

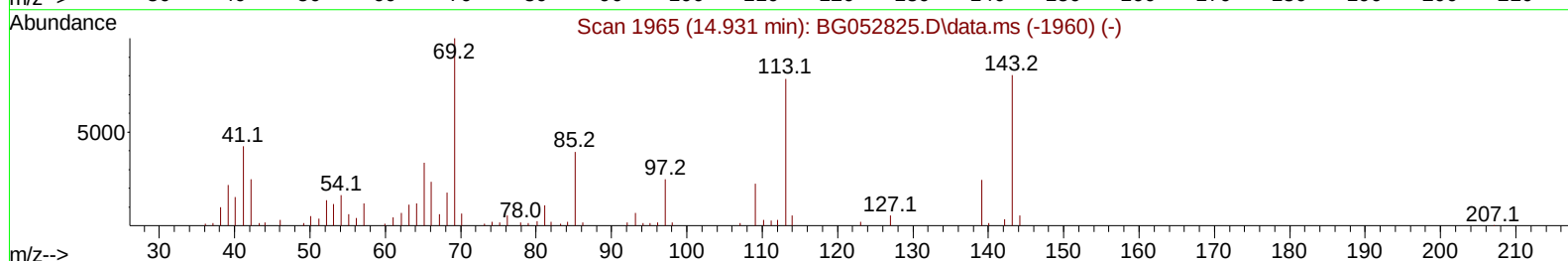
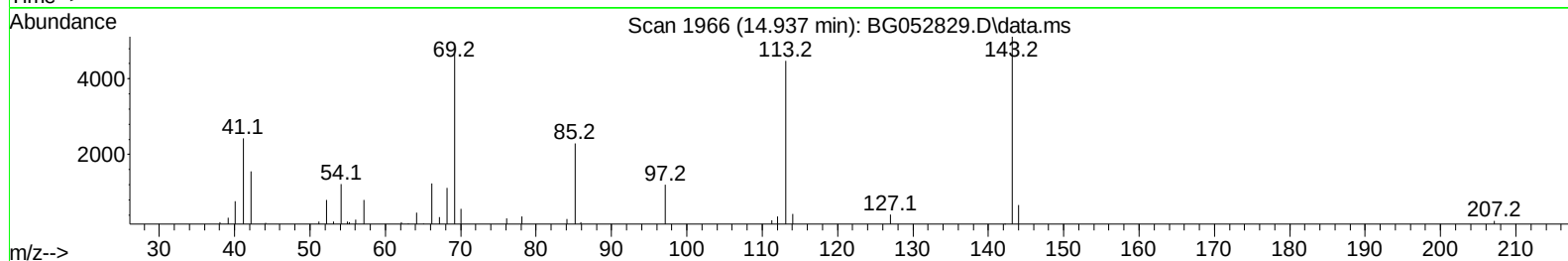
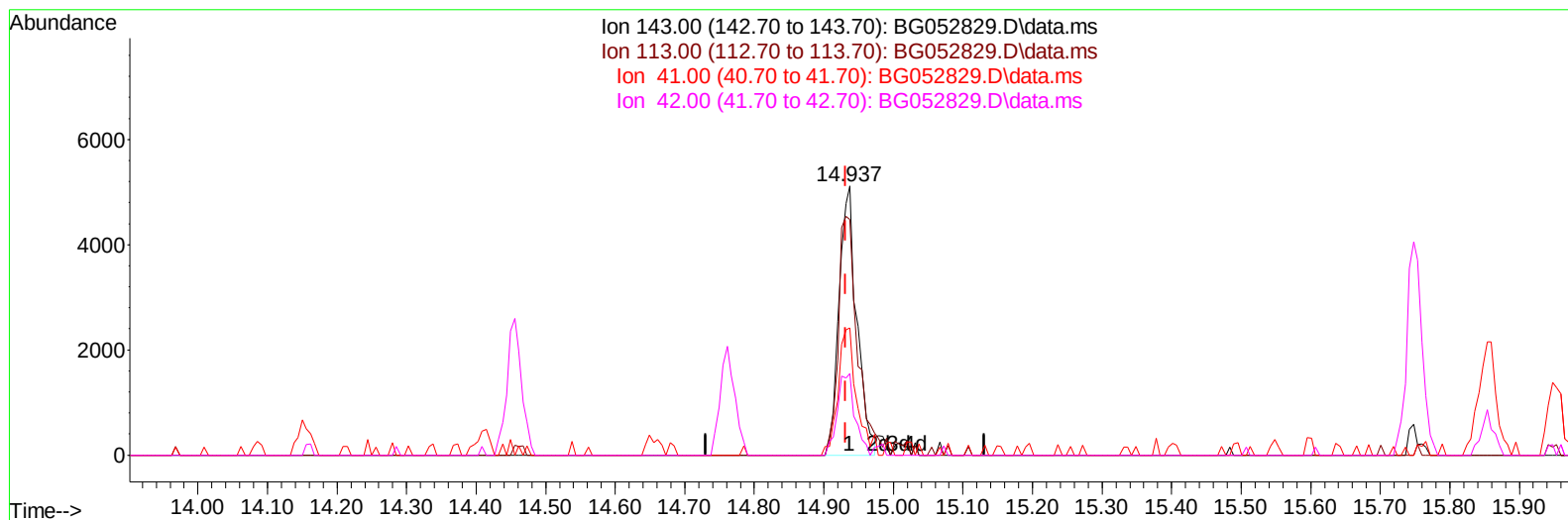
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
Data File : BG052829.D
Acq On : 25 Mar 2022 14:46
Operator : CG/JU
Sample : N2079-19
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW607

Manual IntegrationsAPPROVED

Quant Time: Mar 25 15:19:57 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 23 13:16:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/28/2022
Supervised By :mohammad ahmed 03/29/2022



TIC: BG052829.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.937min (+ 0.006) 8.14 ng/ul

response 9150

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	87.68
41.00	47.20	47.35
42.00	29.40	30.28

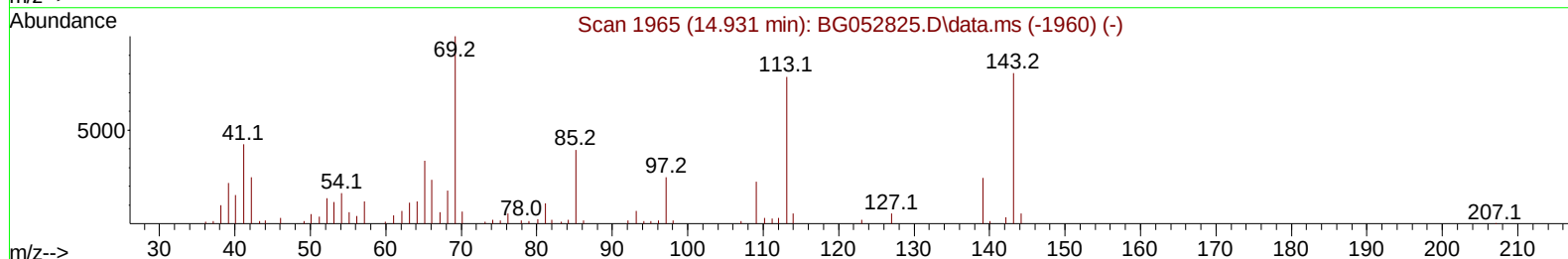
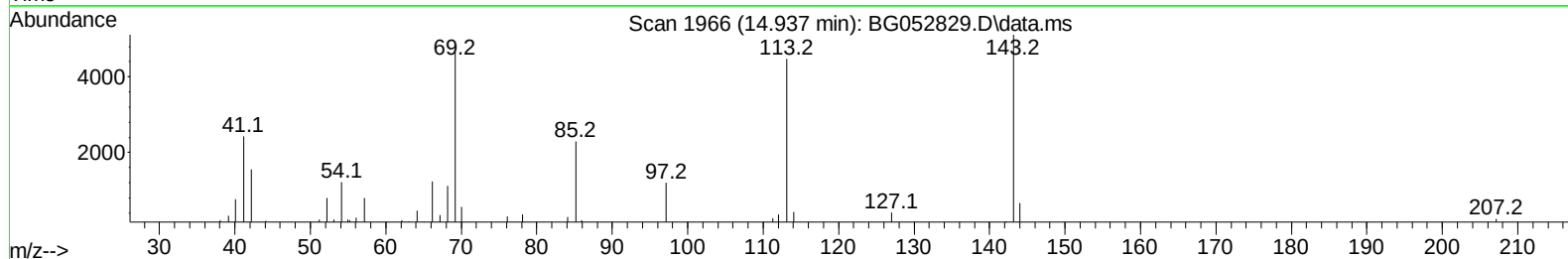
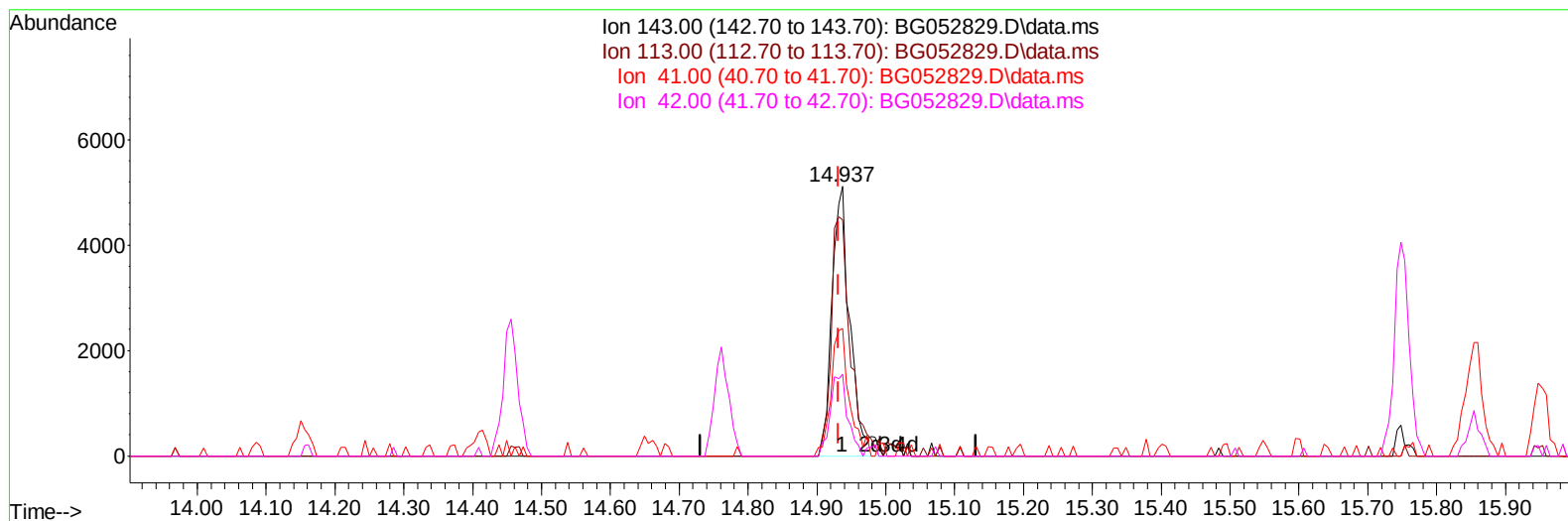
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TIC: BG052829.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.937min (+ 0.006) 8.42 ng/ul m

response 9463

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	87.68
41.00	47.20	47.35
42.00	29.40	30.28

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.146	152	22872	20.000	ng/ul	0.00
20) Naphthalene-d8	10.954	136	104714	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.761	164	87040	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.504	188	219850	20.000	ng/ul	0.00
79) Chrysene-d12	21.793	240	238869	20.000	ng/ul	0.00
88) Perylene-d12	25.106	264	236981	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.547	96	2879	4.446	ng/uL	0.00
4) Pyridine-d5	3.976	84	11810	7.306	ng/ul	0.00
7) Phenol-d5	7.289	99	16470	8.040	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.459	67	47156	37.690	ng/ul	0.00
11) 2-Chlorophenol-d4	7.671	132	38603	27.752	ng/ul	0.00
15) 4-Methylphenol-d8	8.834	113	28454	19.175	ng/ul	0.00
21) Nitrobenzene-d5	9.304	128	27998	37.144	ng/ul	0.00
24) 2-Nitrophenol-d4	10.026	143	31145	39.174	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.567	165	54113	31.286	ng/ul	0.00
31) 4-Chloroaniline-d4	11.078	131	69707	30.493	ng/ul	0.00
46) Dimethylphthalate-d6	14.156	166	248632	37.183	ng/ul	0.00
49) Acenaphthylene-d8	14.456	160	286219	35.225	ng/ul	0.00
54) 4-Nitrophenol-d4	14.937	143	9463m	8.417	ng/ul	0.00
60) Fluorene-d10	15.748	176	215001	36.619	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.854	200	43519	38.897	ng/ul	0.00
73) Anthracene-d10	17.604	188	391663	38.258	ng/ul	0.00
81) Pyrene-d10	19.884	212	486035	36.372	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.883	264	480411	39.234	ng/ul	0.00

Target Compounds Qvalue

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

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