

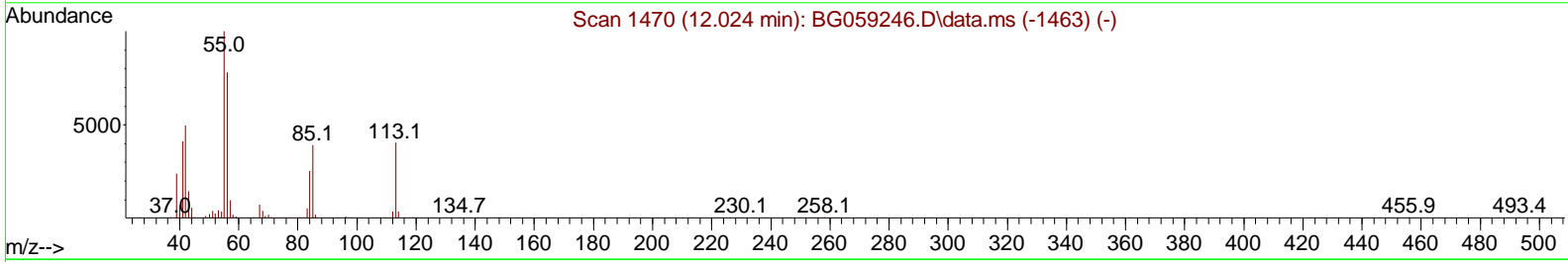
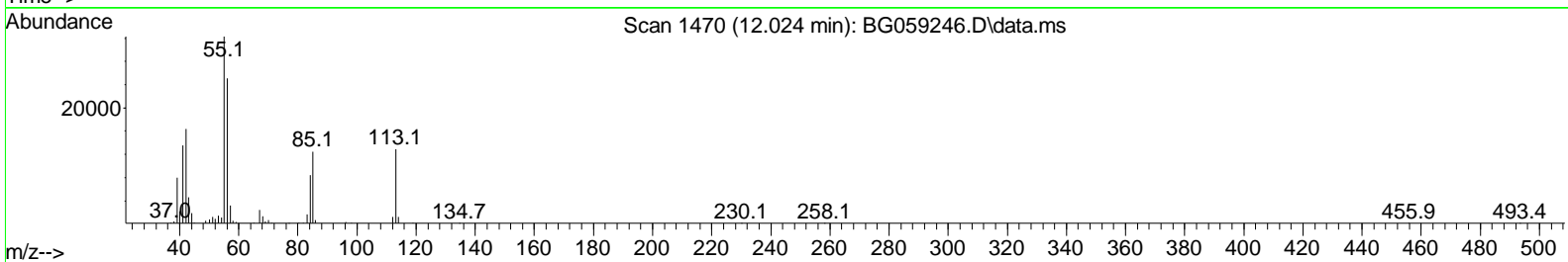
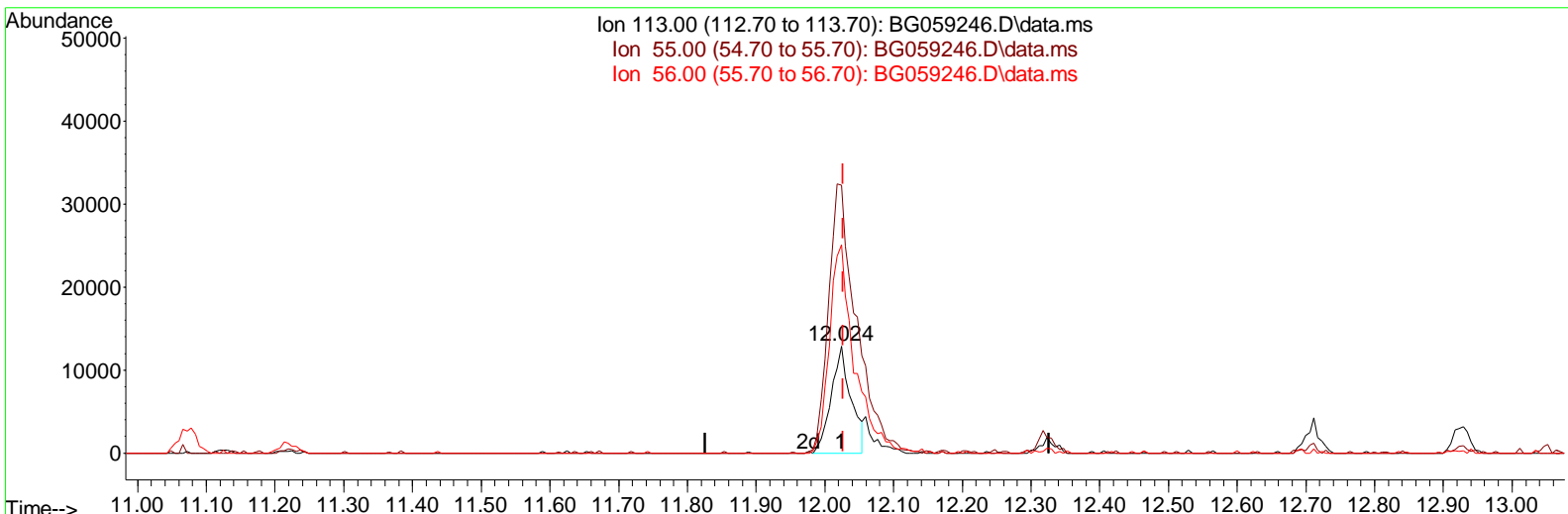
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG100223\  
 Data File : BG059246.D  
 Acq On : 2 Oct 2023 12:01  
 Operator : MA/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

**Instrument :**  
 BNA\_G  
**LabSampleId :**  
 SSTDCCC020

**Manual Integrations APPROVED**

Reviewed By :Yogesh Patel 10/03/2023  
 Supervised By :mohammad ahmed 10/04/2023

Quant Time: Oct 02 23:53:43 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



TIC: BG059246.D\data.ms

(34) Caprolactam

12.024min (-0.003) 17.75 ng/ul

response 25724

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	264.10	250.31
56.00	209.90	194.65
0.00	0.00	0.00

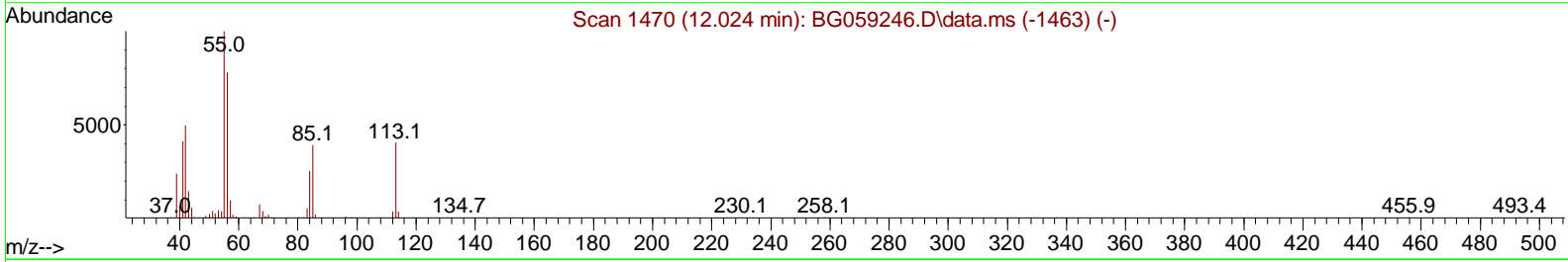
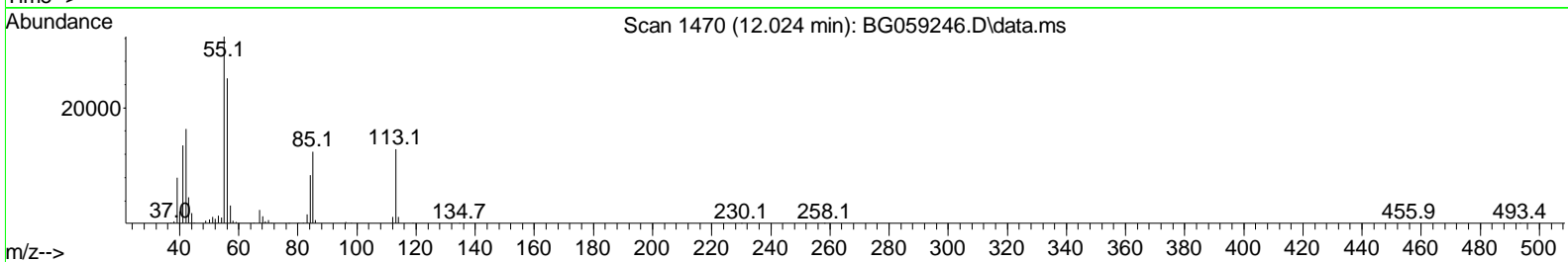
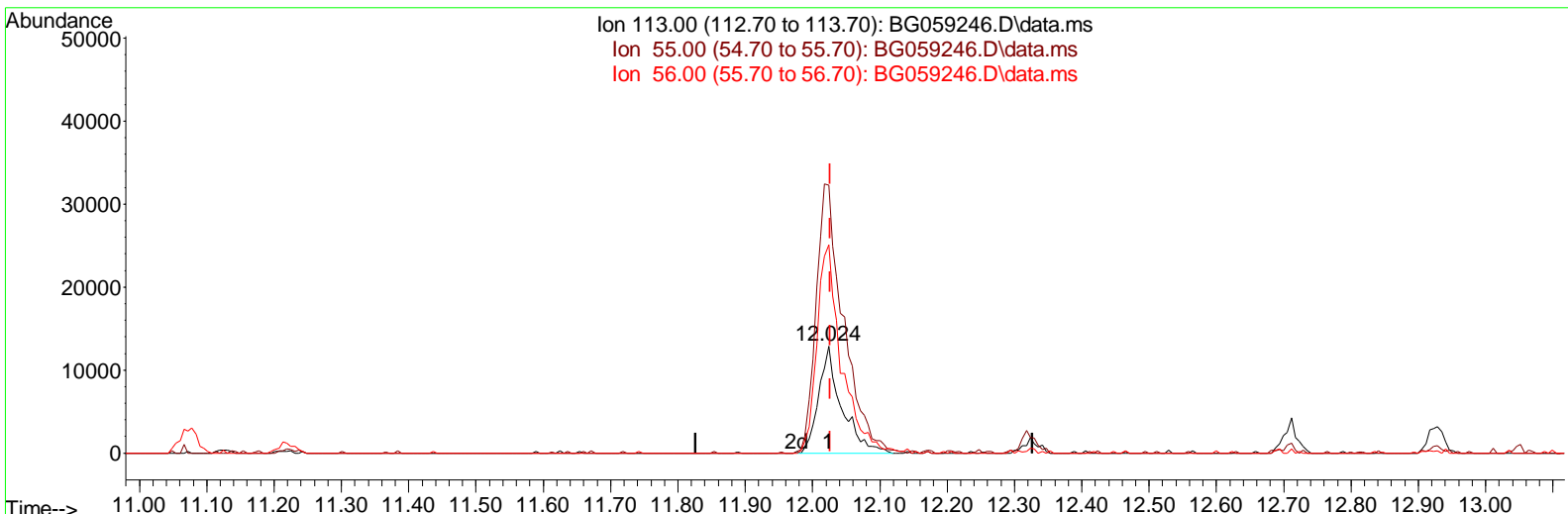
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 Supervised By :mohammad ahmed 10/04/2023



TIC: BG059246.D\data.ms

(34) Caprolactam

12.024min (-0.003) 20.96 ng/ul m

response 30378

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	264.10	250.31
56.00	209.90	194.65
0.00	0.00	0.00

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 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

**Instrument :**  
 BNA\_G  
**LabSampled :**  
 SSTDCCC020

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 10/03/2023  
 Supervised By :mohammad ahmed 10/04/2023

Quant Time: Oct 02 23:57:29 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG092723.MA.M  
 Quant Title : SVOA CALI BRATI ON  
 QLast Update : Thu Sep 28 04:33:25 2023  
 Response via : Initial Calibration

Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
<b>Internal Standards</b>						
1) 1,4-Di chlorobenzene-d4	8.228	152	57426	20.000	ng/ul	0.00
20) Naphthalene-d8	11.072	136	296164	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.868	164	201606	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.611	188	457983	20.000	ng/ul	0.00
79) Chrysene-d12	21.918	240	408563	20.000	ng/ul	0.00
88) Perylene-d12	25.396	264	482605	20.000	ng/ul	-0.01
<b>System Monitoring Compounds</b>						
3) 1,4-Dioxane-d8	3.528	96	12854	8.448	ng/uL	0.00
4) Pyridine-d5	3.957	84	101901	22.953	ng/ul	0.00
7) Phenol-d5	7.359	99	131332	22.847	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.553	67	79074	21.534	ng/ul	0.00
11) 2-Chlorophenol-d4	7.747	132	92602	23.965	ng/ul	0.00
15) 4-Methylphenol-d8	8.928	113	99966	22.498	ng/ul	0.00
21) Nitrobenzene-d5	9.427	128	46615	22.305	ng/ul	0.00
24) 2-Nitrophenol-d4	10.144	143	50933	24.825	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.678	165	95756	22.720	ng/ul	0.00
31) 4-Chloroaniline-d4	11.219	131	146311	21.456	ng/ul	0.00
46) Dimethylphthalate-d6	14.262	166	305913	20.987	ng/ul	0.00
49) Acenaphthylene-d8	14.568	160	363939	19.892	ng/ul	0.00
54) 4-Nitrophenol-d4	15.056	143	49975	19.634	ng/ul	0.00
60) Fluorene-d10	15.855	176	279153	20.122	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.966	200	45561	19.354	ng/ul	0.00
73) Anthracene-d10	17.711	188	420859	20.168	ng/ul	0.00
81) Pyrene-d10	19.985	212	471364	20.602	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.156	264	465746	20.083	ng/ul	0.00
<b>Target Compounds</b>						
2) 1,4-Dioxane	3.563	88	13938	7.797	ng/uL	98
5) Pyridine	3.980	79	100351	22.259	ng/ul	90
6) Benzaldehyde	7.376	77	62026	24.571	ng/ul	90
8) Phenol	7.382	94	129511	22.649	ng/ul	93
10) Bis(2-Chloroethyl)ether	7.647	93	103221	22.010	ng/ul	96
12) 2-Chlorophenol	7.782	128	93949	23.060	ng/ul	97
13) 2-Methylphenol	8.657	108	91907	21.273	ng/ul	90
14) 2,2'-oxybis(1-Chloropr...	8.740	45	221720	21.386	ng/ul	98
16) Acetophenone	9.074	105	156049	22.676	ng/ul	97
17) N-Nitrosodipropylamine	9.039	70	77789	22.285	ng/ul	99
18) 4-Methylphenol	8.992	108	106191	23.199	ng/ul	88
19) Hexachloroethane	9.315	117	36983	23.042	ng/ul	98
22) Nitrobenzene	9.474	77	109498	21.225	ng/ul	94
23) Isophorone	9.979	82	234411	21.130	ng/ul	99
25) 2-Nitrophenol	10.185	139	51768	22.901	ng/ul	92
26) 2,4-Dimethylphenol	10.208	107	107341	22.207	ng/ul	96
27) Bis(2-Chloroethoxy)met...	10.461	93	149015	21.288	ng/ul #	95
29) 2,4-Dichlorophenol	10.702	162	97263	23.216	ng/ul	91
30) Naphthalene	11.125	128	339933	21.407	ng/ul	96
32) 4-Chloroaniline	11.242	127	142601	21.869	ng/ul	98
33) Hexachlorobutadiene	11.354	225	59483	21.214	ng/ul	99
34) Caprolactam	12.024	113	30378m	20.956	ng/ul	
35) 4-Chloro-3-methylphenol	12.324	107	104057	23.278	ng/ul	95

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**Instrument :**  
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**LabSampleId :**  
 SSTDCCC020

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 10/03/2023  
 Supervised By :mohammad ahmed 10/04/2023

Quant Time: Oct 02 23: 57: 29 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG092723.MA.M  
 Quant Title : SVOA CALI BRATI ON  
 QLast Update : Thu Sep 28 04: 33: 25 2023  
 Response via : Ini tial Cal i brati on

Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
36) 2-Methyl naphthal ene	12. 711	142	222593	21. 398	ng/ul	99
37) 1-Methyl naphthal ene	12. 929	142	231506	21. 927	ng/ul	93
39) 1, 2, 4, 5-Tetrachl oroben. . .	13. 052	216	115365	19. 114	ng/ul	99
40) Hexachl orocycl opentadi ene	13. 005	237	65774	17. 666	ng/ul #	94
41) 2, 4, 6-Tri chl orophenol	13. 299	196	76793	20. 781	ng/ul	89
42) 2, 4, 5-Tri chl orophenol	13. 369	196	82811	21. 356	ng/ul	95
43) 1, 1' -Bi phenyl	13. 704	154	315772	19. 394	ng/ul	96
44) 2-Chl oronaphthal ene	13. 757	162	241796	19. 724	ng/ul	98
45) 2-Ni troani li ne	13. 975	65	77983	22. 552	ng/ul	99
47) Di methyl phthal ate	14. 309	163	298289	19. 980	ng/ul	98
48) 2, 6-Di ni trotol uene	14. 456	165	57753	22. 744	ng/ul	88
50) Acenaphthyl ene	14. 597	152	389634	19. 952	ng/ul	98
51) 3-Ni troani li ne	14. 791	138	62381	22. 757	ng/ul	93
52) Acenaphthene	14. 932	153	269938	20. 138	ng/ul	91
53) 2, 4-Di ni trophenol	14. 985	184	19973	14. 150	ng/ul	87
55) 4-Ni trophenol	15. 073	109	36334	20. 236	ng/ul	89
56) Di benzofuran	15. 261	168	360698	20. 300	ng/ul	97
57) 2, 4-Di ni trotol uene	15. 232	165	82686	21. 992	ng/ul #	99
58) 2, 3, 4, 6-Tetrachl orophenol	15. 473	232	76682	21. 845	ng/ul	95
59) Di ethyl phthal ate	15. 655	149	296662	20. 554	ng/ul	94
61) Fl uorene	15. 908	166	300327	20. 130	ng/ul	97
62) 4-Chl orophenyl -phenyl e. . .	15. 890	204	153536	20. 644	ng/ul	98
63) 4-Ni troani li ne	15. 955	138	55052	21. 823	ng/ul	91
66) 4, 6-Di ni tro-2-methyl ph. . .	15. 978	198	49579	19. 974	ng/ul #	91
67) N-Ni trosodi phenyl ami ne	16. 113	169	247003	20. 233	ng/ul	98
68) 4-Bromophenyl -phenyl ether	16. 789	248	90125	20. 313	ng/ul	93
69) Hexachl orobenzene	16. 877	284	102242	20. 398	ng/ul	95
70) Atrazi ne	17. 047	200	96728	20. 837	ng/ul	93
71) Pentachl orophenol	17. 235	266	57505	19. 934	ng/ul	98
72) Phenanthrene	17. 658	178	513758	20. 441	ng/ul	99
74) Anthracene	17. 747	178	518352	19. 973	ng/ul	98
75) 1, 2, 3, 4-Tetrachl oroben. . .	13. 663	216	119820	19. 853	ng/uL	99
76) Pentachl orobenzene	15. 156	250	119443	20. 043	ng/uL	99
77) Carbazol e	18. 029	167	438587	20. 735	ng/ul	99
78) Di -n-butyl phthal ate	18. 528	149	522087	20. 822	ng/ul	99
80) Fl uoranthene	19. 650	202	578213	20. 514	ng/ul	97
82) Pyrene	20. 015	202	613951	20. 565	ng/ul	97
83) Butyl benzyl phthal ate	20. 872	149	224768	21. 349	ng/ul	97
84) 3, 3' -Di chl orobenzi di ne	21. 812	252	200974	21. 241	ng/ul	100
85) Benzo(a)anthracene	21. 895	228	585241	20. 349	ng/ul	100
86) Bi s(2-ethyl hexyl )phtha. . .	21. 724	149	337019	21. 081	ng/ul	100
87) Chrysene	21. 965	228	548539	20. 194	ng/ul	99
89) Di -n-octyl phthal ate	23. 017	149	569259	20. 281	ng/ul	100
90) Benzo(b)fl uoranthene	24. 268	252	588333	19. 904	ng/ul	97
91) Benzo(k)fl uoranthene	24. 345	252	609595	20. 354	ng/ul	98
93) Benzo(a)pyrene	25. 232	252	568977	20. 291	ng/ul	98
94) I ndeno(1, 2, 3-cd)pyrene	29. 404	276	636006	20. 037	ng/ul	99
95) Di benzo(a, h)anthracene	29. 492	278	510089	19. 761	ng/ul	98
96) Benzo(g, h, i )peryl ene	30. 690	276	505887	19. 601	ng/ul	97

(#) = qual i fier out of range (m) = manual i ntegrati on (+) = signal s summed

**Instrument :**

BNA\_G

**LabSampleId :**

SSTDCCC020

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 10/03/2023

Supervised By :mohammad ahmed 10/04/2023

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