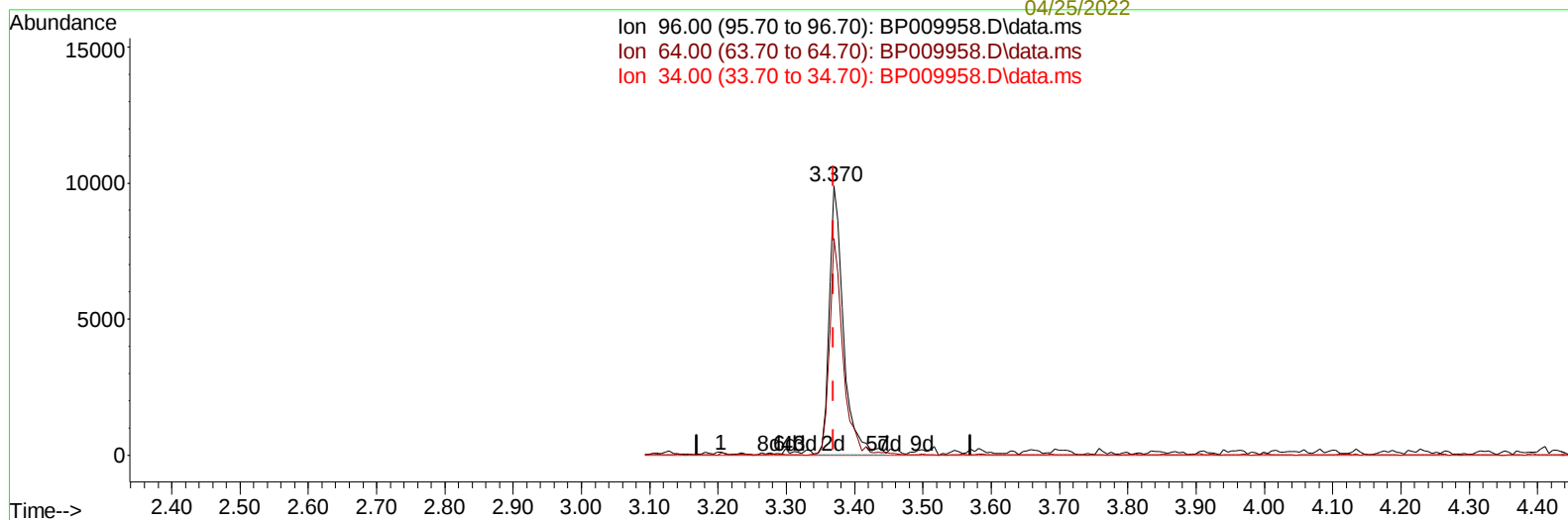
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(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 5.41 ng/uL m

response 14069

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.58#
34.00	0.00	0.00
0.00	0.00	0.00

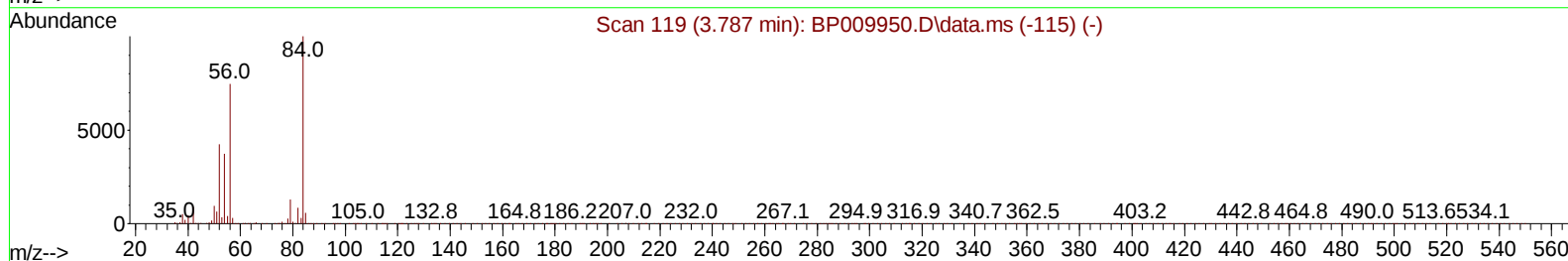
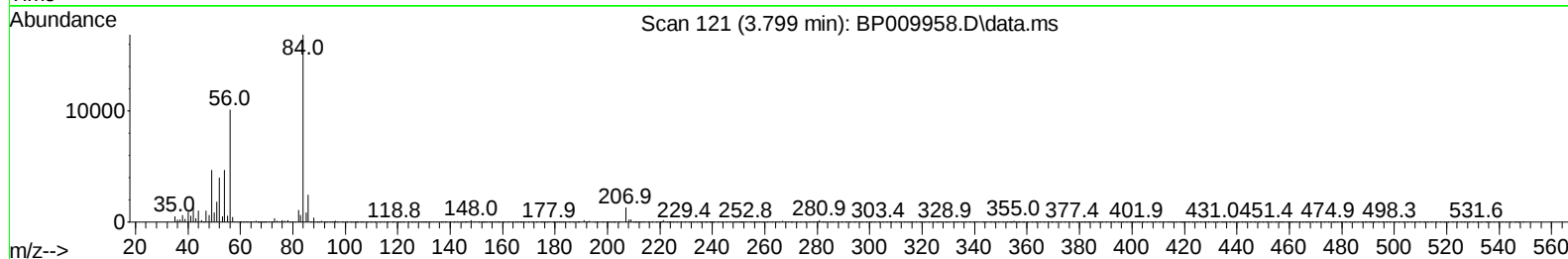
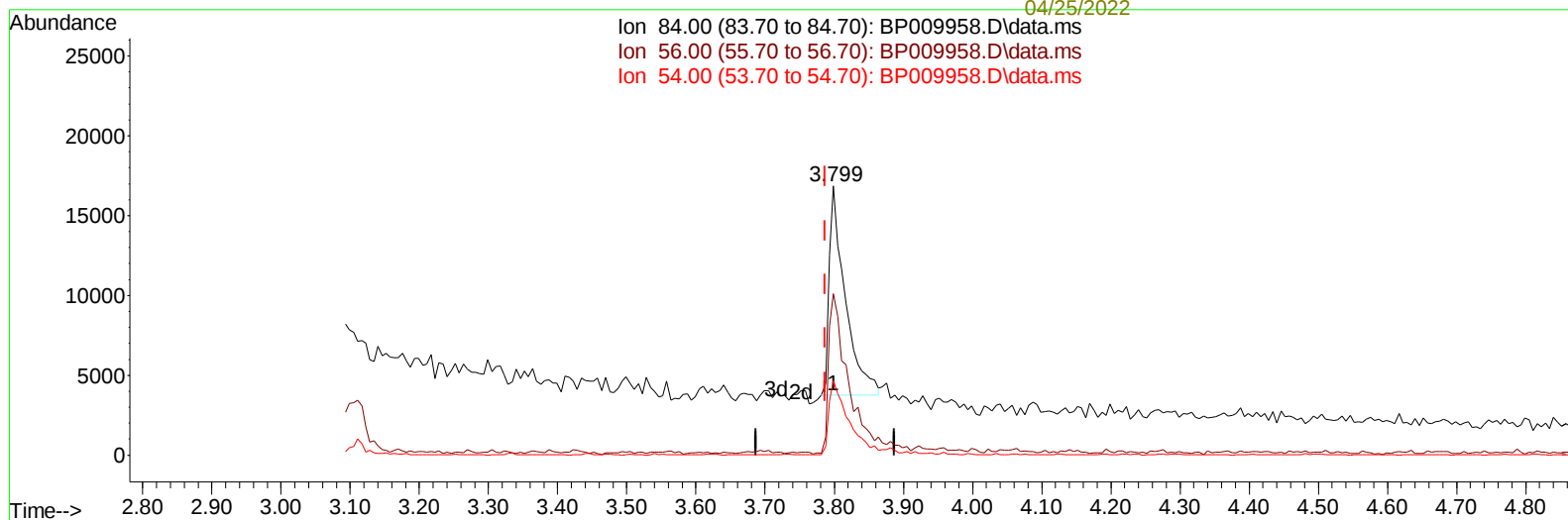
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(4) Pyridine-d5 (S)

3.799min (+ 0.012) 2.95 ng/u1

response 21105

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	60.00#
54.00	24.80	27.69
0.00	0.00	0.00

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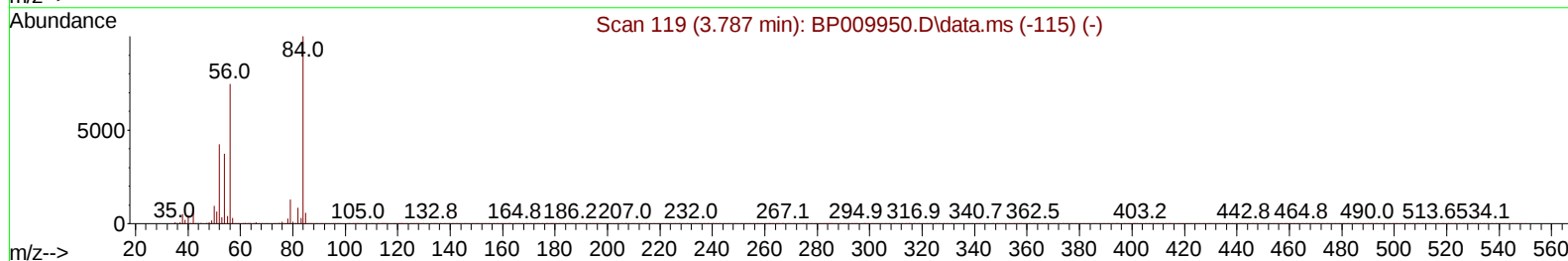
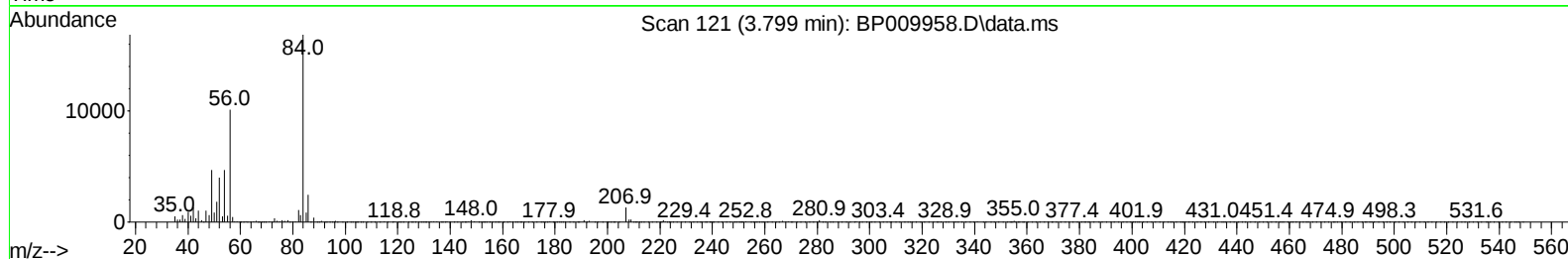
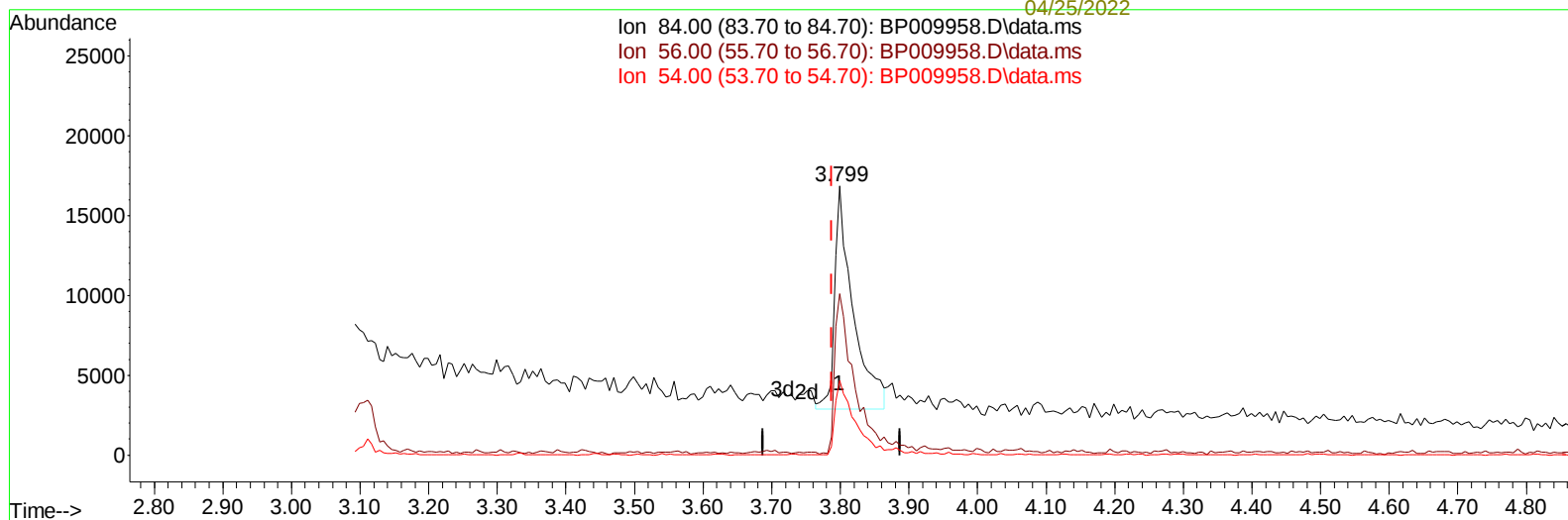
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TIC: BP009958.D\data.ms

(4) Pyridine-d5 (S)

3.799min (+ 0.012) 3.64 ng/ul m

response 26056

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	60.00#
54.00	24.80	27.69
0.00	0.00	0.00

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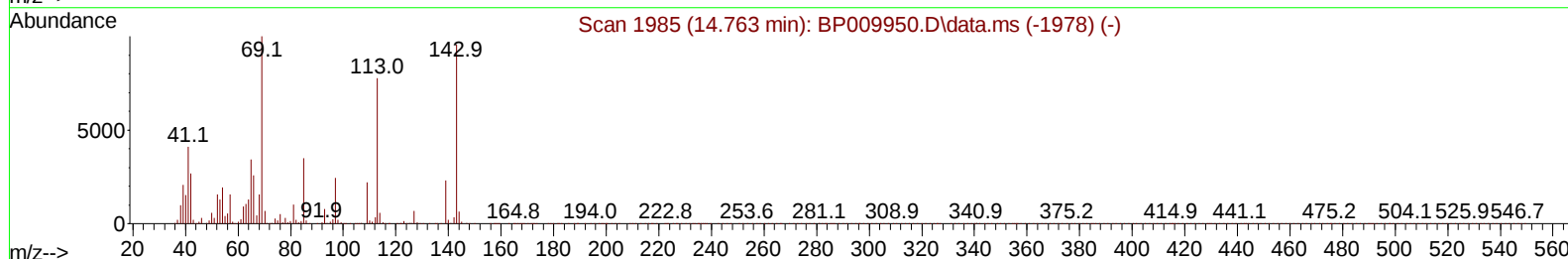
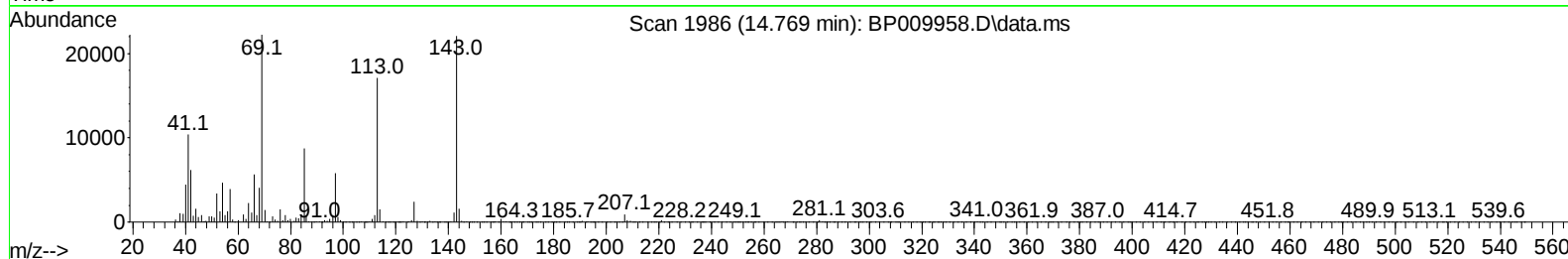
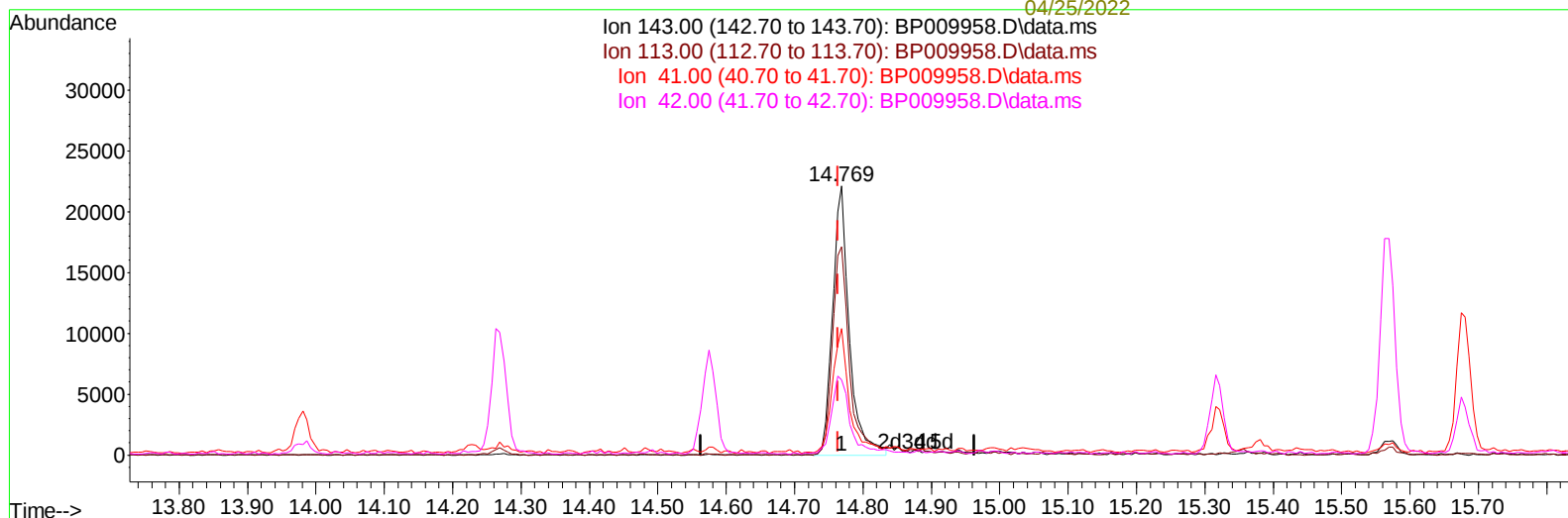
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(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 6.76 ng/u1

response 37335

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.41#
41.00	23.20	46.94#
42.00	18.00	28.05#

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.946	152		115957	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136		526895	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164		387729	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188		845272	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240		856950	20.000	ng/ul	# 0.00
88) Perylene-d12	23.863	264		821854	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.370	96		14069m	5.415	ng/uL	0.00
4) Pyridine-d5	3.799	84		26056m	3.637	ng/ul	0.01
7) Phenol-d5	7.099	99		80284	7.886	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67		228073	37.632	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132		214981	27.995	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113		158856	18.692	ng/ul	0.00
21) Nitrobenzene-d5	9.105	128		146994	38.687	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143		156673	37.610	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165		266127	30.530	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131		255323	20.565	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166		1135297	37.276	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160		1237107	34.285	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143		38369m	6.947	ng/ul	0.00
60) Fluorene-d10	15.569	176		993868	38.791	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200		184789	36.903	ng/ul	0.00
73) Anthracene-d10	17.428	188		1739887	42.991	ng/ul	0.00
81) Pyrene-d10	19.657	212		2145724	45.846	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264		1877302	43.863	ng/ul	0.00

Target Compounds Qvalue

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

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