

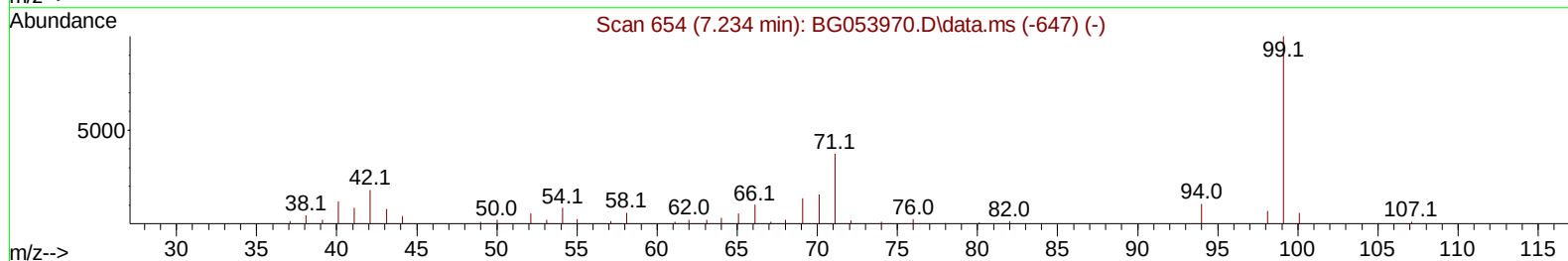
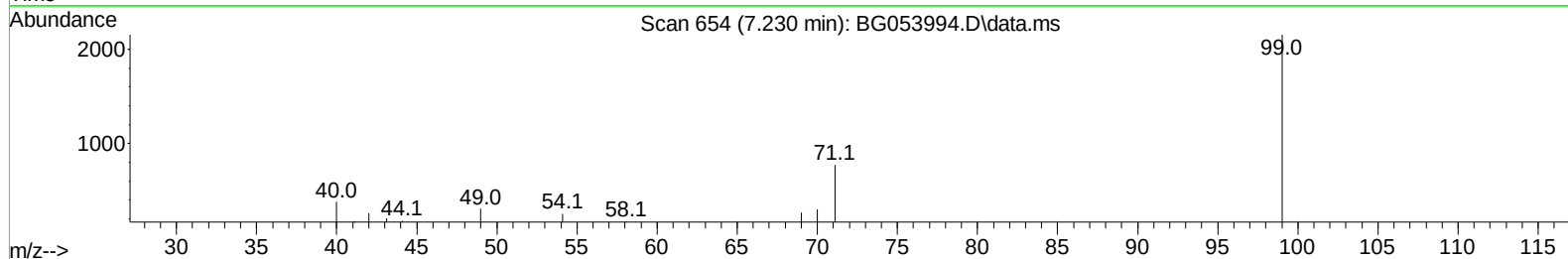
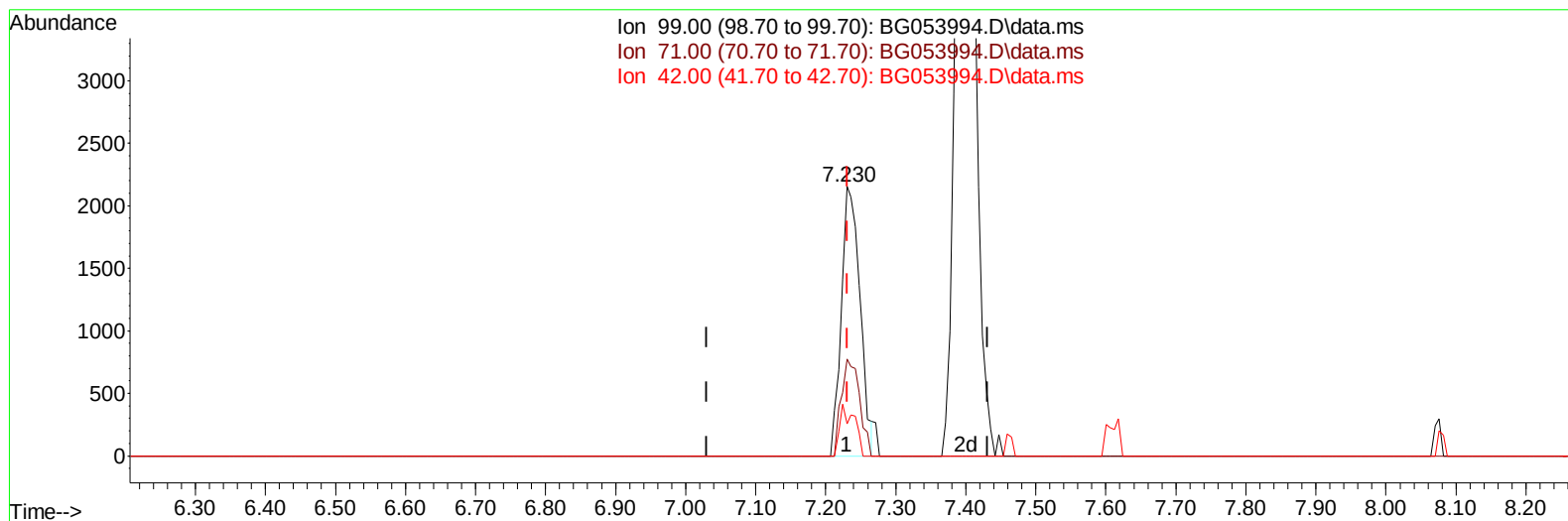
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062122\
 Data File : BG053994.D
 Acq On : 22 Jun 2022 1:51
 Operator : CG/JU
 Sample : N3400-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AD4

Manual IntegrationsAPPROVED

Quant Time: Jun 22 03:10:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 20 15:29:13 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/22/2022
 Supervised By :mohammad ahmed 06/28/2022



TIC: BG053994.D\data.ms

(7) Phenol-d5 (S)

7.230min (-0.001) 2.42 ng/u1

response 4016

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	39.00	35.96
42.00	22.60	11.97#
0.00	0.00	0.00

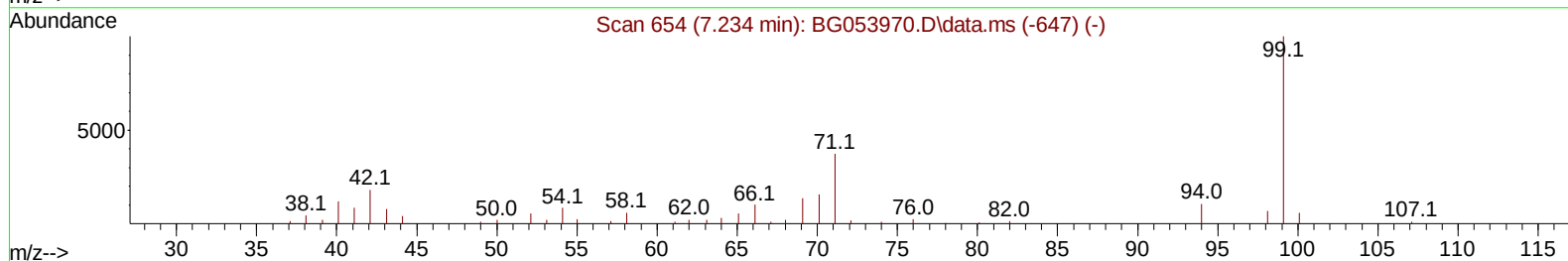
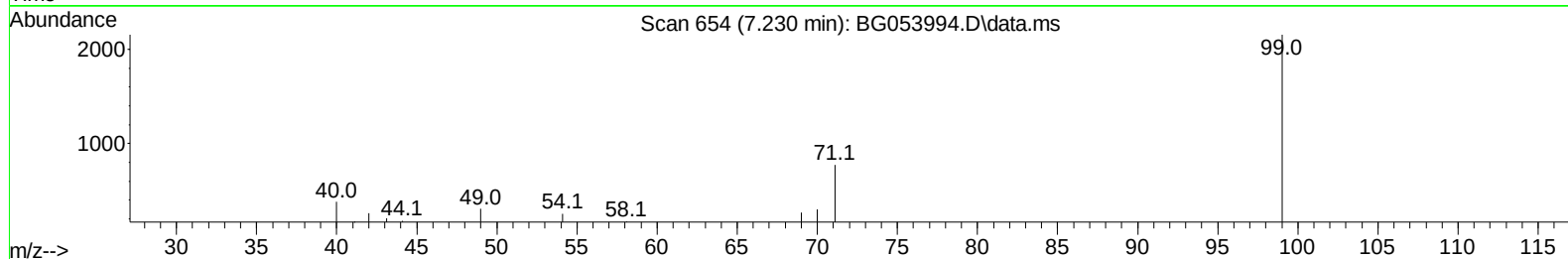
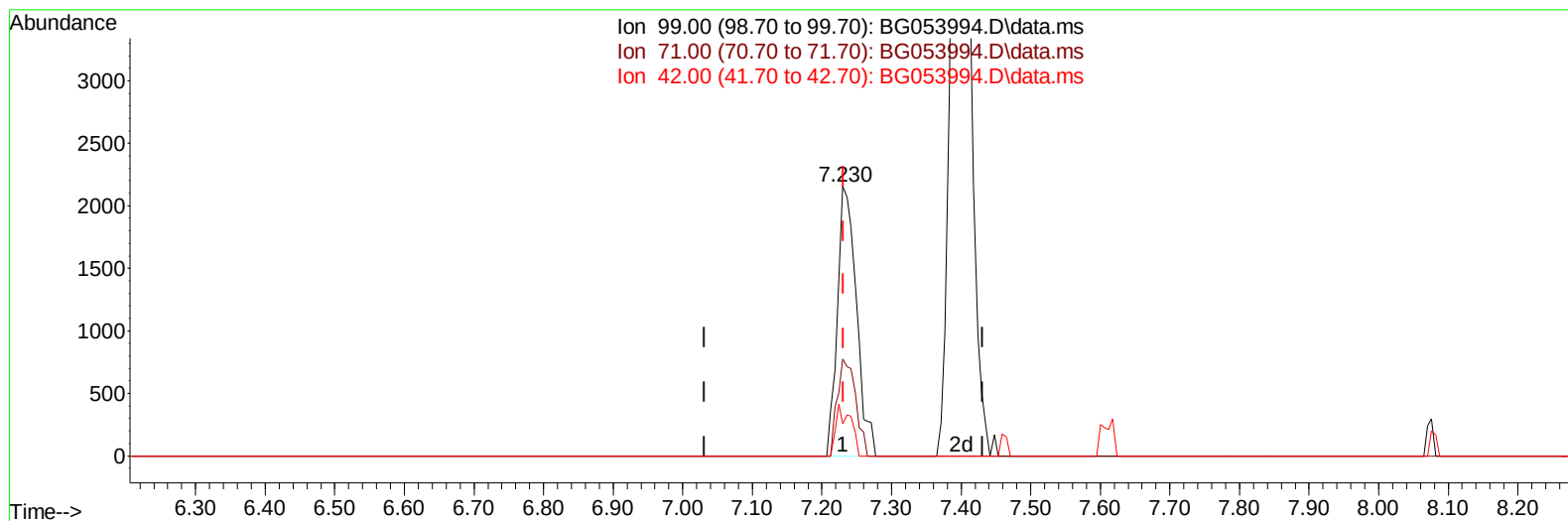
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TIC: BG053994.D\data.ms

(7) Phenol-d5 (S)

7.230min (-0.001) 2.48 ng/ul m

response 4111

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	39.00	35.96
42.00	22.60	11.97#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.082	152	23175	20.000	ng/ul	0.00
20) Naphthalene-d8	10.885	136	92191	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	69158	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.442	188	187275	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.719	240	211441	20.000	ng/ul	# 0.00
88) Perylene-d12	24.980	264	214064	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.505	96	1939	3.981	ng/uL	0.00
4) Pyridine-d5	3.916	84	14125	10.662	ng/ul	0.00
7) Phenol-d5	7.230	99	4111m	2.476	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.395	67	12990	12.970	ng/ul	0.00
11) 2-Chlorophenol-d4	7.612	132	5969	4.508	ng/ul	0.00
15) 4-Methylphenol-d8	8.775	113	5605	3.959	ng/ul	0.00
21) Nitrobenzene-d5	9.239	128	8366	12.226	ng/ul	0.00
24) 2-Nitrophenol-d4	9.968	143	2914	4.008	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.509	165	7200	4.476	ng/ul	0.00
31) 4-Chloroaniline-d4	11.014	131	41539	20.440	ng/ul	0.00
46) Dimethylphthalate-d6	14.098	166	74312	13.659	ng/ul	0.00
49) Acenaphthylene-d8	14.392	160	81296	13.742	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	840	1.128	ng/ul	0.02
60) Fluorene-d10	15.691	176	67079	13.726	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.802	200	1718	1.682	ng/ul	0.00
73) Anthracene-d10	17.541	188	123503	14.603	ng/ul	0.00
81) Pyrene-d10	19.827	212	169731	15.396	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.751	264	166604	15.419	ng/ul	0.00

Target Compounds Qvalue

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

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