

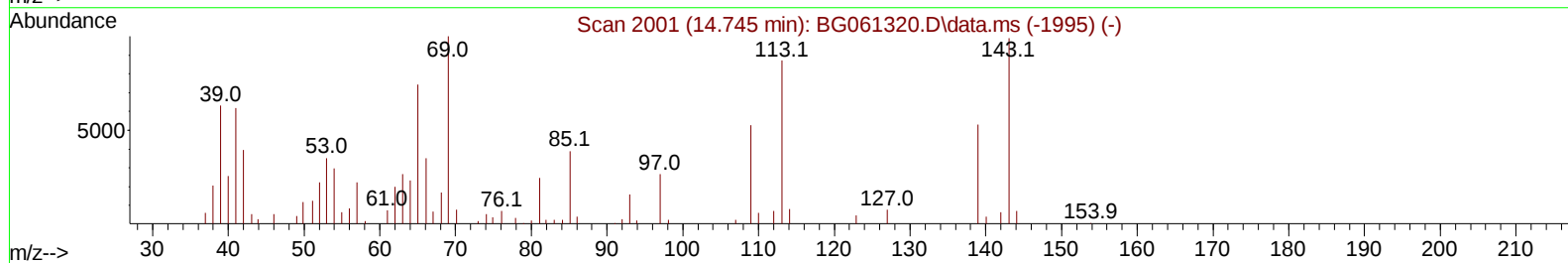
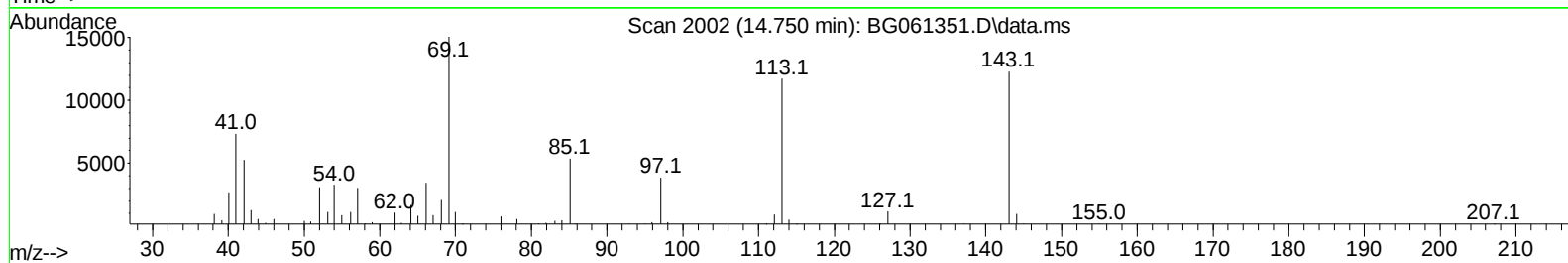
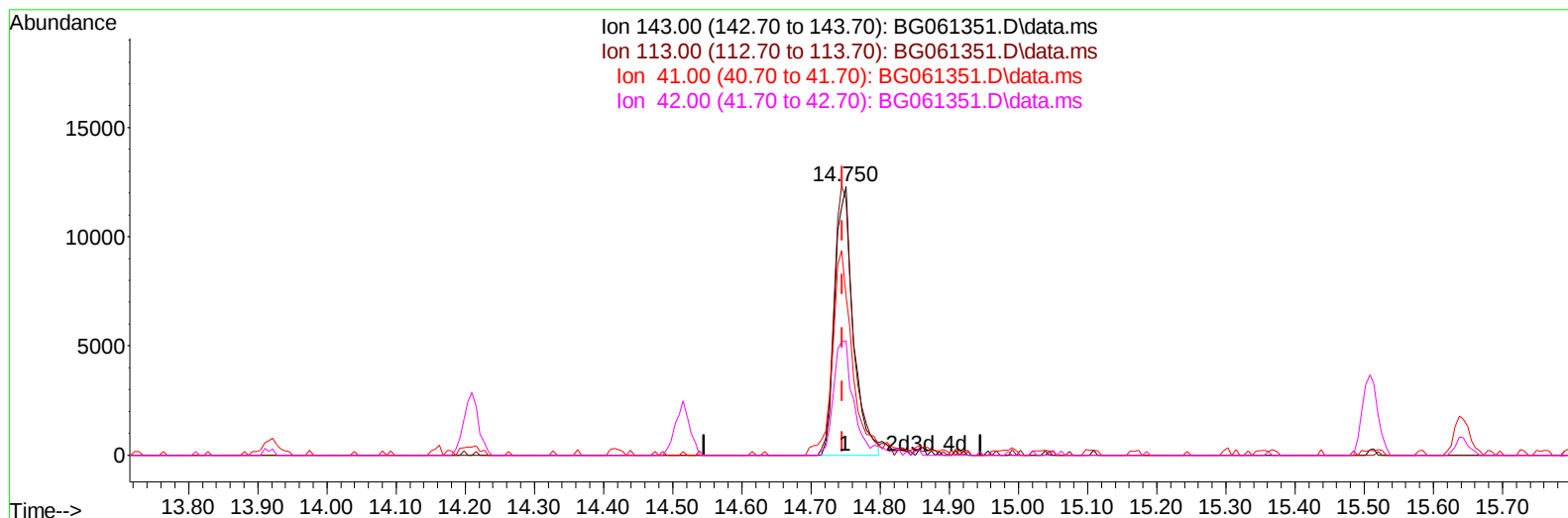
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052924\
Data File : BG061351.D
Acq On : 30 May 2024 7:02
Operator : MA/JU
Sample : P2571-18
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH6C2

Manual IntegrationsAPPROVED

Quant Time: May 30 07:38:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 21 15:23:39 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/30/2024
Supervised By :mohammad ahmed 06/01/2024



TIC: BG061351.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.750min (+ 0.005) 27.34 ng/u1

response 22746

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	122.20	95.36#
41.00	77.10	59.45#
42.00	49.50	42.62

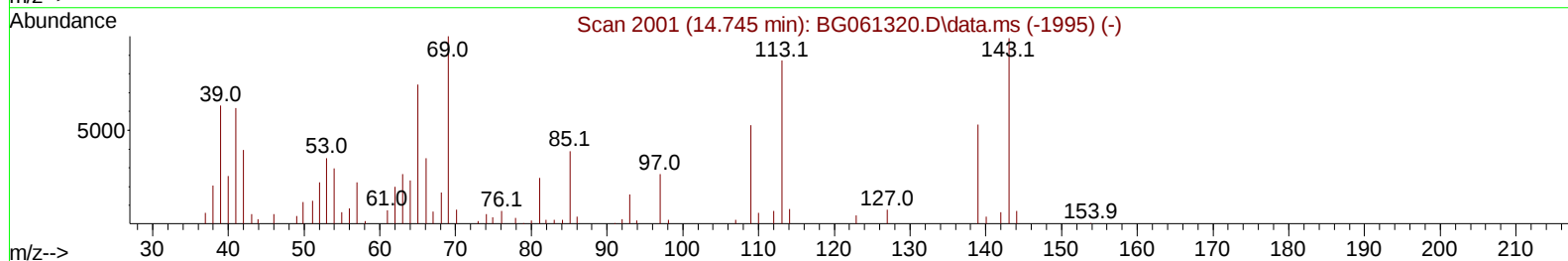
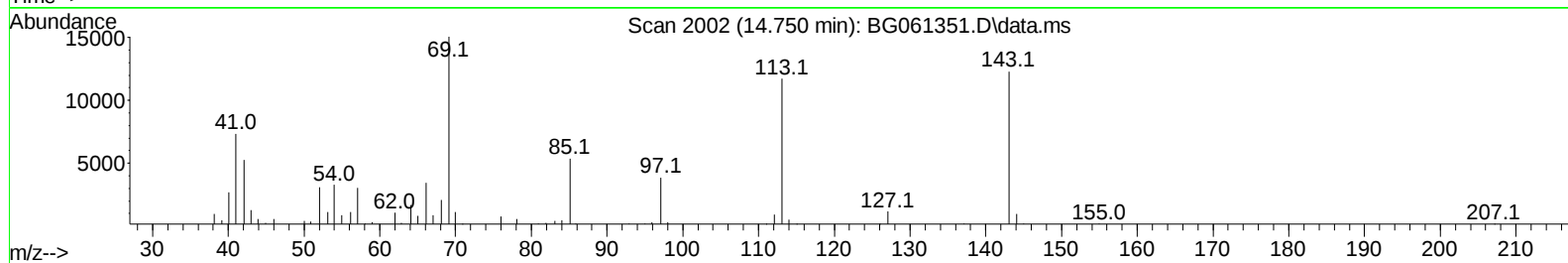
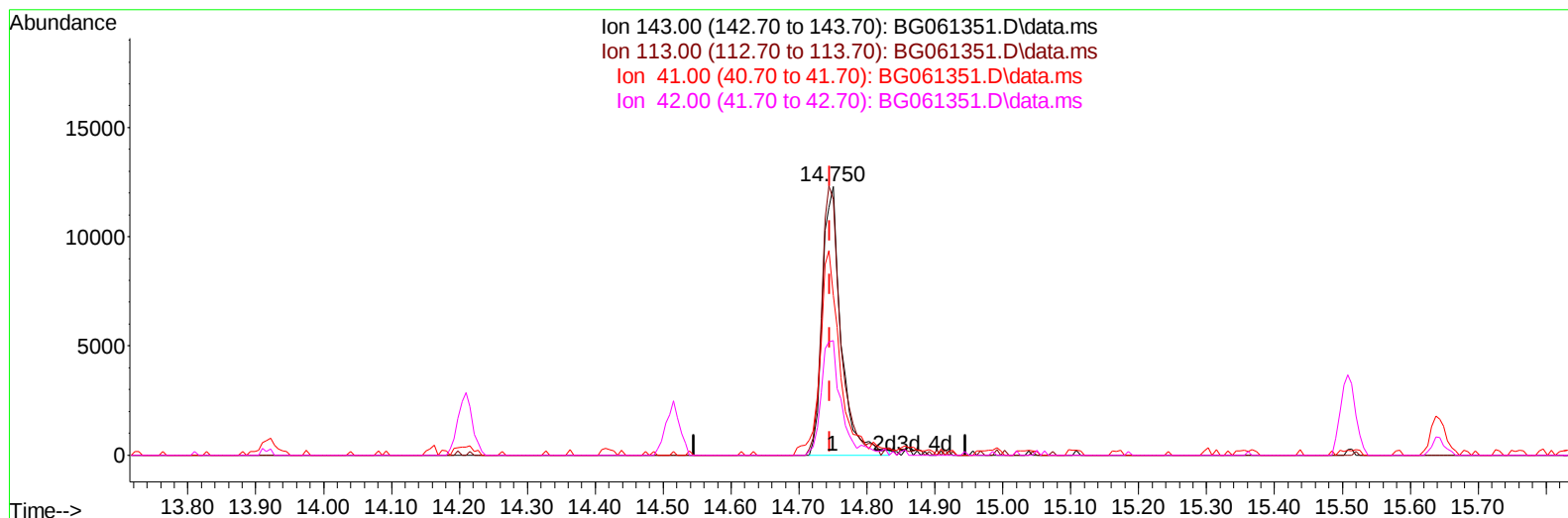
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TIC: BG061351.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.750min (+ 0.005) 28.23 ng/ul m

response 23485

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	122.20	95.36#
41.00	77.10	59.45#
42.00	49.50	42.62

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.882	152	27466	20.000	ng/ul	0.00
20) Naphthalene-d8	10.673	136	116977	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.515	164	86770	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.265	188	196416	20.000	ng/ul	0.00
79) Chrysene-d12	21.513	240	208524	20.000	ng/ul	0.00
88) Perylene-d12	24.597	264	233088	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	2439	4.970	ng/uL	0.00
4) Pyridine-d5	3.745	84	41120	23.551	ng/ul	0.00
7) Phenol-d5	7.065	99	65534	29.055	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.212	67	40224	28.373	ng/ul	0.00
11) 2-Chlorophenol-d4	7.418	132	50656	30.346	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	55854	29.041	ng/ul	0.00
21) Nitrobenzene-d5	9.039	128	27715	30.773	ng/ul	0.00
24) 2-Nitrophenol-d4	9.762	143	30916	29.330	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.308	165	63335	30.821	ng/ul	0.00
31) 4-Chloroaniline-d4	10.814	131	74156	27.584	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	208120	30.926	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	234214	33.476	ng/ul	0.00
54) 4-Nitrophenol-d4	14.750	143	23485m	28.227	ng/ul	0.00
60) Fluorene-d10	15.508	176	187167	32.214	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.643	200	21135	17.791	ng/ul	0.00
73) Anthracene-d10	17.365	188	288365	33.301	ng/ul	0.00
81) Pyrene-d10	19.656	212	354592	33.118	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.392	264	387911	34.617	ng/ul	0.00
Target Compounds						
					Qvalue	
80) Fluoranthene	19.321	202	34812	2.710	ng/ul	99
82) Pyrene	19.686	202	36662	2.750	ng/ul	98
85) Benzo(a)anthracene	21.495	228	23472	1.631	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.407	149	10030	1.479	ng/ul#	93
87) Chrysene	21.560	228	23291	1.758	ng/ul	97
90) Benzo(b)fluoranthene	23.622	252	28043	2.021	ng/ul#	97
93) Benzo(a)pyrene	24.451	252	18800	1.427	ng/ul#	92

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

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