

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010070.D
 Acq On : 22 Apr 2022 23:55
 Operator : CG/JU
 Sample : N2475-05
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

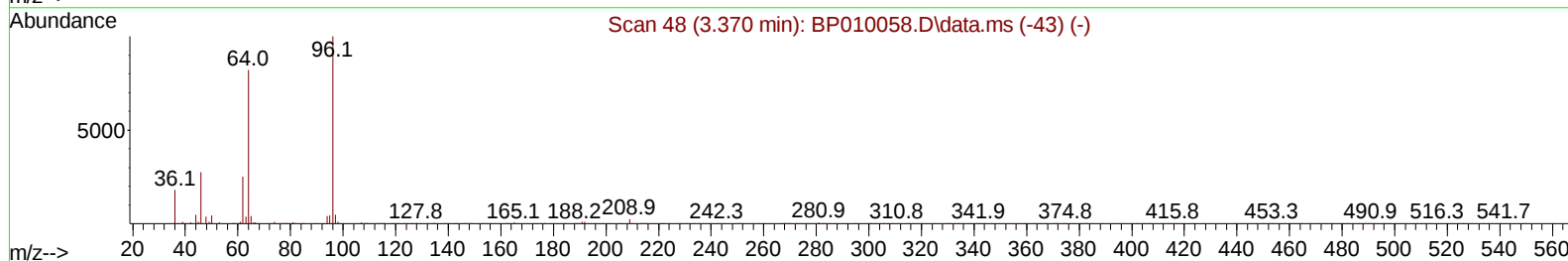
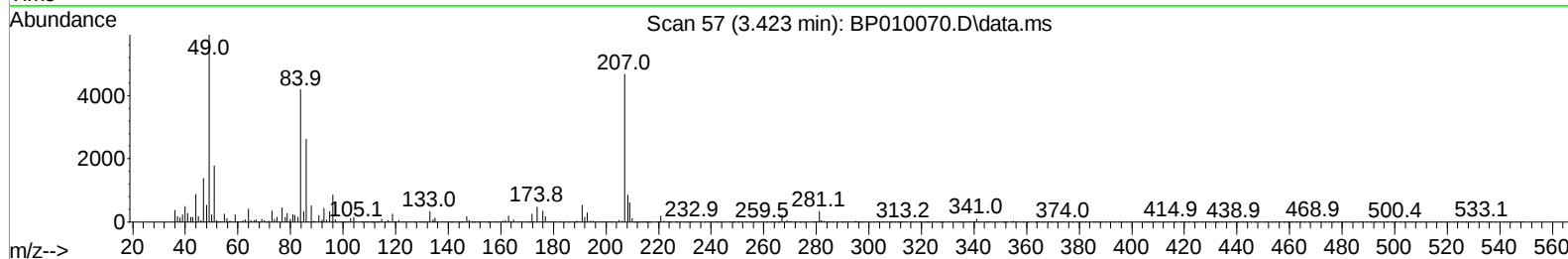
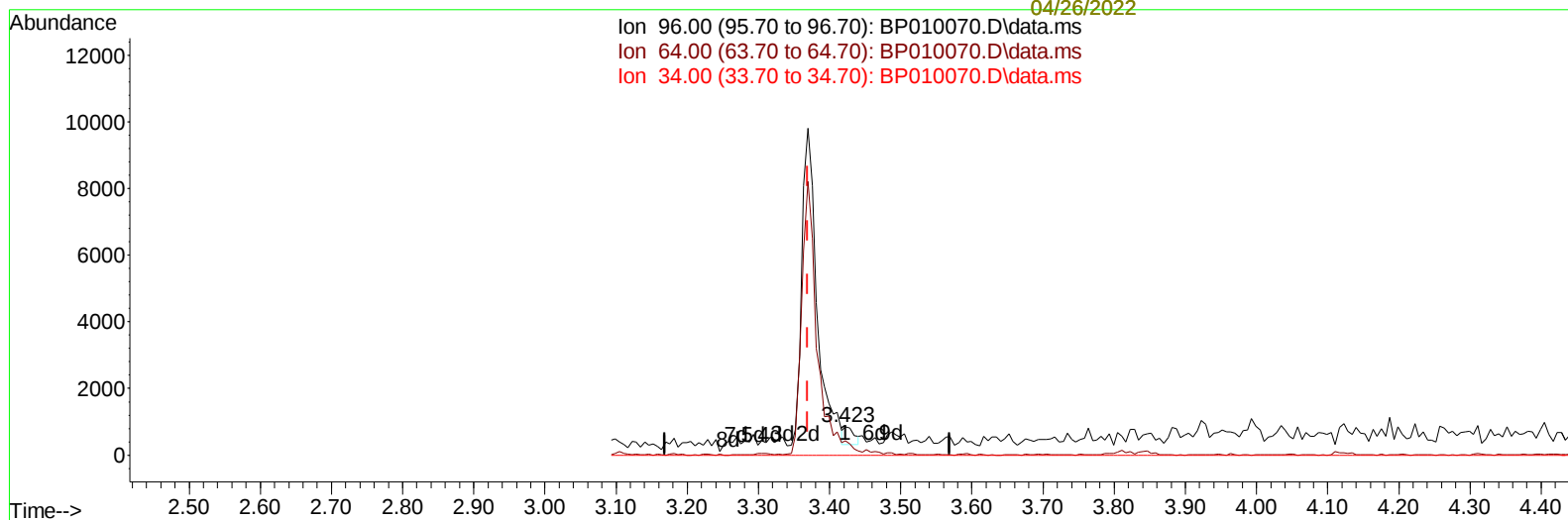
Instrument :
 BNA_P
 ClientSampleId :
 A0BN6

Manual IntegrationsAPPROVED

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 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut
 Upadhyay
 04/25/2022

Supervised By :mohammad
 ahmed
 04/26/2022



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

3.423min (+ 0.054) 0.17 ng/uL

response 550

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	49.42
34.00	0.00	0.00
0.00	0.00	0.00

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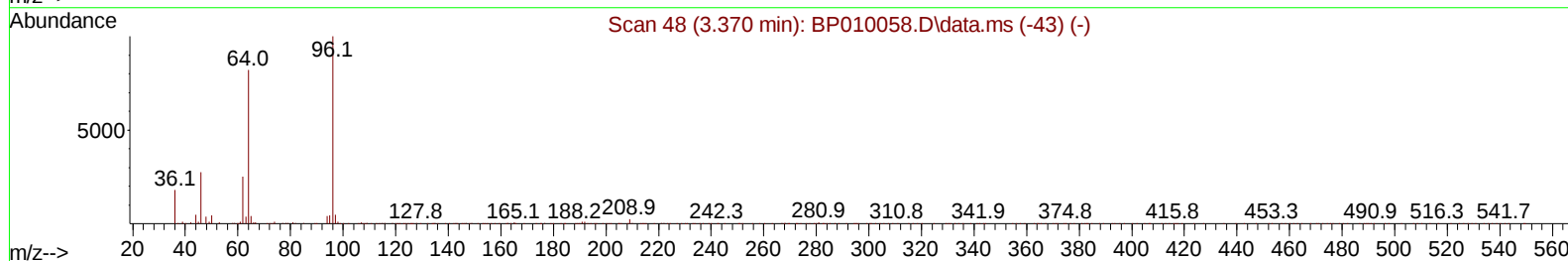
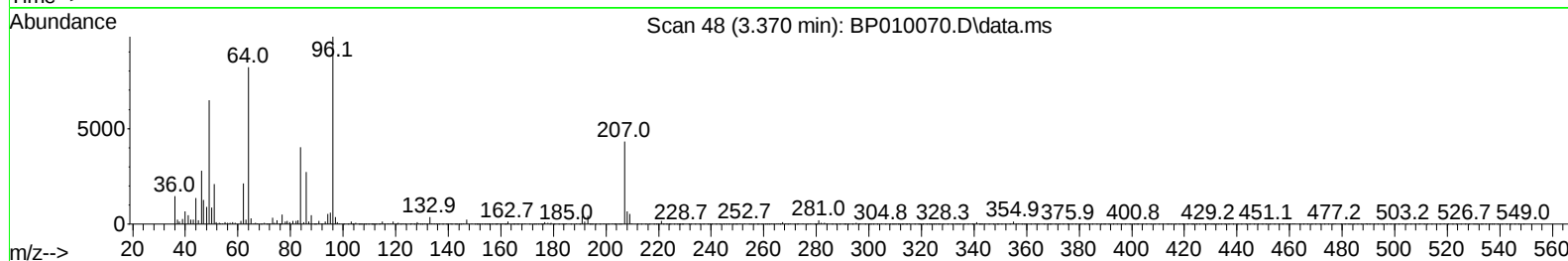
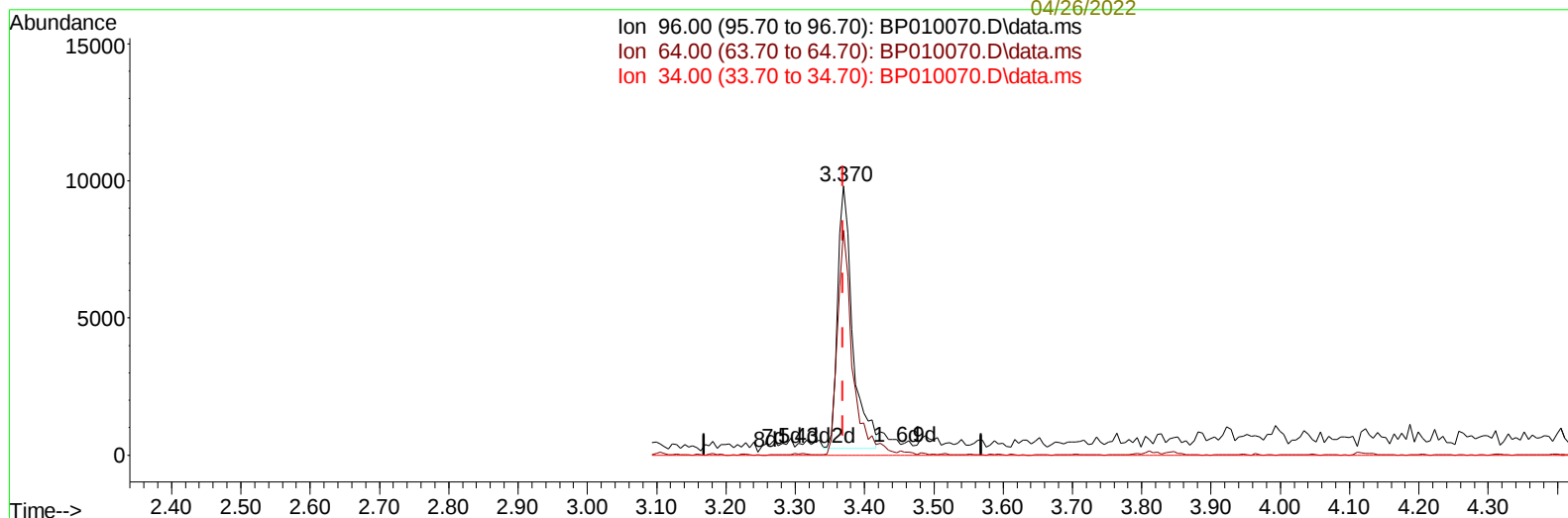
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(3) 1,4-Dioxane-d8 (S)

3.370min (+ 0.001) 4.42 ng/uL m

response 14392

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.67#
34.00	0.00	0.00
0.00	0.00	0.00

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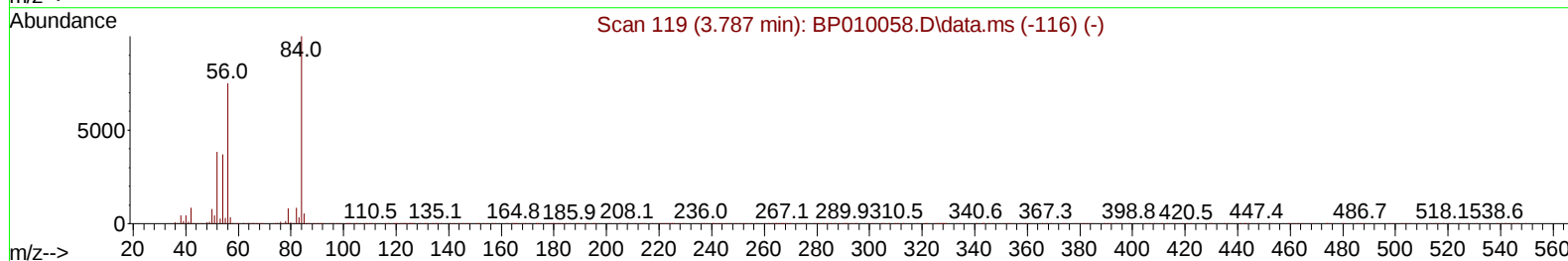
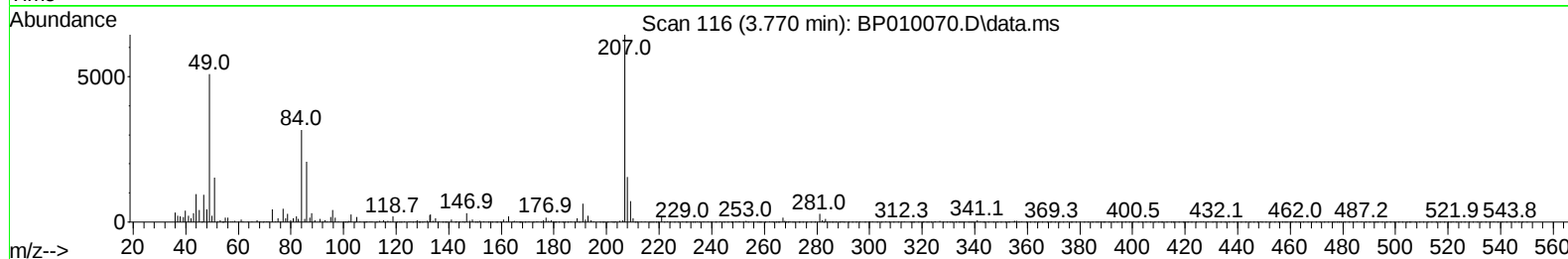
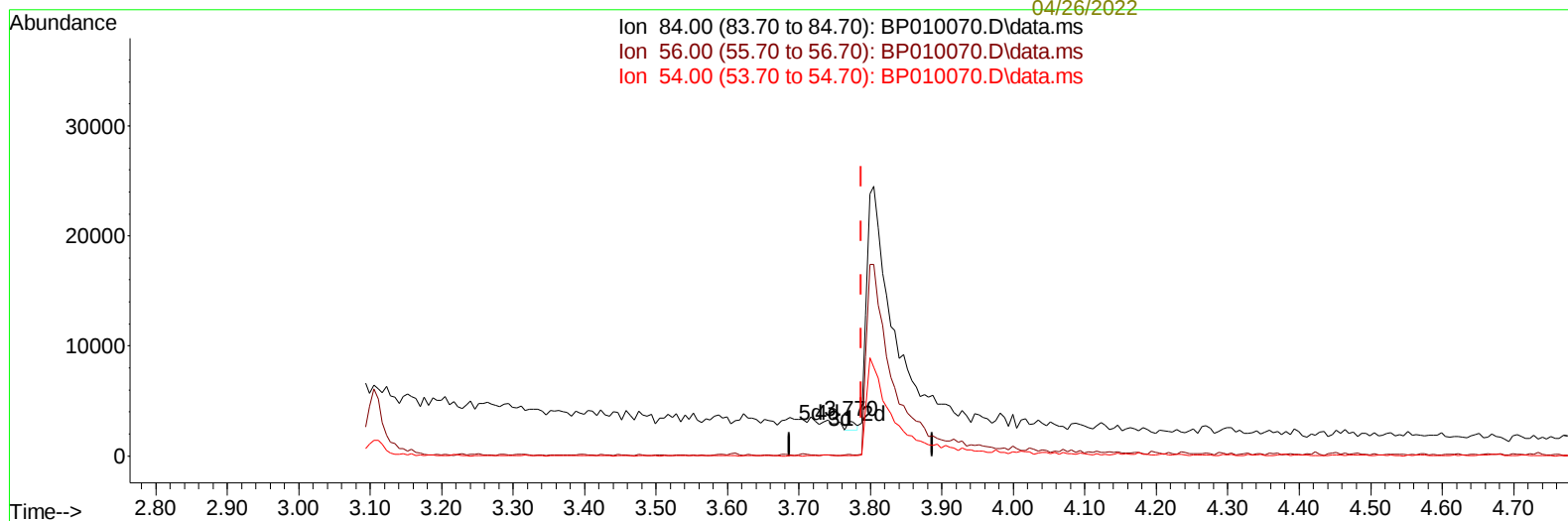
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(4) Pyridine-d5 (S)

3.770min (-0.017) 0.08 ng/u1

response 729

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	4.74#
54.00	24.80	0.38#
0.00	0.00	0.00

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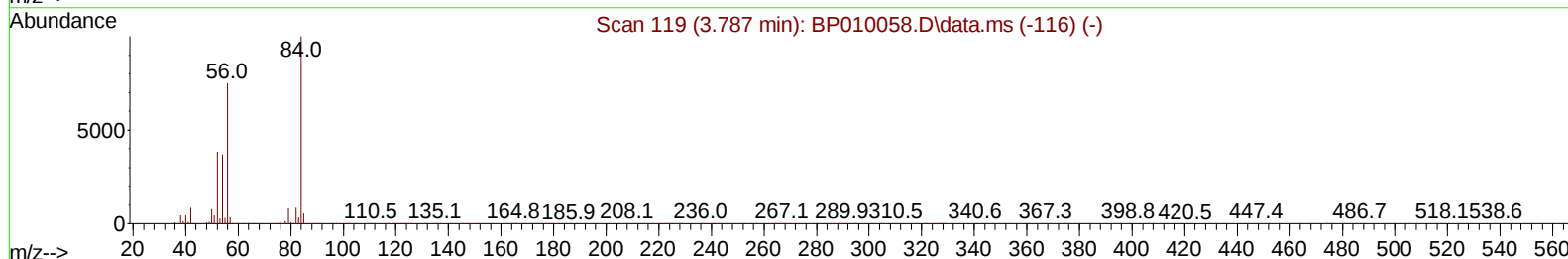
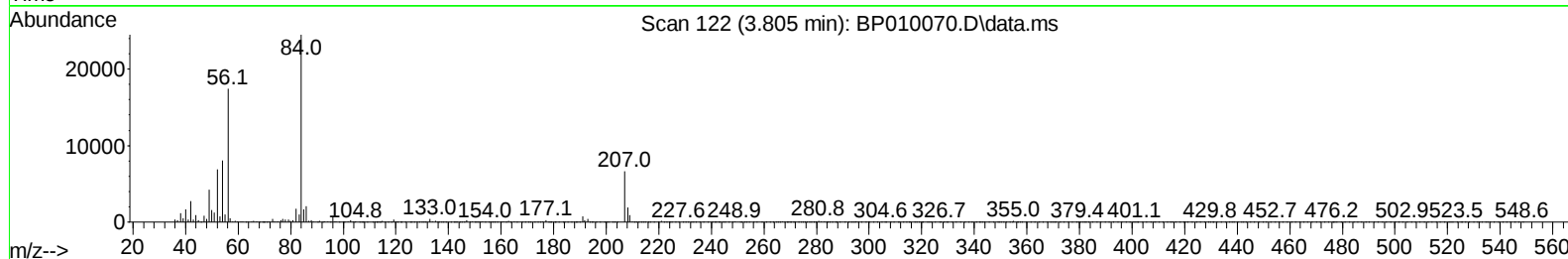
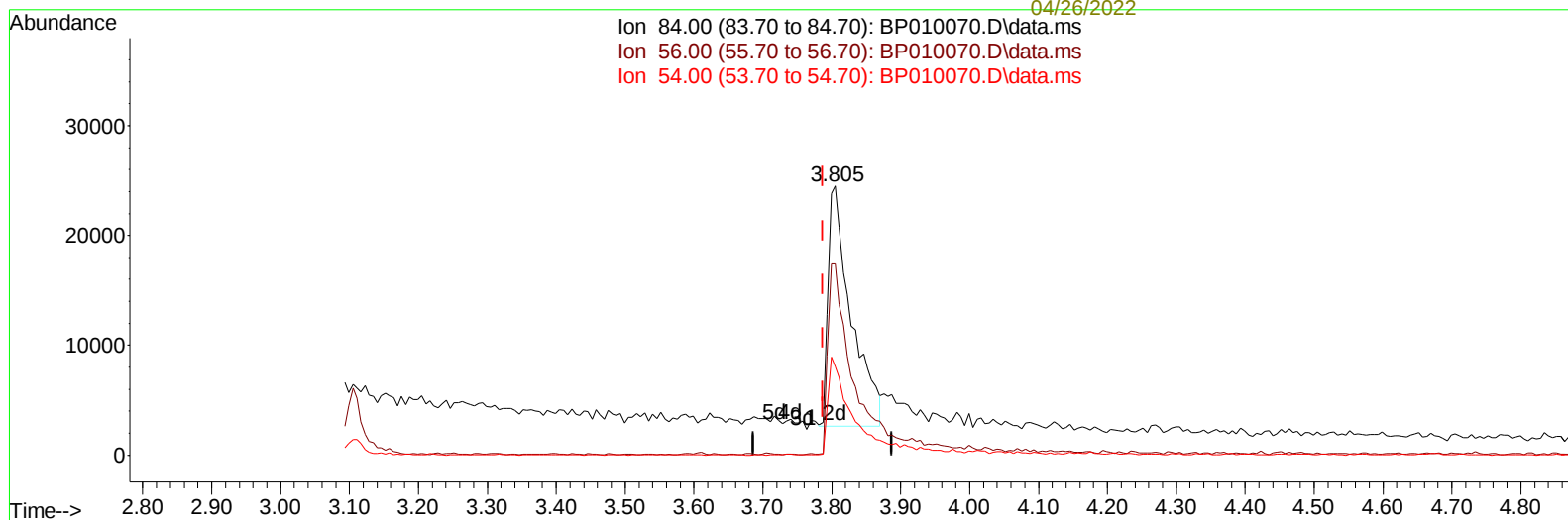
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(4) Pyridine-d5 (S)

3.805min (+ 0.018) 5.56 ng/u1 m

response 50937

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	71.10#
54.00	24.80	32.98#
0.00	0.00	0.00

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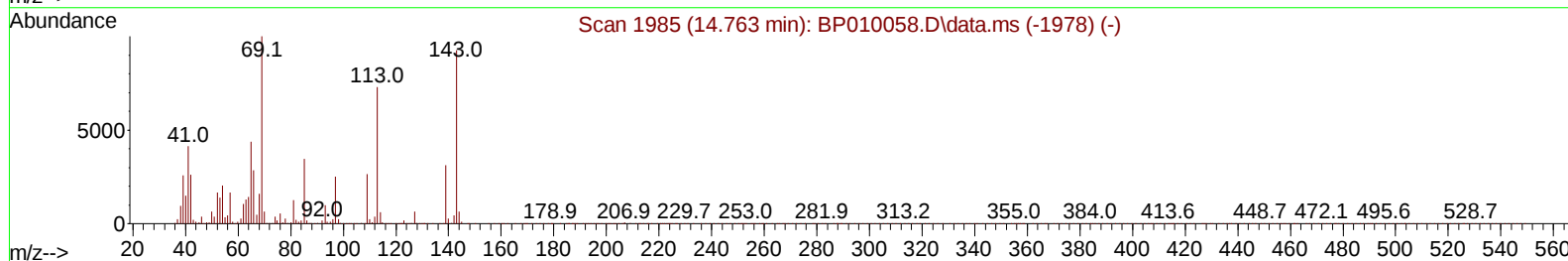
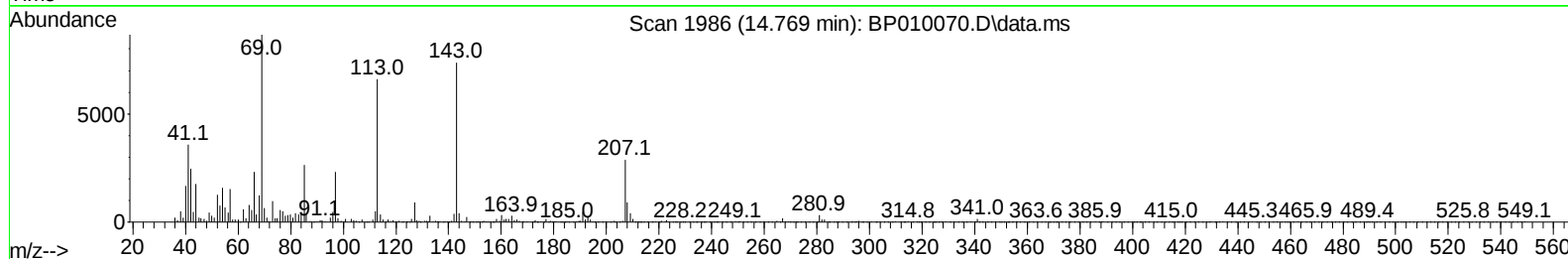
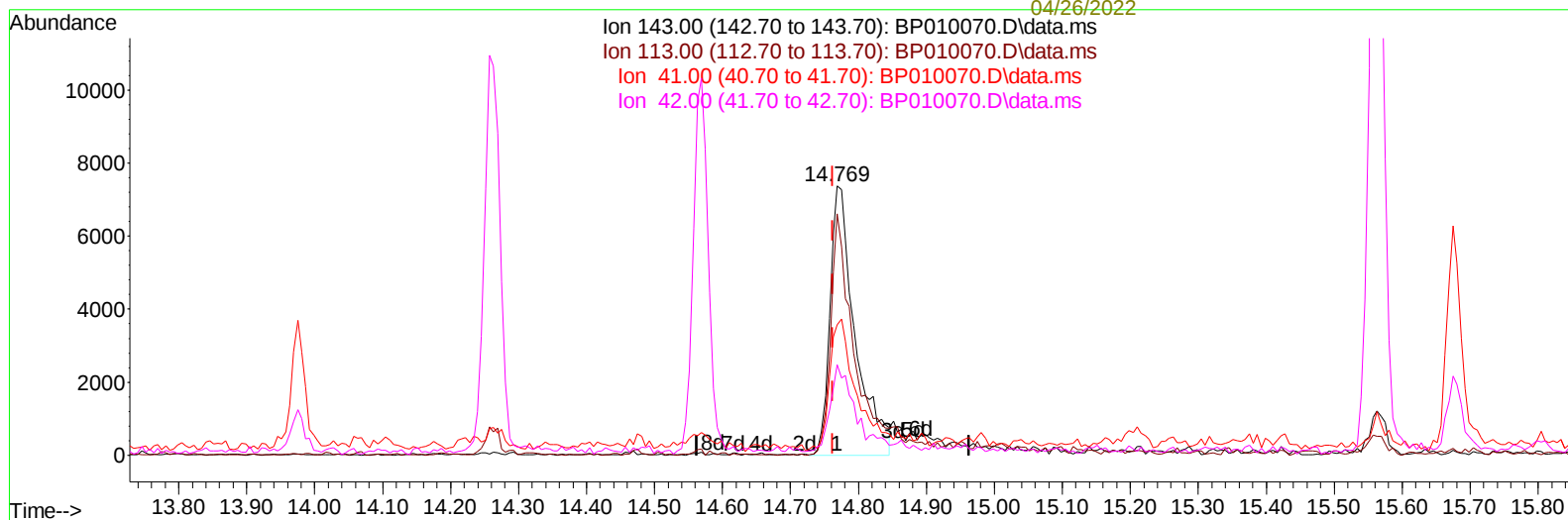
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(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 2.68 ng/ul m

response 19007

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	89.60#
41.00	23.20	48.43#
42.00	18.00	33.51#

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.940	152		141881	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136		676701	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164		494690	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188		1084514	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240		833137	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264		709272	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.370	96		14392m	4.423	ng/uL	0.00
4) Pyridine-d5	3.805	84		50937m	5.565	ng/ul	0.02
7) Phenol-d5	7.099	99		74995	5.811	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67		234924	30.337	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132		224585	23.498	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113		162682	15.215	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128		161363	29.981	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143		154733	26.324	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165		295411	25.816	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131		343116	20.912	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166		1214389	31.558	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160		1358354	29.166	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143		19007m	2.683	ng/ul	0.00
60) Fluorene-d10	15.563	176		1071306	32.364	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200		96429	14.062	ng/ul	0.00
73) Anthracene-d10	17.422	188		1709259	32.577	ng/ul	0.00
81) Pyrene-d10	19.651	212		1965730	41.488	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264		1281531	34.465	ng/ul	0.00

Target Compounds Qvalue

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

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