

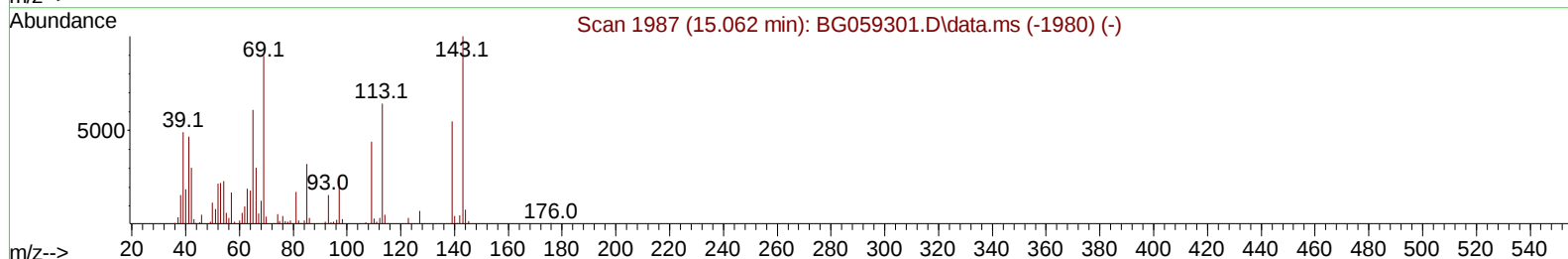
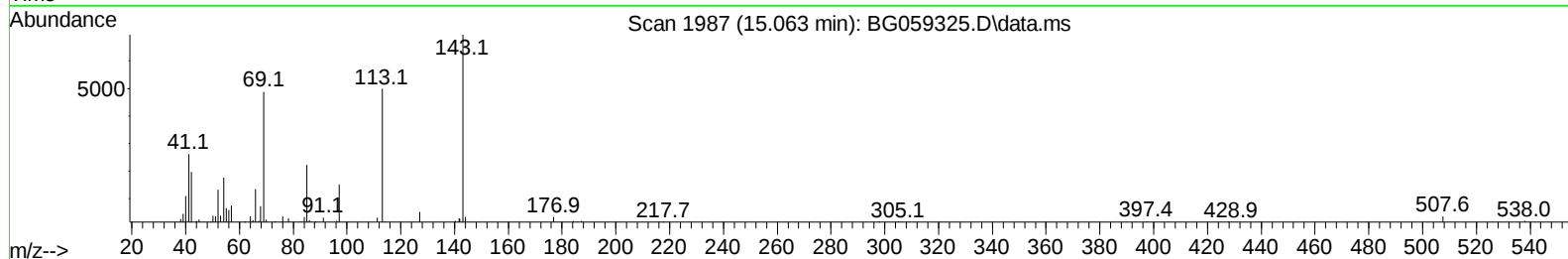
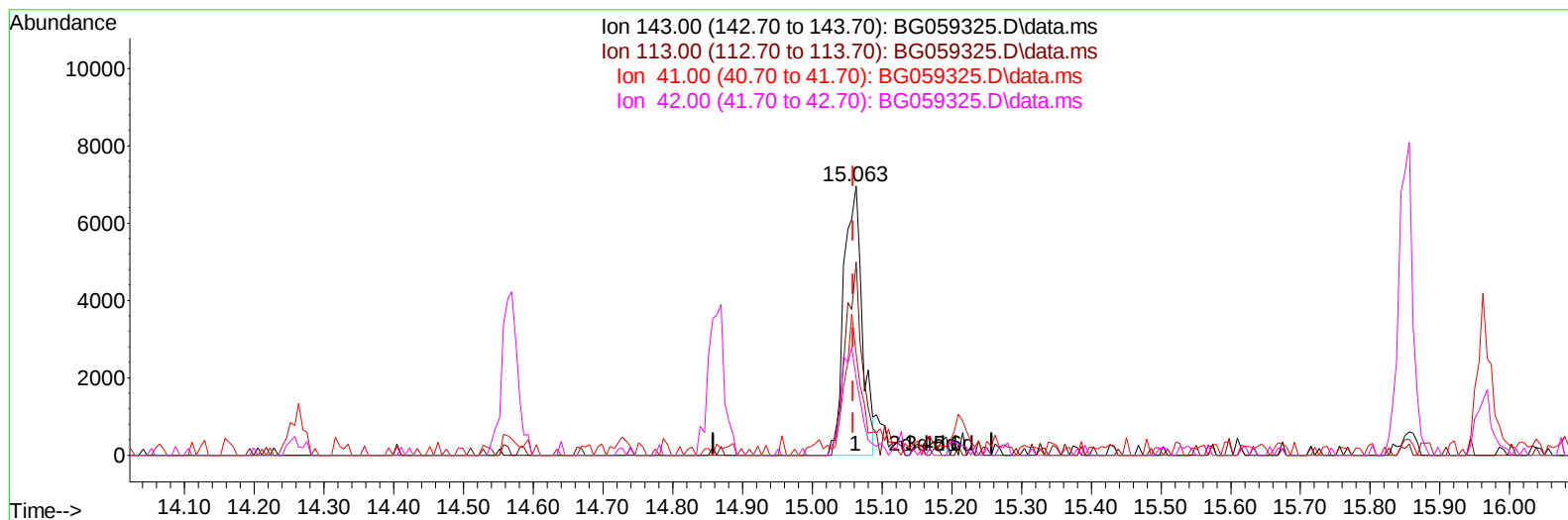
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100523\
Data File : BG059325.D
Acq On : 6 Oct 2023 11:14
Operator : MA/JU
Sample : 04592-05 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZY3

Manual IntegrationsAPPROVED

Quant Time: Oct 06 22:28:57 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 28 04:33:25 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/07/2023
Supervised By :mohammad ahmed 10/07/2023



TIC: BG059325.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.063min (+ 0.004) 6.45 ng/u1

response 12631

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.10	71.70
41.00	57.70	37.69#
42.00	37.50	28.37#

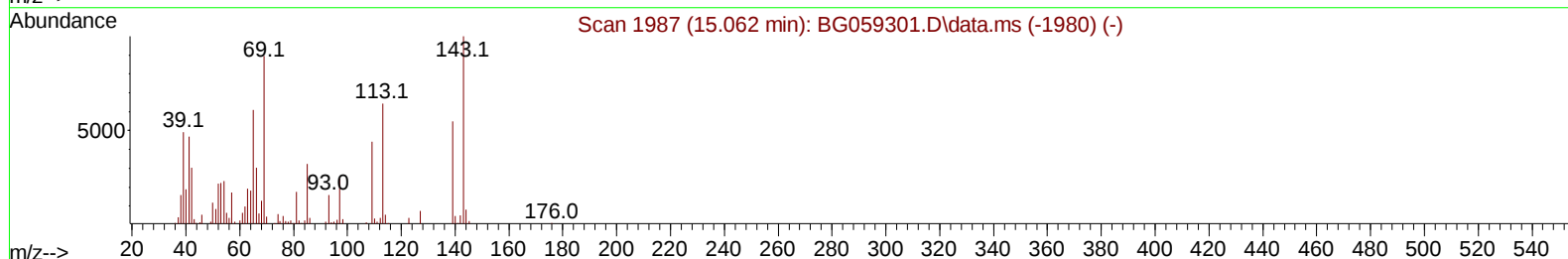
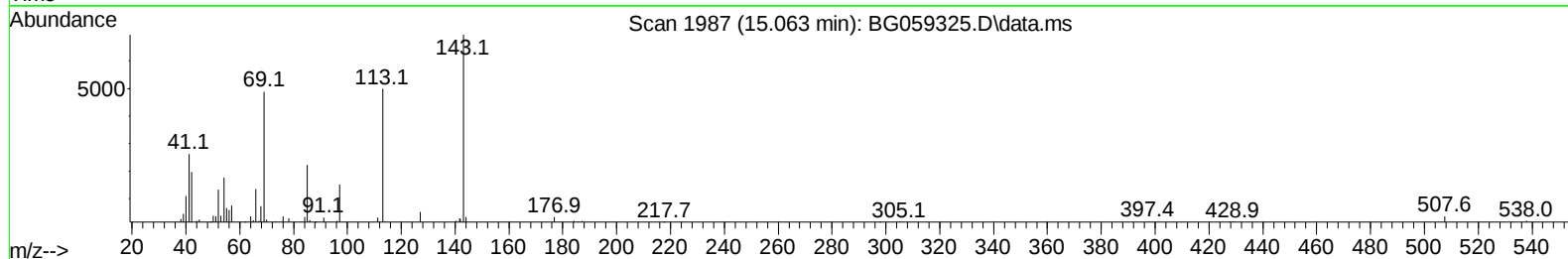
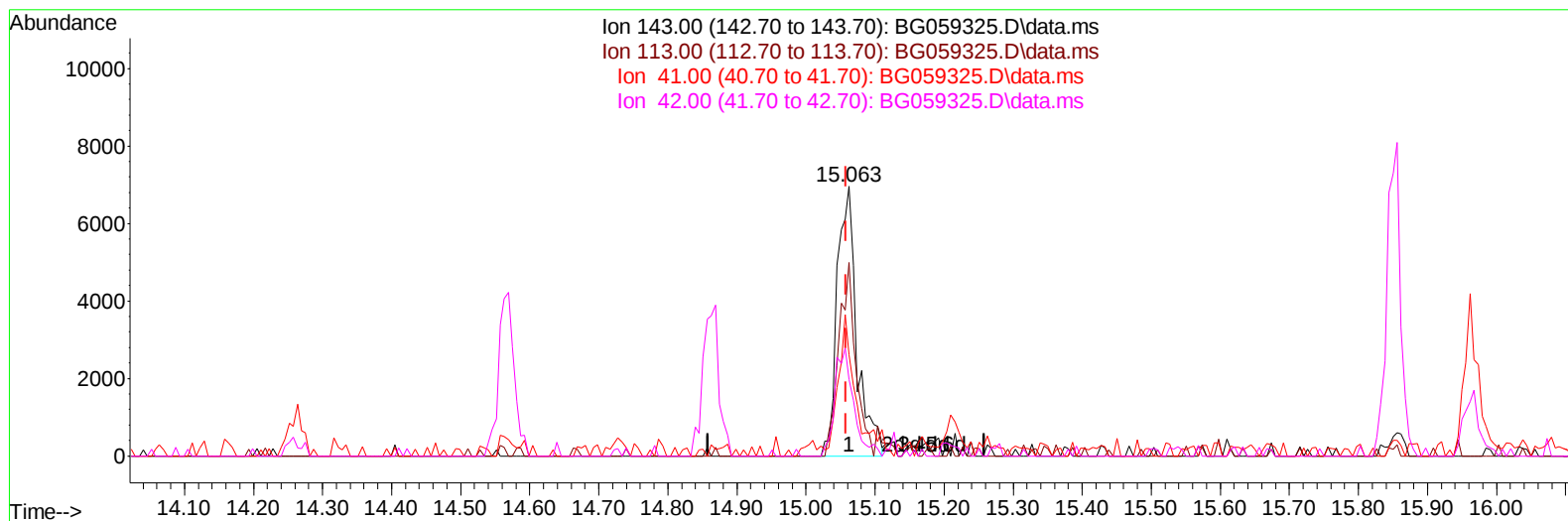
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TIC: BG059325.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.063min (+ 0.004) 6.96 ng/ul m

response 13633

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.10	71.70
41.00	57.70	37.69#
42.00	37.50	28.37#

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 EZYY3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/07/2023
 Supervised By :mohammad ahmed 10/07/2023

Quant Time: Oct 06 22:44:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 28 04:33:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.229	152	56658	20.000	ng/ul	0.00
20) Naphthalene-d8	11.073	136	262501	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.863	164	155068	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.613	188	350852	20.000	ng/ul	0.00
79) Chrysene-d12	21.913	240	319596	20.000	ng/ul	0.00
88) Perylene-d12	25.398	264	373174	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.529	96	3026	2.016	ng/uL	0.00
4) Pyridine-d5	3.970	84	46620	10.643	ng/ul	0.00
7) Phenol-d5	7.360	99	41228	7.270	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.554	67	106133	29.295	ng/ul	0.00
11) 2-Chlorophenol-d4	7.748	132	112905	29.616	ng/ul	0.00
15) 4-Methylphenol-d8	8.929	113	70782	16.146	ng/ul	0.00
21) Nitrobenzene-d5	9.428	128	72964	39.390	ng/ul	0.00
24) 2-Nitrophenol-d4	10.151	143	82936	45.608	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.674	165	129080	34.554	ng/ul	0.00
31) 4-Chloroaniline-d4	11.220	131	186802	30.907	ng/ul	0.00
46) Dimethylphthalate-d6	14.258	166	404441	36.074	ng/ul	0.00
49) Acenaphthylene-d8	14.569	160	504507	35.851	ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	13633m	6.963	ng/ul	0.00
60) Fluorene-d10	15.850	176	393825	36.907	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.967	200	56957	31.582	ng/ul	0.00
73) Anthracene-d10	17.712	188	614486	38.438	ng/ul	0.00
81) Pyrene-d10	19.986	212	673731	37.643	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.157	264	670685	37.400	ng/ul	0.00

Target Compounds Qvalue

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

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