

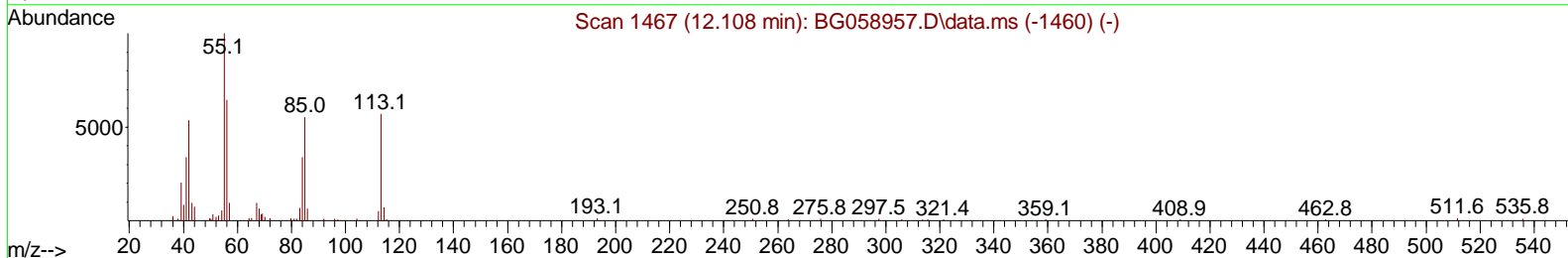
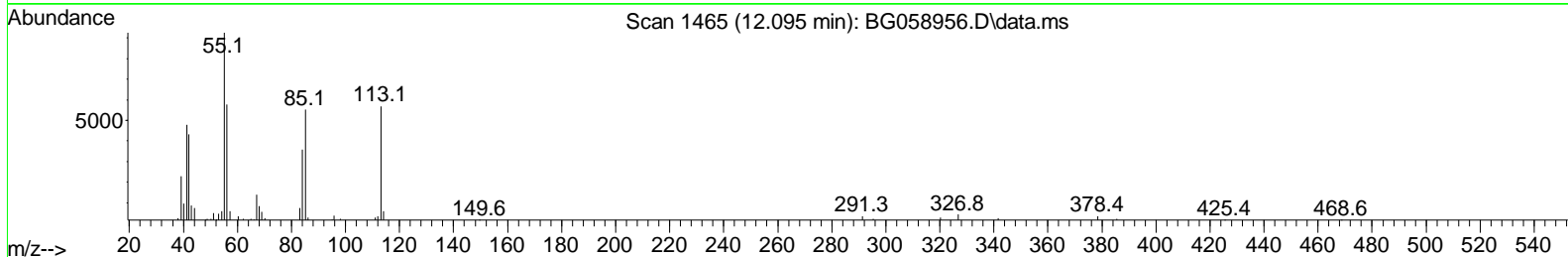
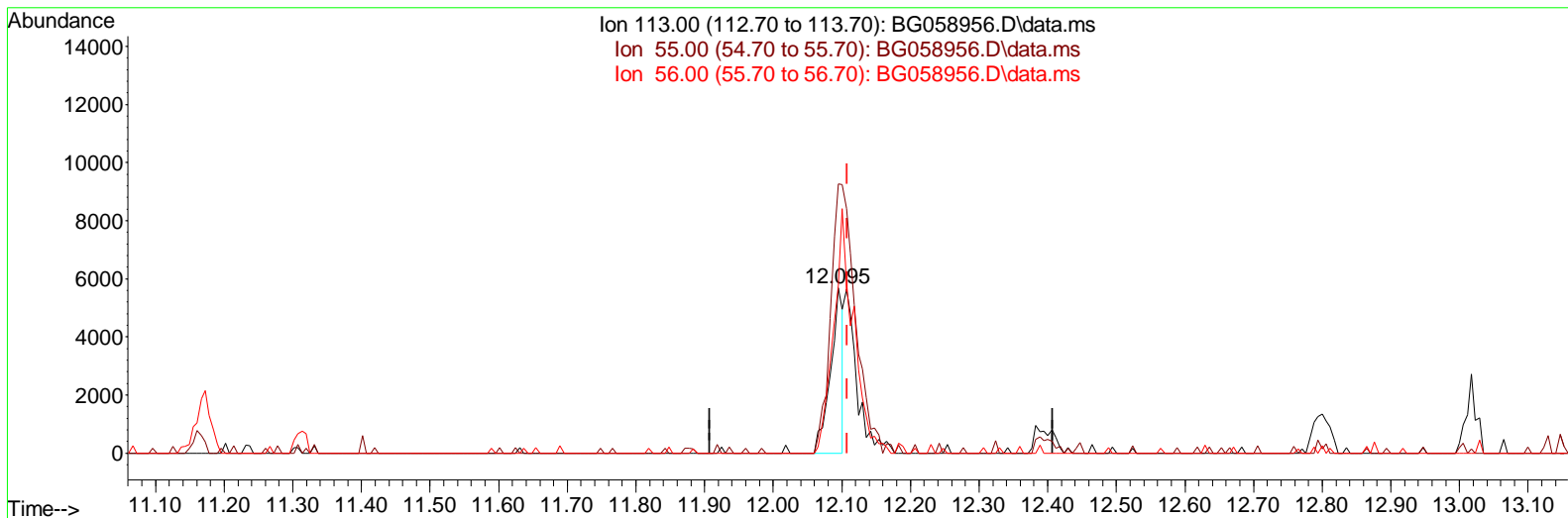
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
ClientSampleId :
 SSTD010408

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 09/11/2023
 Supervised By :mohammad ahmed 09/13/2023

Quant Time: Sep 10 23:05:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Sep 10 22:58:52 2023
 Response via : Initial Calibration



TIC: BG058956.D\data.ms

(34) Caprolactam

12.095min (-0.013) 5.29 ng/ul

response 7249

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	175.60	162.91
56.00	113.00	101.72
0.00	0.00	0.00

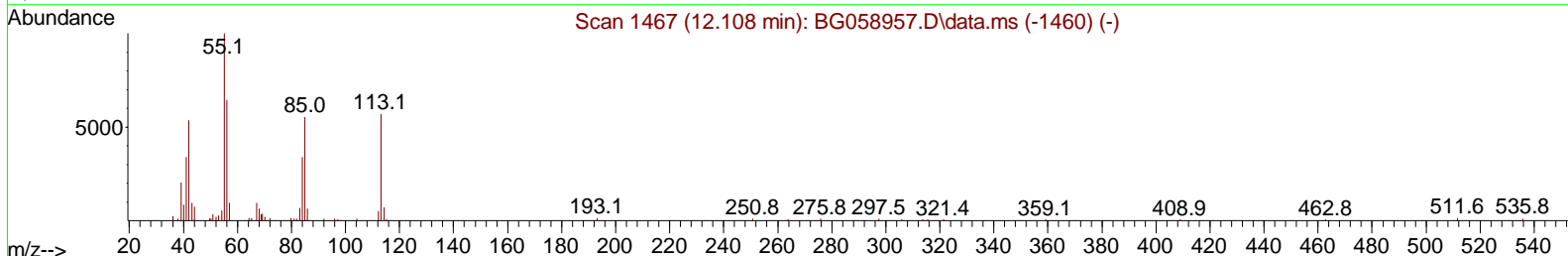
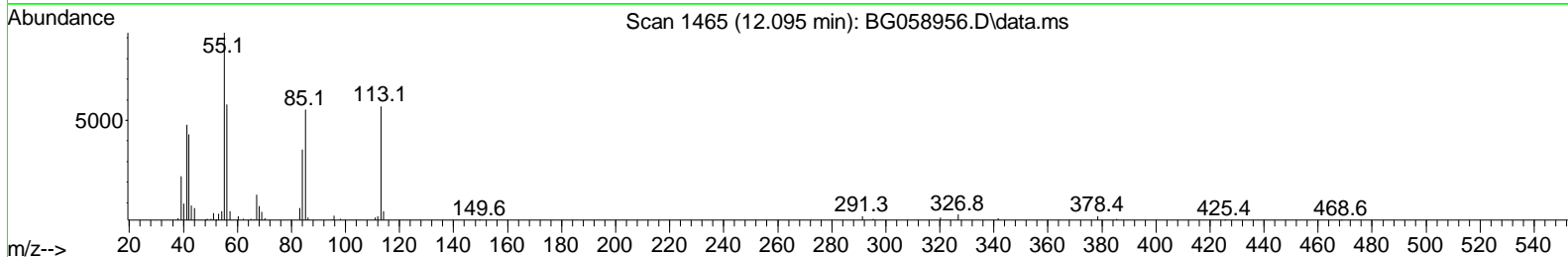
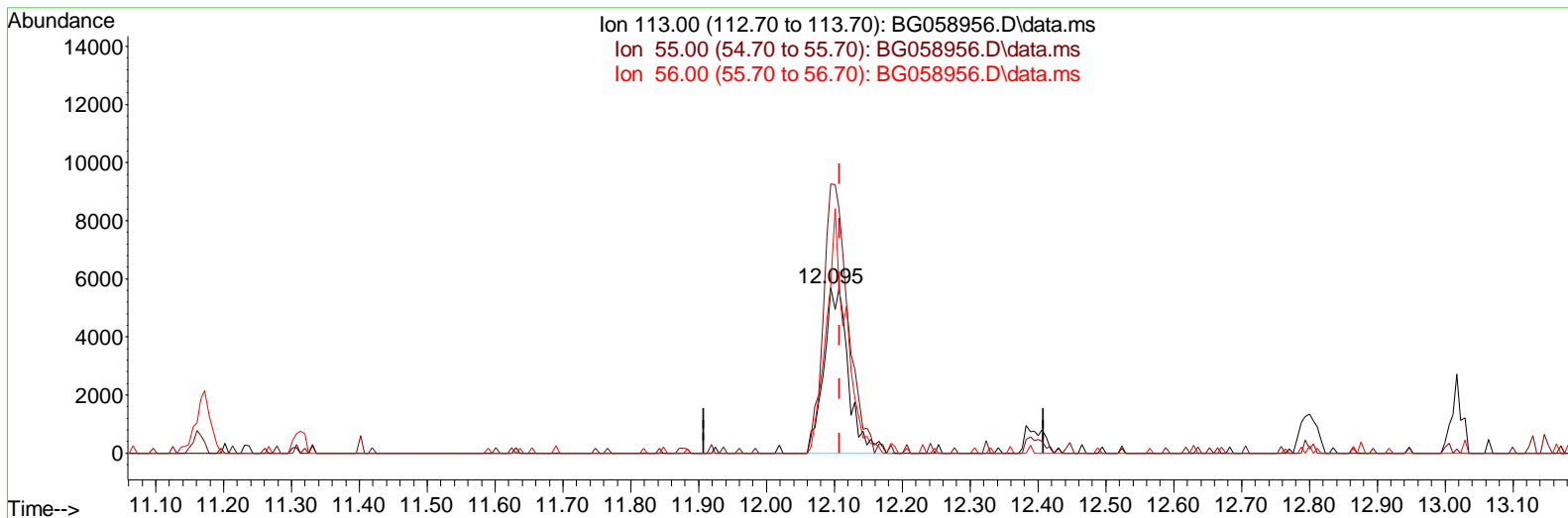
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
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Manual Integrations APPROVED

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 Supervised By : mohammad ahmed 09/13/2023

Quant Time: Sep 10 23:05:36 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Sep 10 22:58:52 2023
 Response via : Initial Calibration



TIC: BG058956.D\data.ms

(34) Caprolactam

12.095min (-0.013) 10.42 ng/ul m

response 14275

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	175.60	162.91
56.00	113.00	101.72
0.00	0.00	0.00

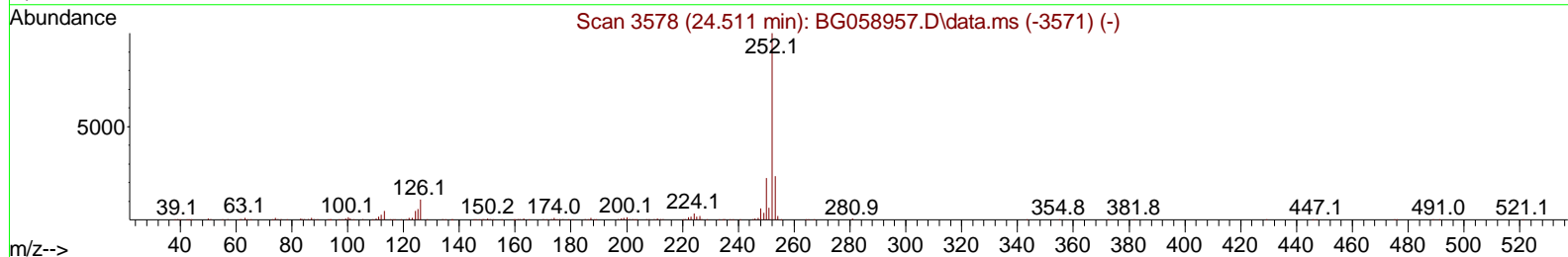
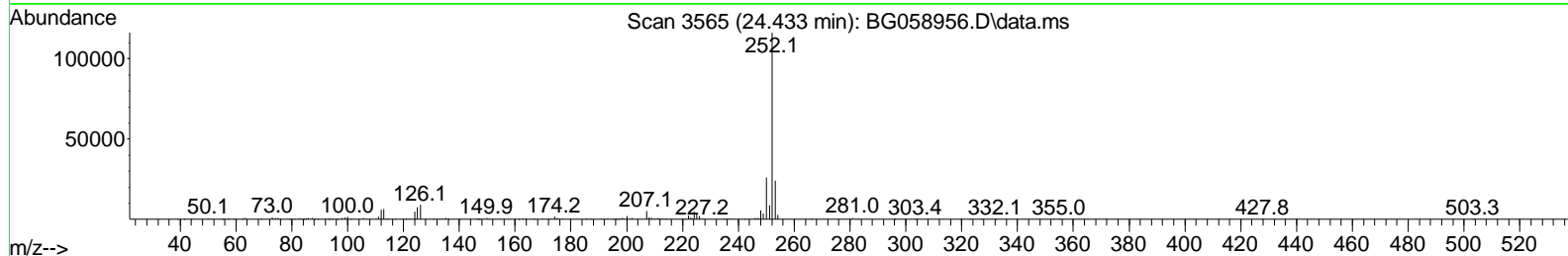
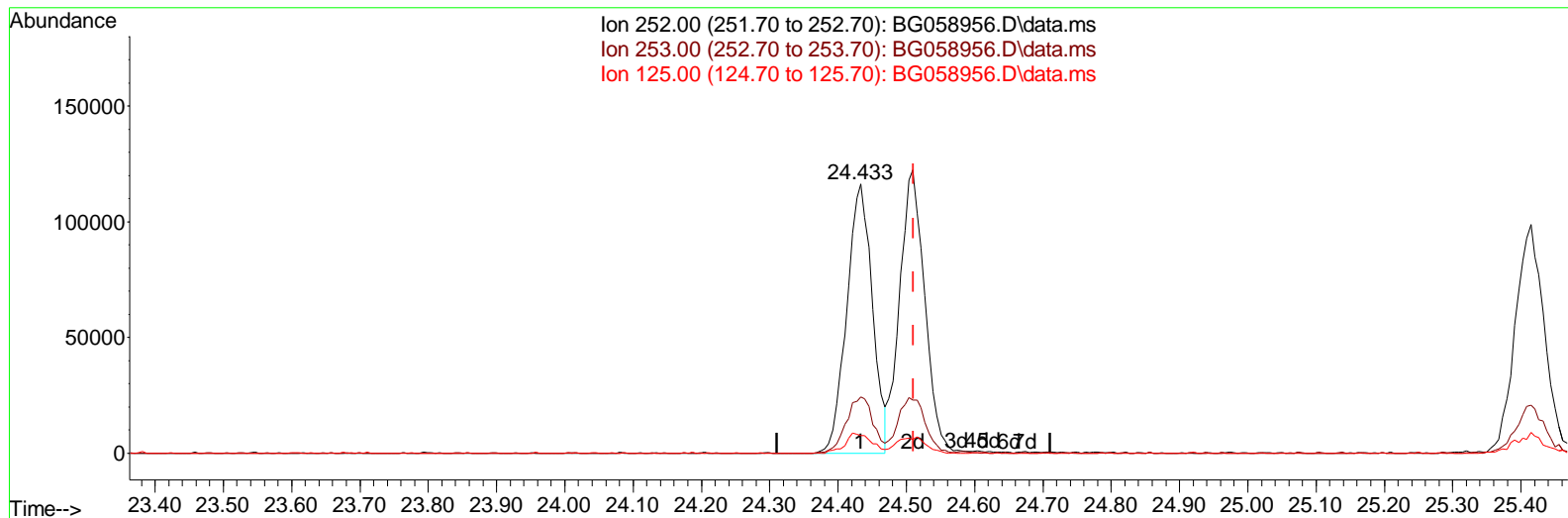
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
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 ALS Vial : 34 Sample Multiplier: 1

Instrument :
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Manual Integrations APPROVED

Quant Time: Sep 10 23:05:36 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Sep 10 22:58:52 2023
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Reviewed By :Yogesh Patel 09/11/2023
 Supervised By :mohammad ahmed 09/13/2023



TIC: BG058956.D\data.ms

(91) Benzo(k)fluoranthene

24.433min (-0.078) 9.33 ng/ul

response 302744

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	23.90	20.78
125.00	6.20	6.50
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
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 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

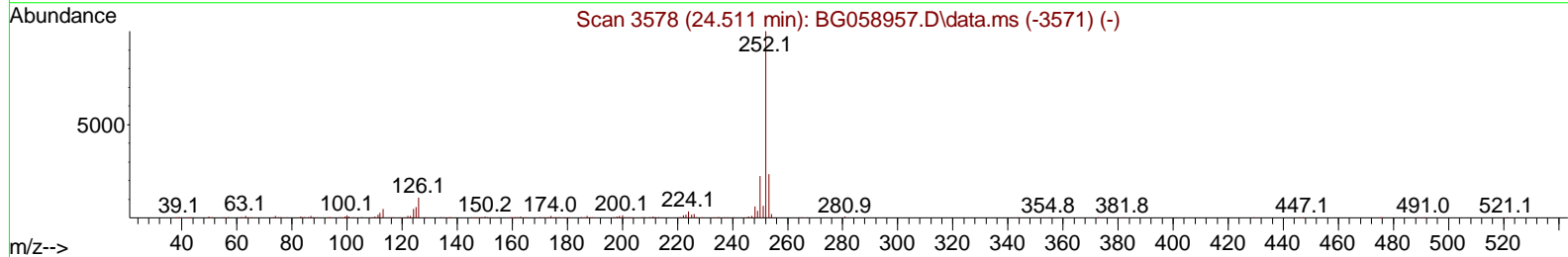
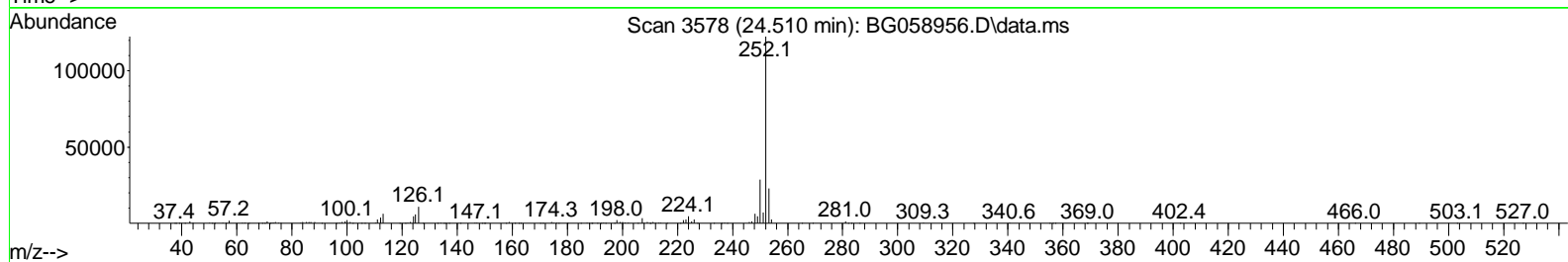
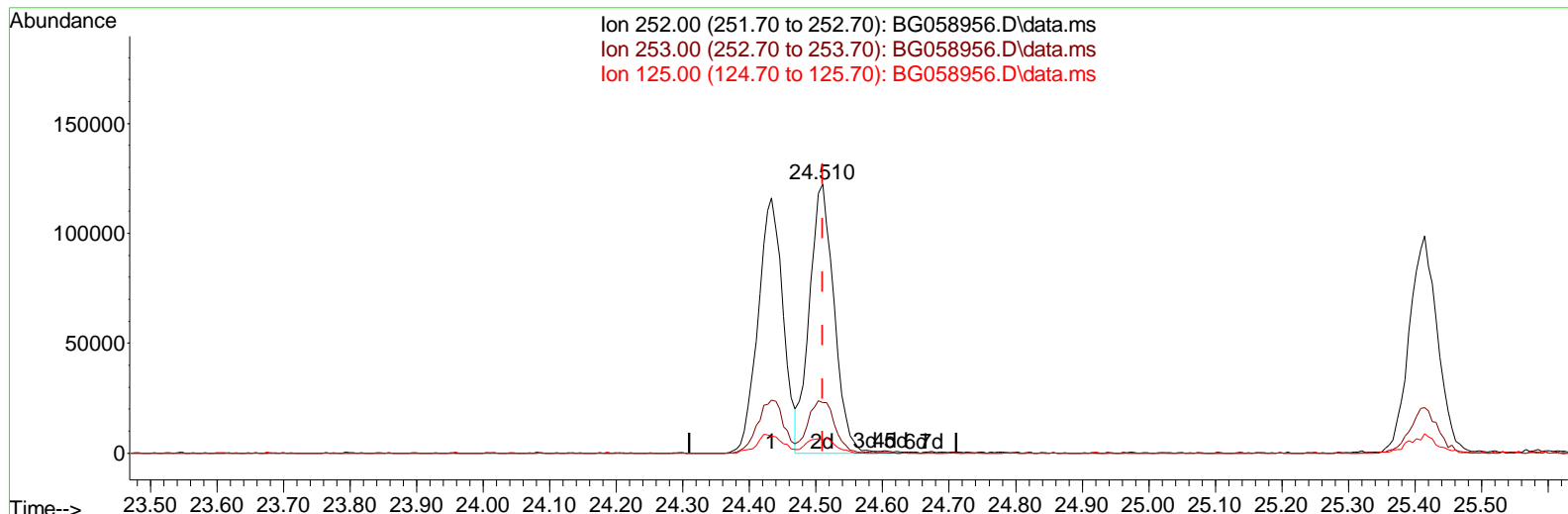
SSTD010408

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/11/2023

Supervised By :mohammad ahmed 09/13/2023

Quant Time: Sep 10 23:05:36 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Sep 10 22:58:52 2023
 Response via : Initial Calibration



TIC: BG058956.D\data.ms

(91) Benzo(k)fluoranthene

24.510min (-0.001) 9.68 ng/ul m

response 314310

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	23.90	18.73#
125.00	6.20	4.81#
0.00	0.00	0.00

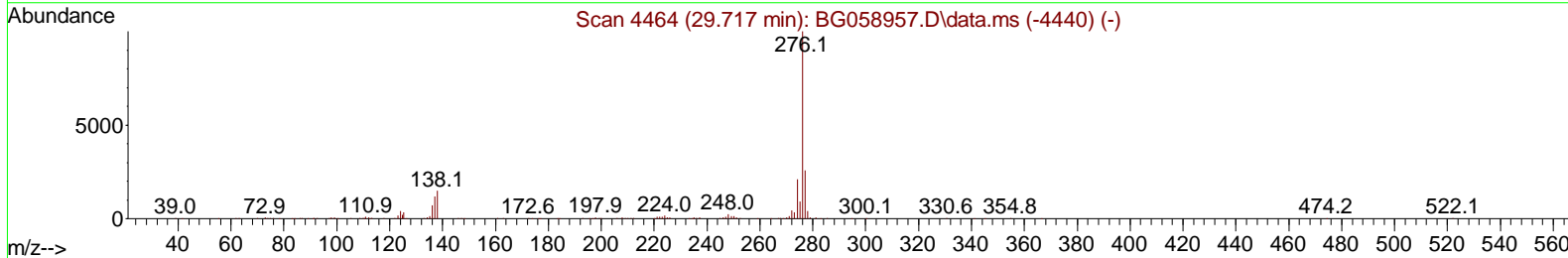
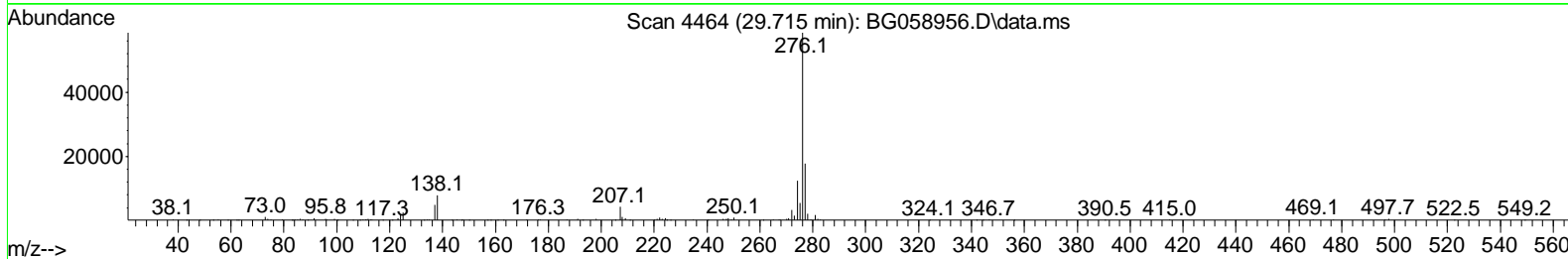
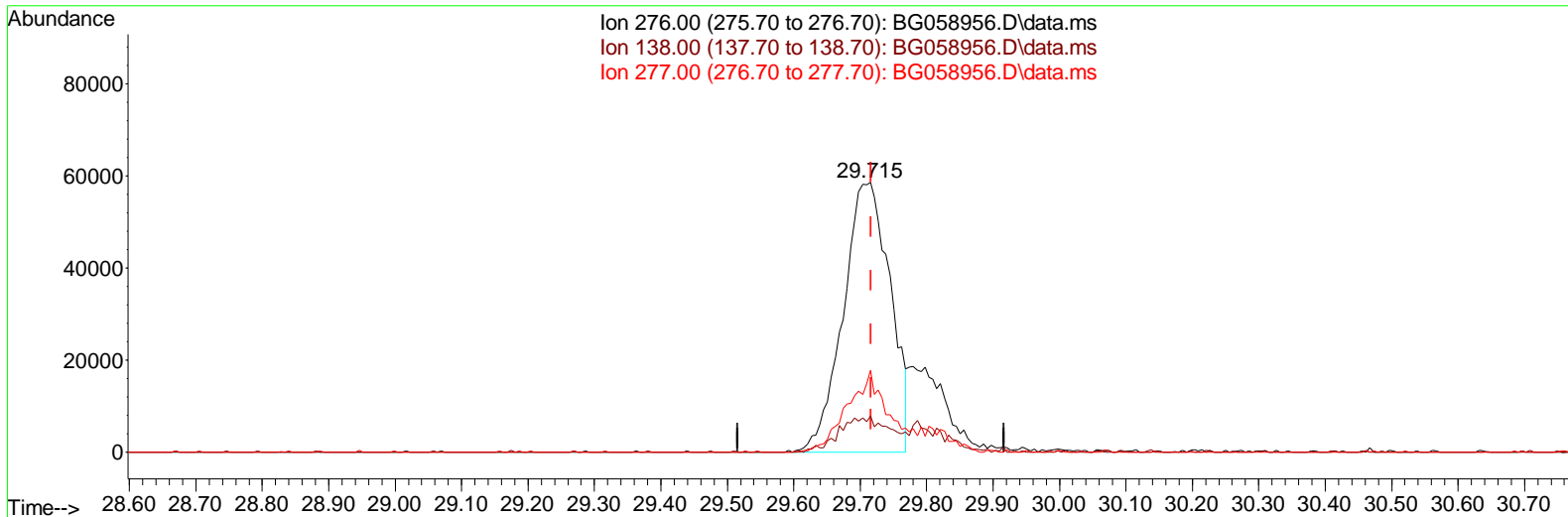
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
ClientSampleId :
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Manual IntegrationsAPPROVED

Quant Time: Sep 10 23:05:36 2023
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Reviewed By :Yogesh Patel 09/11/2023
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TIC: BG058956.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

29.715min (-0.001) 7.63 ng/ul

response 287318

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	14.80	13.29
277.00	25.90	30.36
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

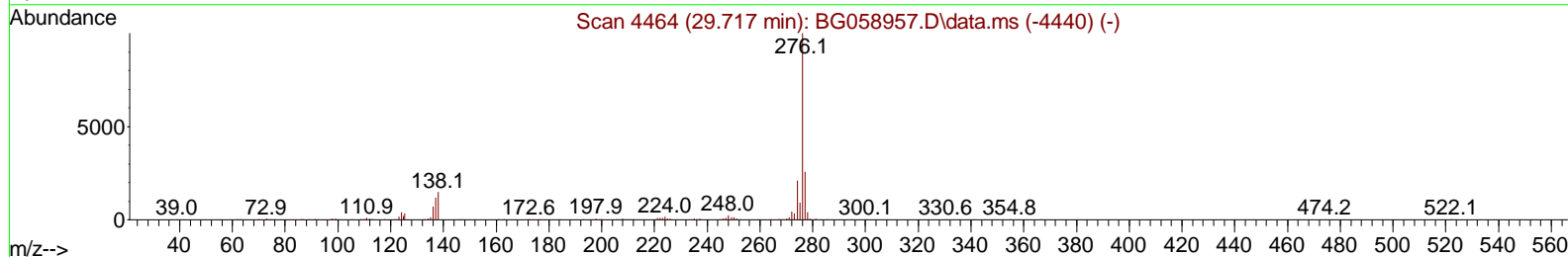
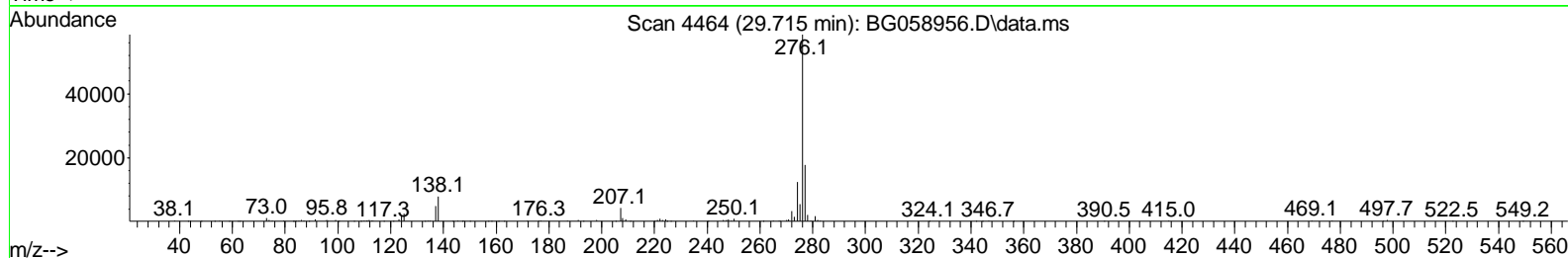
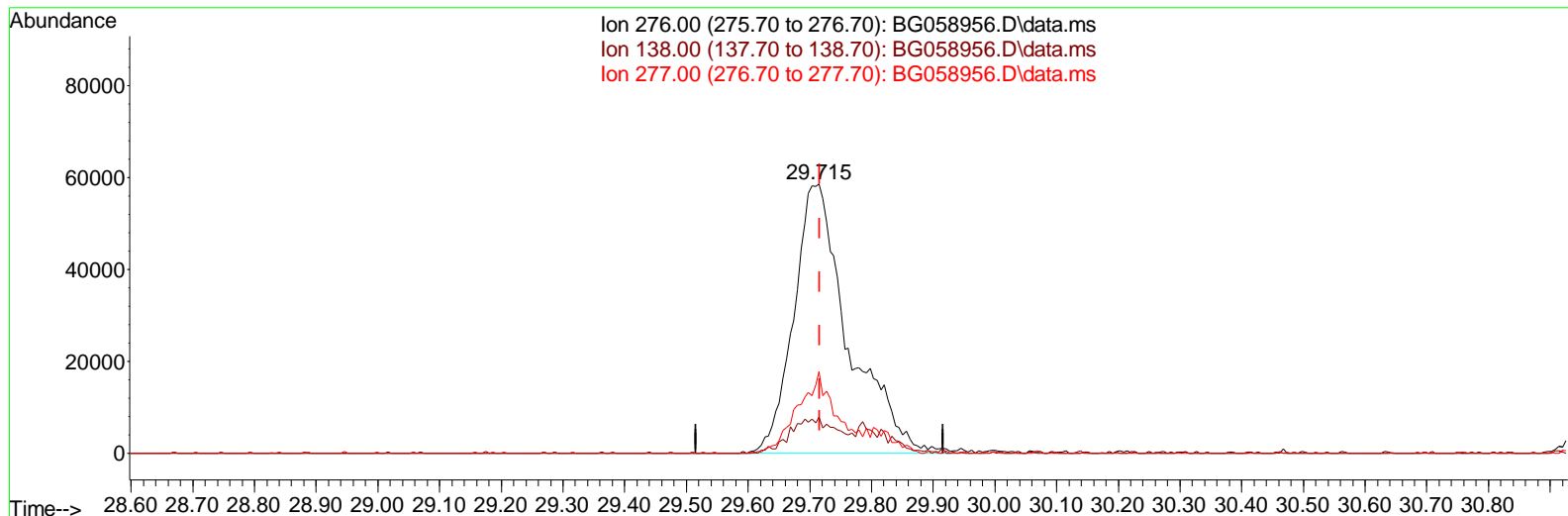
SSTD010408

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 09/11/2023

Supervised By :mohammad ahmed 09/13/2023

Quant Time: Sep 10 23:05:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
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TIC: BG058956.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

29.715min (-0.001) 9.58 ng/ul m

response 360556

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	14.80	13.29
277.00	25.90	30.36
0.00	0.00	0.00

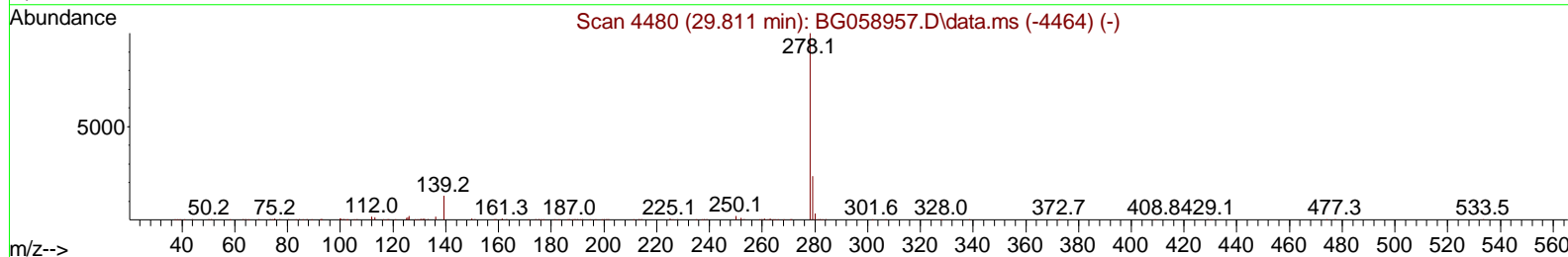
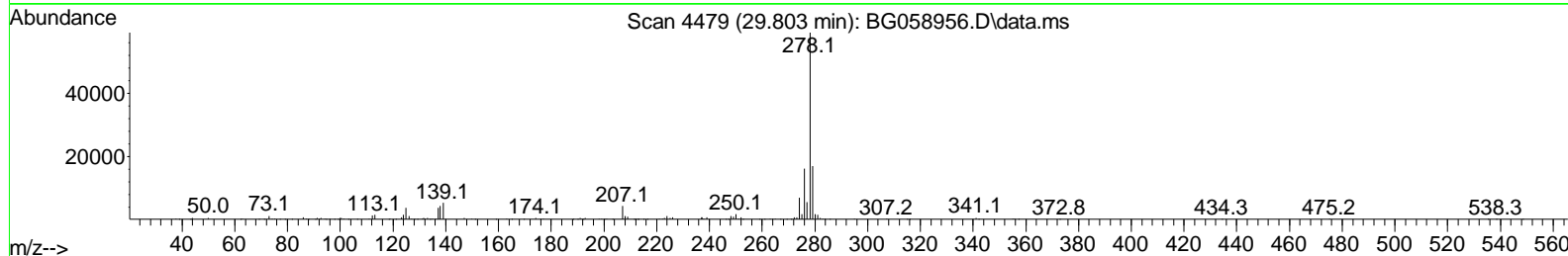
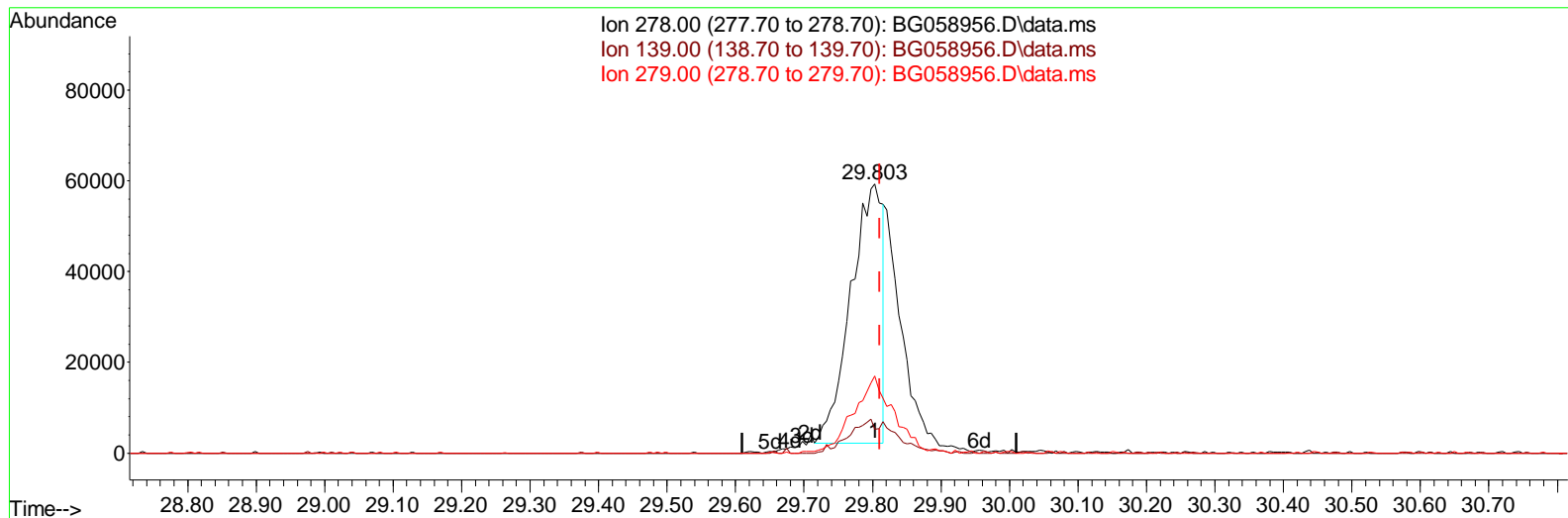
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD010408

Manual Integrations APPROVED

Quant Time: Sep 10 23:24:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Sep 10 22:58:52 2023
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Reviewed By :Yogesh Patel 09/11/2023
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TIC: BG058956.D\data.ms

(95) Dibenzo(a,h)anthracene

29.803min (-0.007) 6.01 ng/ul

response 183881

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	12.30	8.94#
279.00	23.10	28.57#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

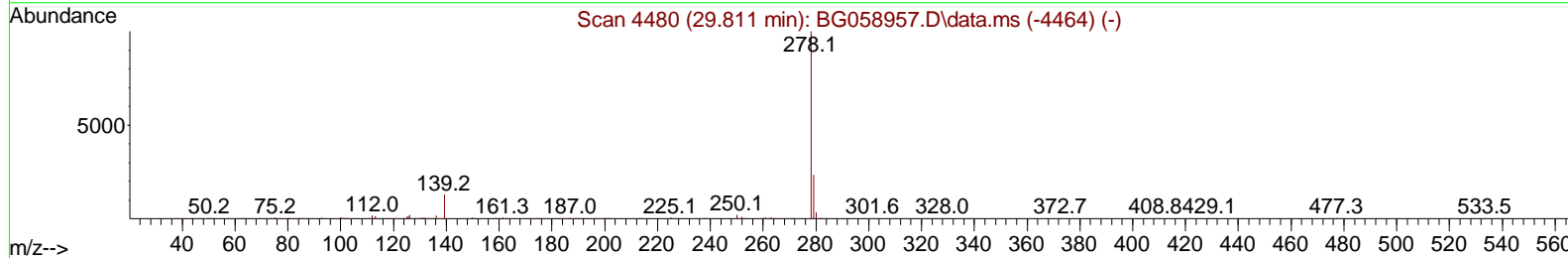
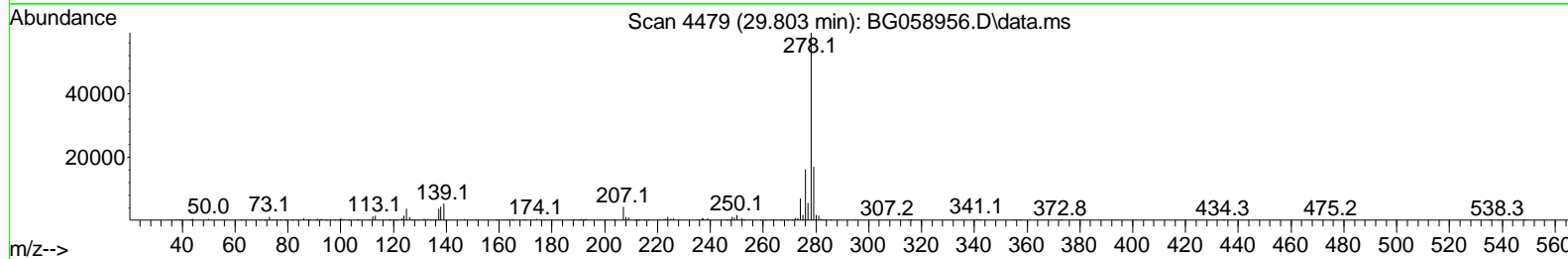
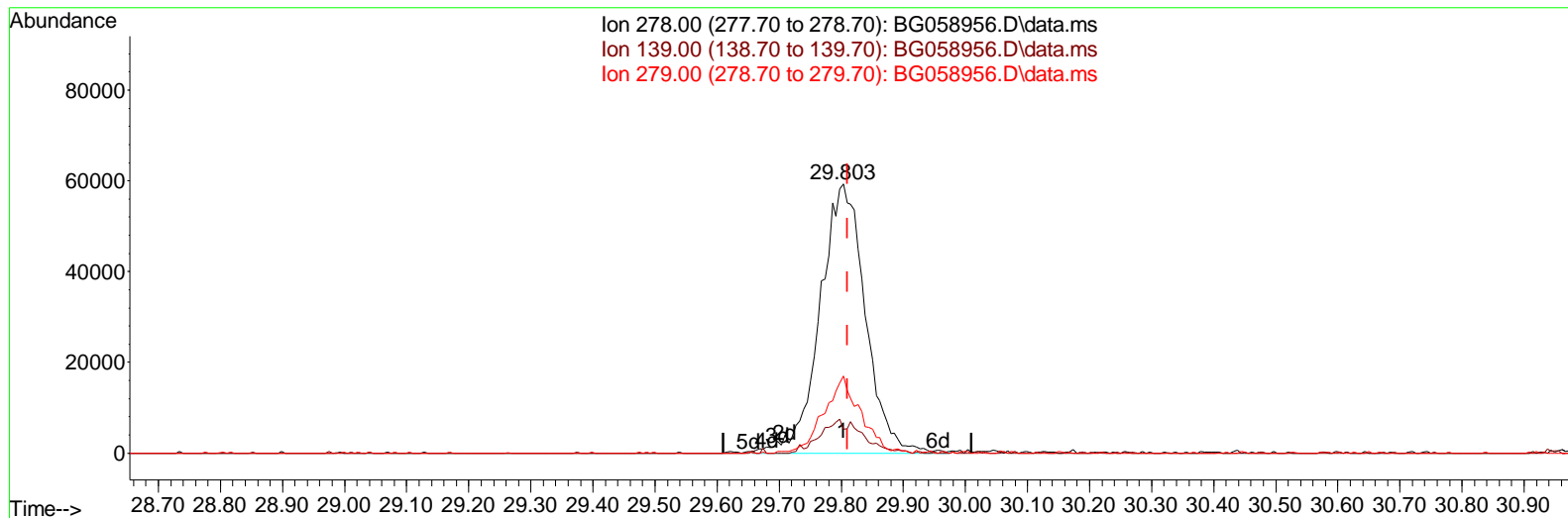
SSTD010408

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/11/2023

Supervised By :mohammad ahmed 09/13/2023

Quant Time: Sep 10 23:24:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Sep 10 22:58:52 2023
 Response via : Initial Calibration



TIC: BG058956.D\data.ms

(95) Dibenzo(a,h)anthracene

29.803min (-0.007) 9.84 ng/ul m

response 301027

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	12.30	8.94#
279.00	23.10	28.57#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

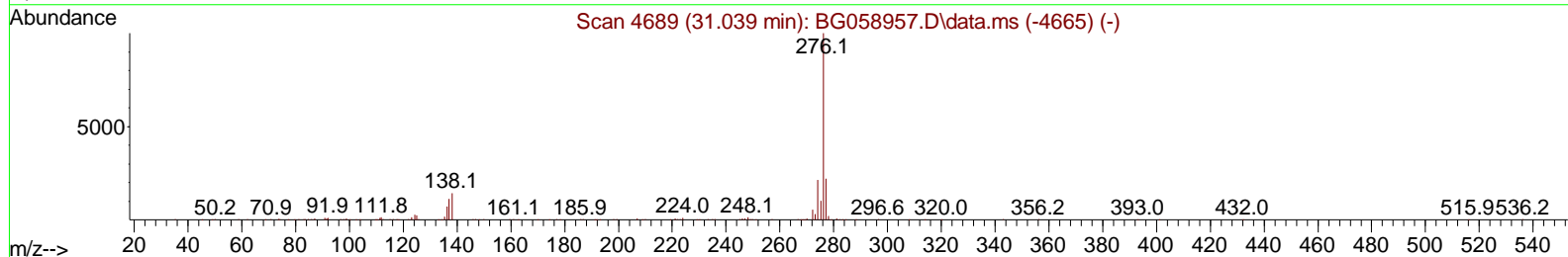
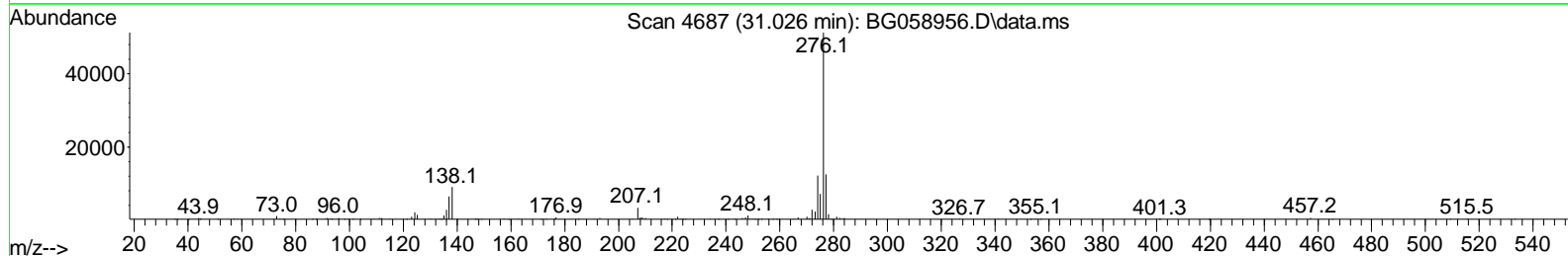
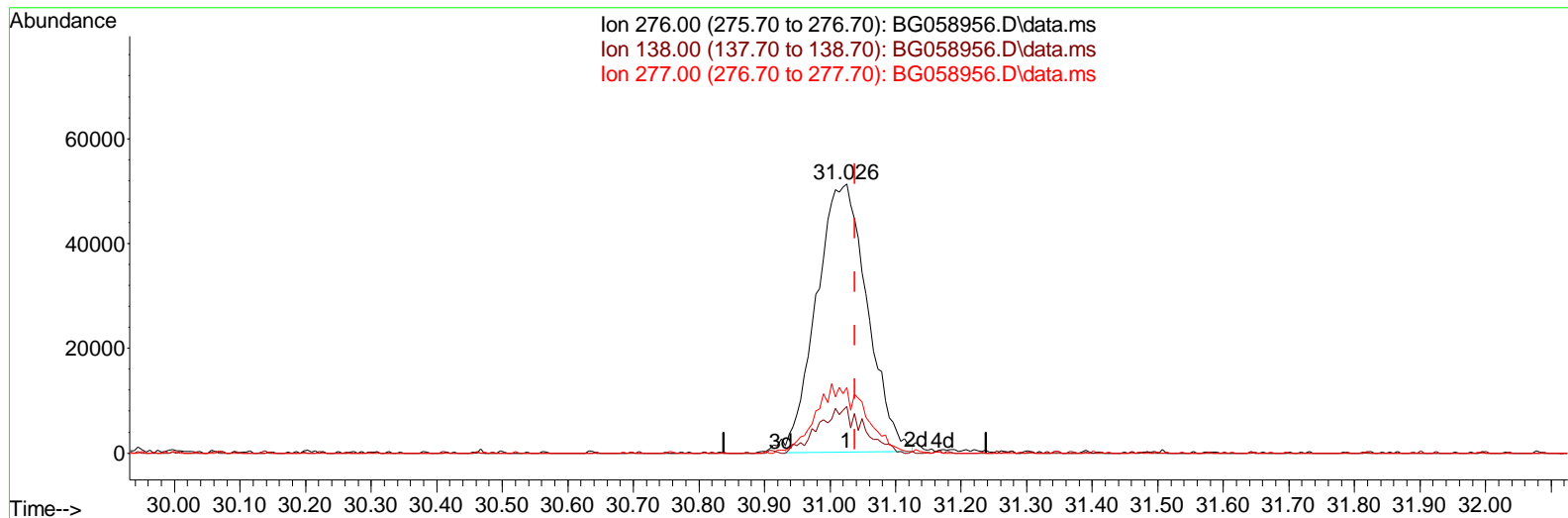
SSTD010408

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/11/2023

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TIC: BG058956.D\data.ms

(96) Benzo(g,h,i)perylene

31.026min (-0.013) 9.21 ng/ul

response 275242

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	14.30	17.45#
277.00	21.80	24.38
0.00	0.00	0.00

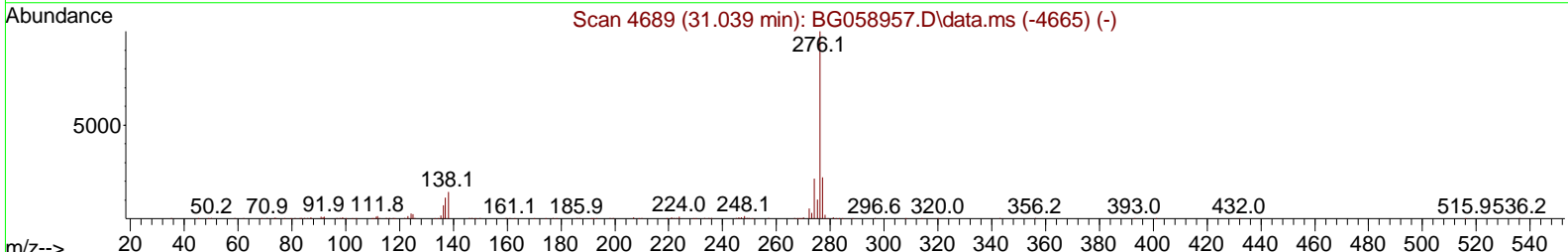
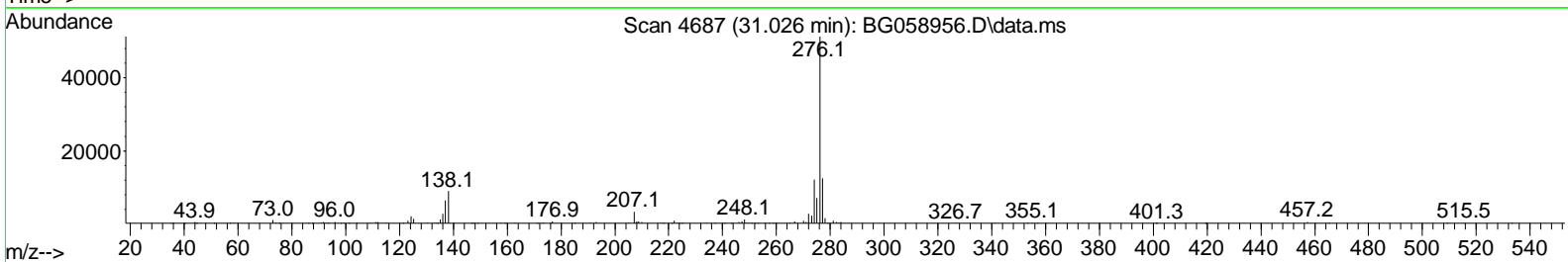
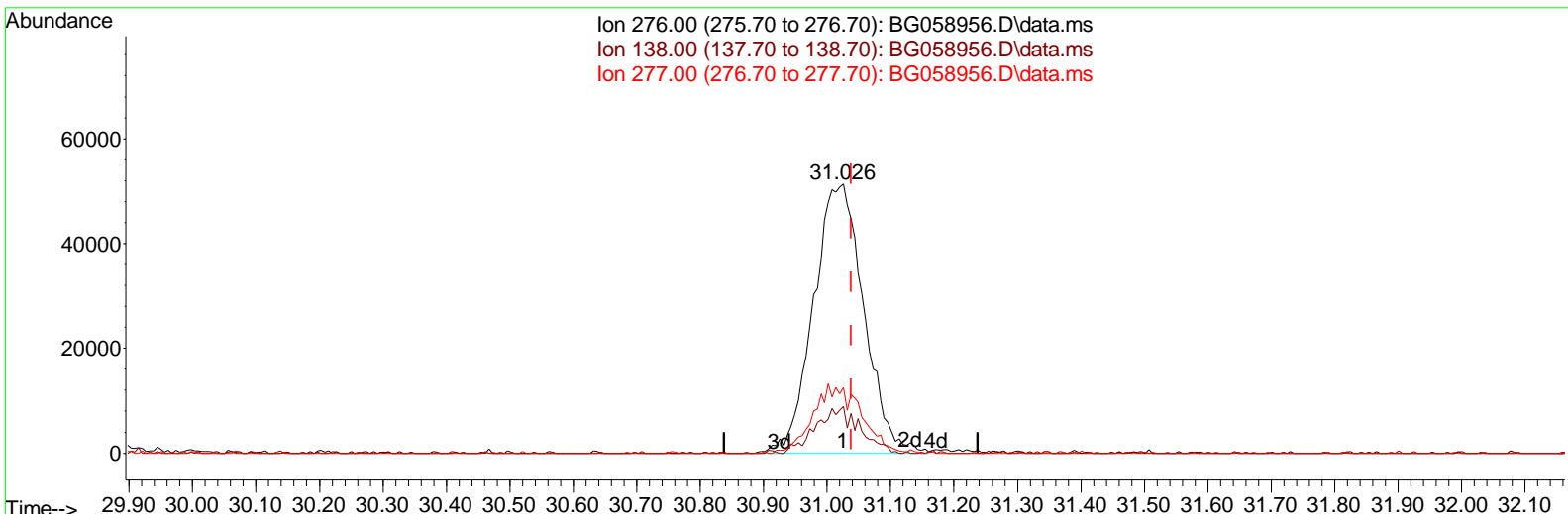
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SSTD01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
ClientSampleId :
 SSTD010408

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 09/11/2023
 Supervised By :mohammad ahmed 09/13/2023

Quant Time: Sep 10 23:24:43 2023
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TIC: BG058956.D\data.ms

(96) Benzo(g,h,i)perylene

31.026min (-0.013) 9.44 ng/ul m

response	282046	
Ion	Exp%	Act%
276.00	100.00	100.00
138.00	14.30	17.45#
277.00	21.80	24.38
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
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 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
 SSTD010408

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/11/2023
 Supervised By :mohammad ahmed 09/13/2023

Quant Time: Sep 10 23:25:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Sep 10 22:58:52 2023
 Response via : Initial Calibration

Compound	R.T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.323	152	51068	20.000	ng/ul	0.00
20) Naphthalene-d8	11.166	136	233405	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.950	164	175666	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.700	188	460073	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.024	240	479383	20.000	ng/ul	0.00
88) Perylene-d12	25.585	264	535707	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.640	96	4661	4.378	ng/uL	0.00
4) Pyridine-d5	4.069	84	37979	10.389	ng/ul	0.00
7) Phenol-d5	7.436	99	46510	9.182	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.641	67	29431	10.271	ng/ul	0.00
11) 2-Chlorophenol-d4	7.841	132	30592	10.125	ng/ul	0.00
15) 4-Methylphenol-d8	9.004	113	37107	9.452	ng/ul	0.00
21) Nitrobenzene-d5	9.527	128	14849	10.801	ng/ul	0.00
24) 2-Nitrophenol-d4	10.232	143	16180	10.743	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.761	165	37611	8.608	ng/ul	0.00
31) 4-Chloroaniline-d4	11.307	131	55038	9.820	ng/ul	0.00
46) Dimethylphthalate-d6	14.345	166	136012	10.030	ng/ul	0.00
49) Acenaphthylene-d8	14.651	160	147484	10.062	ng/ul	0.00
54) 4-Nitrophenol-d4	15.109	143	23591	13.168	ng/ul	0.00
60) Fluorene-d10	15.937	176	119290	9.334	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.037	200	19513	9.316	ng/ul	0.00
73) Anthracene-d10	17.794	188	199492	9.532	ng/ul	0.00
81) Pyrene-d10	20.068	212	249046	8.958	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.338	264	251983	9.549	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.681	88	6555	5.055	ng/uL	89
5) Pyridine	4.092	79	39542	10.275	ng/ul	91
6) Benzaldehyde	7.471	77	31895	10.683	ng/ul #	88
8) Phenol	7.459	94	50909	9.638	ng/ul #	94
10) Bis(2-Chloroethyl)ether	7.741	93	38629	9.861	ng/ul #	85
12) 2-Chlorophenol	7.870	128	30741	10.091	ng/ul #	79
13) 2-Methylphenol	8.740	108	38578	9.754	ng/ul	89
14) 2,2'-oxybis(1-Chloropr...	8.828	45	47914	10.627	ng/ul #	94
16) Acetophenone	9.163	105	61477	8.783	ng/ul	87
17) N-Nitrosodimethylamine	9.128	70	33775	9.226	ng/ul	96
18) 4-Methylphenol	9.069	108	41946	9.692	ng/ul	99
19) Hexachloroethane	9.416	117	13938	11.210	ng/ul #	75
22) Nitrobenzene	9.562	77	46945	9.984	ng/ul	94
23) Isophorone	10.074	82	100818	8.594	ng/ul	99
25) 2-Nitrophenol	10.273	139	16661	10.256	ng/ul #	88
26) 2,4-Dimethylphenol	10.291	107	47140	9.333	ng/ul #	86
27) Bis(2-Chloroethoxy)met...	10.555	93	54845	9.028	ng/ul	93
29) 2,4-Dichlorophenol	10.790	162	35210	8.501	ng/ul #	91
30) Naphthalene	11.219	128	115314	9.146	ng/ul	96
32) 4-Chloroaniline	11.331	127	50499	9.499	ng/ul	94
33) Hexachlorobutadiene	11.454	225	26457	6.523	ng/ul #	90
34) Caprolactam	12.095	113	14275m	10.419	ng/ul	
35) 4-Chloro-3-methylphenol	12.394	107	43450	8.929	ng/ul	92

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
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 Misc :
 ALS Vial : 34 Sample Multi plier: 1

Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/11/2023
 Supervised By :mohammad ahmed 09/13/2023

Quant Time: Sep 10 23: 25: 16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALI BRATI ON
 QLast Update : Sun Sep 10 22: 58: 52 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	12.800	142	92026	9.641	ng/ul	99
37) 1-Methyl naphthal ene	13.017	142	89111	9.128	ng/ul	95
39) 1, 2, 4, 5-Tetrachl oroben. . .	13.135	216	52291	7.978	ng/ul	93
40) Hexachl orocycl opentadi ene	13.094	237	59481	12.609	ng/ul #	96
41) 2, 4, 6-Tri chl orophenol	13.376	196	34657	9.099	ng/ul	99
42) 2, 4, 5-Tri chl orophenol	13.440	196	38655	9.465	ng/ul	96
43) 1, 1' -Bi phenyl	13.787	154	125783	10.523	ng/ul #	96
44) 2-Chl oronaphthal ene	13.840	162	93439	9.802	ng/ul	97
45) 2-Ni troani li ne	14.051	65	27644	13.115	ng/ul	94
47) Di methyl phthal ate	14.392	163	126381	9.167	ng/ul #	99
48) 2, 6-Di ni trotol uene	14.533	165	22014	11.250	ng/ul #	88
50) Acenaphthyl ene	14.680	152	162660	9.945	ng/ul #	95
51) 3-Ni troani li ne	14.868	138	22241	11.279	ng/ul	85
52) Acenaphthene	15.009	153	107001	10.048	ng/ul	95
53) 2, 4-Di ni trophenol	15.062	184	21440	12.952	ng/ul	91
55) 4-Ni trophenol	15.127	109	26516	11.833	ng/ul #	90
56) Di benzofuran	15.344	168	156218	9.826	ng/ul	98
57) 2, 4-Di ni trotol uene	15.303	165	31877	11.292	ng/ul #	91
58) 2, 3, 4, 6-Tetrachl orophenol	15.550	232	37883	8.780	ng/ul	95
59) Di ethyl phthal ate	15.738	149	130051	10.495	ng/ul	98
61) Fl uorene	15.990	166	132313	9.803	ng/ul	96
62) 4-Chl orophenyl -phenyl e. . .	15.978	204	70444	8.700	ng/ul	90
63) 4-Ni troani li ne	16.025	138	22134	12.064	ng/ul	88
66) 4, 6-Di ni tro-2-methyl ph. . .	16.055	198	20786	9.453	ng/ul #	92
67) N-Ni trosodi phenyl ami ne	16.190	169	114238	10.097	ng/ul	92
68) 4-Bromophenyl -phenyl ether	16.872	248	47395	8.484	ng/ul	94
69) Hexachl orobenzene	16.960	284	55658	8.634	ng/ul	89
70) Atrazi ne	17.124	200	51598	11.359	ng/ul	96
71) Pentachl orophenol	17.312	266	34306	8.665	ng/ul #	83
72) Phenanthrene	17.741	178	230436	10.317	ng/ul	96
74) Anthracene	17.829	178	237183	10.401	ng/ul	98
75) 1, 2, 3, 4-Tetrachl oroben. . .	13.746	216	57606	8.623	ng/uL#	86
76) Pentachl orobenzene	15.238	250	56653	8.414	ng/uL	95
77) Carbazol e	18.105	167	206201	11.913	ng/ul	97
78) Di -n-butyl phthal ate	18.617	149	225807	14.512	ng/ul	98
80) Fl uoranthene	19.733	202	285195	8.441	ng/ul	97
82) Pyrene	20.097	202	298184	8.810	ng/ul #	96
83) Butyl benzyl phthal ate	20.961	149	89241	15.207	ng/ul	99
84) 3, 3' -Di chl orobenzi di ne	21.913	252	107061	10.225	ng/ul	93
85) Benzo(a)anthracene	22.001	228	309589	9.486	ng/ul	99
86) Bi s(2-ethyl hexyl)phtha. . .	21.848	149	133486	14.582	ng/ul	98
87) Chrysene	22.071	228	285207	9.328	ng/ul	99
89) Di -n-octyl phthal ate	23.182	149	220411	13.175	ng/ul	100
90) Benzo(b)fl uoranthene	24.433	252	302744	9.384	ng/ul	97
91) Benzo(k)fl uoranthene	24.510	252	314310m	9.683	ng/ul	
93) Benzo(a)pyrene	25.414	252	288513	9.558	ng/ul #	98
94) I ndeno(1, 2, 3-cd)pyrene	29.715	276	360556m	9.581	ng/ul	
95) Di benzo(a, h)anthracene	29.803	278	301027m	9.837	ng/ul	
96) Benzo(g, h, i)peryl ene	31.026	276	282046m	9.440	ng/ul	

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

Instrument :

BNA_G

ClientSampleId :

SSTD010408

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/11/2023

Supervised By :mohammad ahmed 09/13/2023

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058956.D
 Acq On : 10 Sep 2023 17:29
 Operator : MA/JU
 Sample : SST01008
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :

BNA_G

Client Sample Id :

SST010408

Manual Integrations APPROVED

Reviewed By : Yogesh Patel 09/11/2023

Supervised By : mohammad ahmed 09/13/2023

Quant Time: Sep 10 23:25:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Sep 10 22:58:52 2023
 Response via : Initial Calibration

