

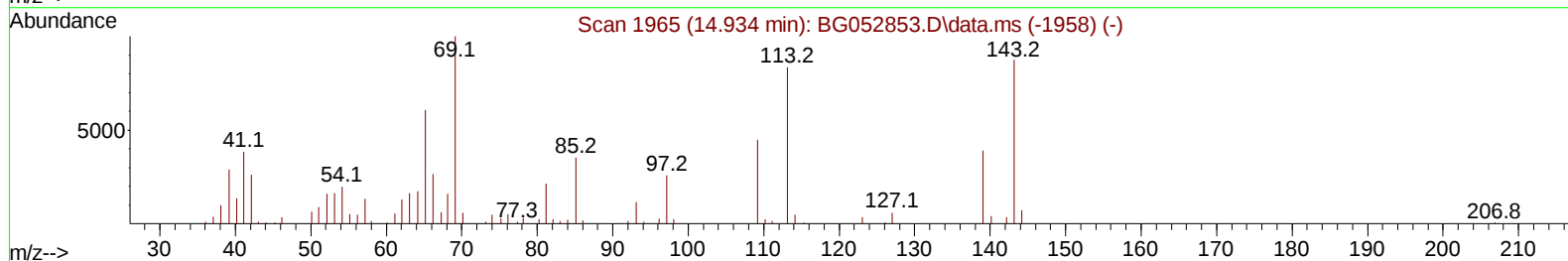
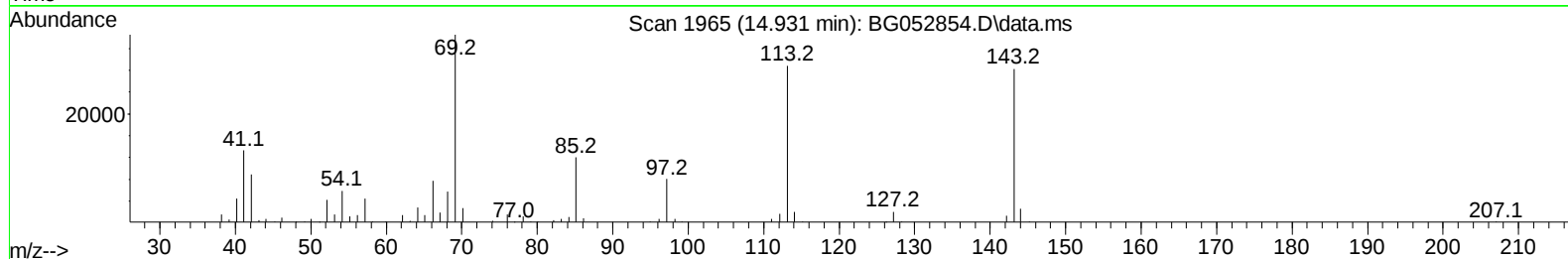
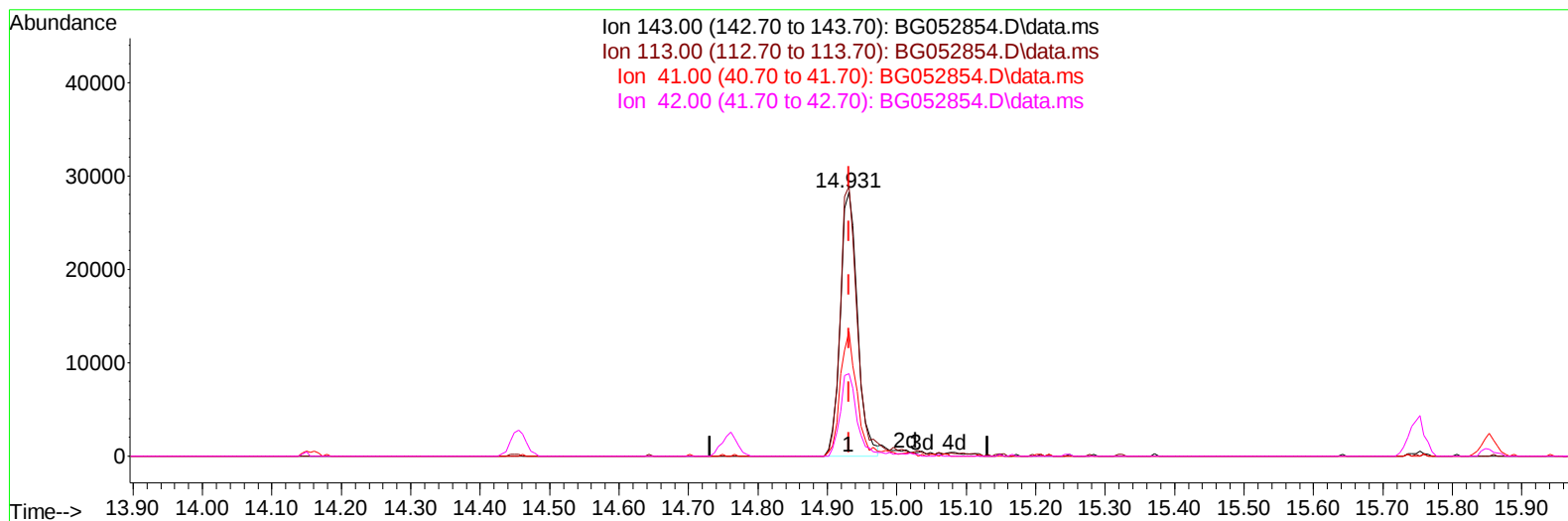
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032522\
Data File : BG052854.D
Acq On : 26 Mar 2022 8:17
Operator : CG/JU
Sample : PB143577BL
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK577

Manual IntegrationsAPPROVED

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

Quant Time: Mar 28 00:57:23 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



TIC: BG052854.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (-0.000) 33.25 ng/u1

response 48305

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	102.12
41.00	47.20	47.07
42.00	29.40	31.26

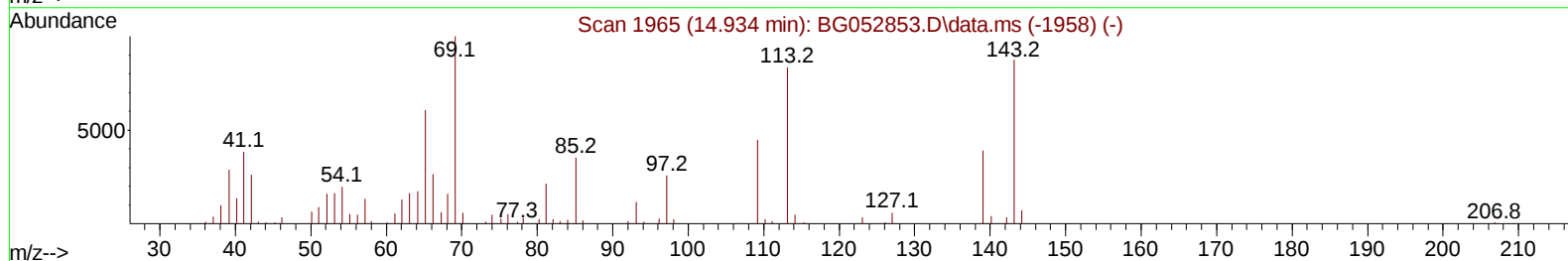
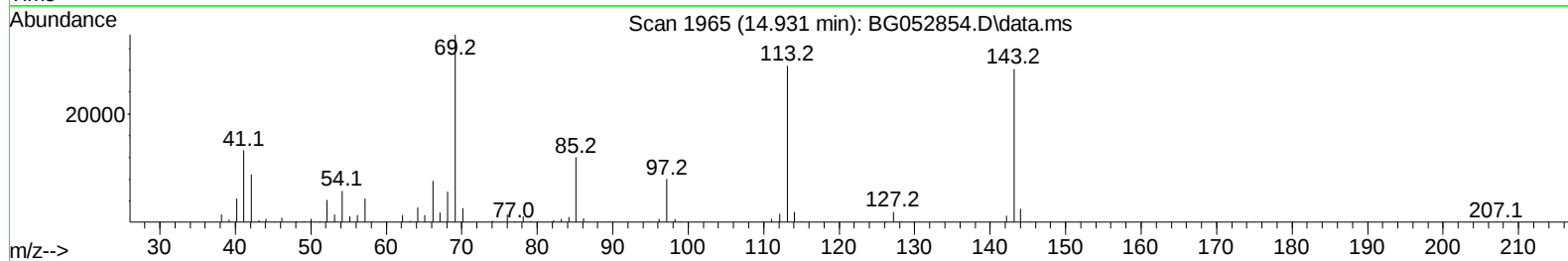
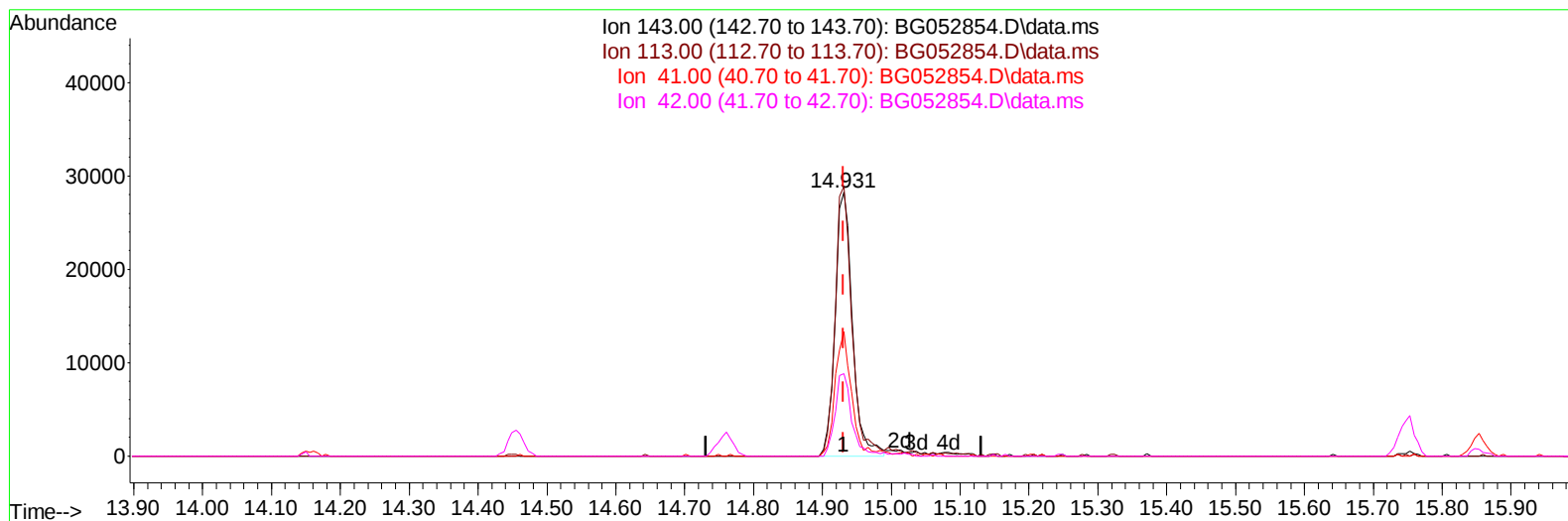
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TIC: BG052854.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.931min (-0.000) 33.89 ng/ul m

response 49240

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.70	102.12
41.00	47.20	47.07
42.00	29.40	31.26

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.140	152	28907	20.000	ng/ul	0.00
20) Naphthalene-d8	10.954	136	133715	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.760	164	112474	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.504	188	284673	20.000	ng/ul	0.00
79) Chrysene-d12	21.792	240	307482	20.000	ng/ul	0.00
88) Perylene-d12	25.105	264	299472	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.558	96	4572	5.587	ng/uL	0.00
4) Pyridine-d5	3.969	84	62439	30.562	ng/ul	0.00
7) Phenol-d5	7.282	99	88428	34.154	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.453	67	54584	34.519	ng/ul	0.00
11) 2-Chlorophenol-d4	7.670	132	58074	33.034	ng/ul	0.00
15) 4-Methylphenol-d8	8.833	113	68533	36.541	ng/ul	0.00
21) Nitrobenzene-d5	9.297	128	33651	34.961	ng/ul	0.00
24) 2-Nitrophenol-d4	10.026	143	37407	36.846	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.560	165	68747	31.126	ng/ul	0.00
31) 4-Chloroaniline-d4	11.077	131	100074	34.283	ng/ul	0.00
46) Dimethylphthalate-d6	14.155	166	289241	33.474	ng/ul	0.00
49) Acenaphthylene-d8	14.455	160	320060	30.483	ng/ul	0.00
54) 4-Nitrophenol-d4	14.931	143	49240m	33.894	ng/ul	0.00
60) Fluorene-d10	15.753	176	245507	32.359	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.853	200	46384	32.018	ng/ul	0.00
73) Anthracene-d10	17.604	188	434832	32.802	ng/ul	0.00
81) Pyrene-d10	19.883	212	543318	31.586	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.882	264	522046	33.737	ng/ul	0.00

Target Compounds Qvalue

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

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