

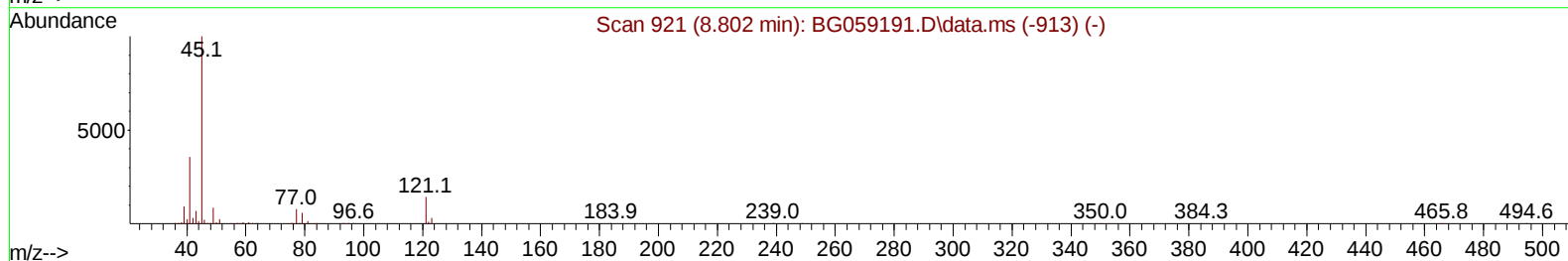
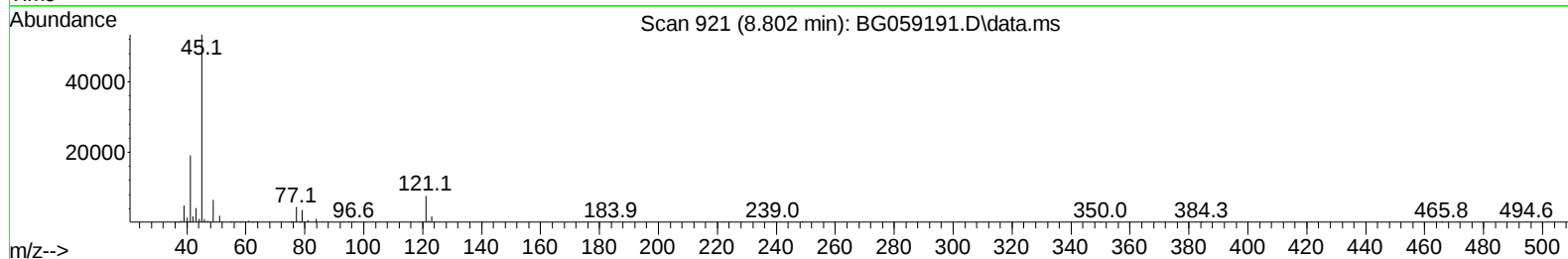
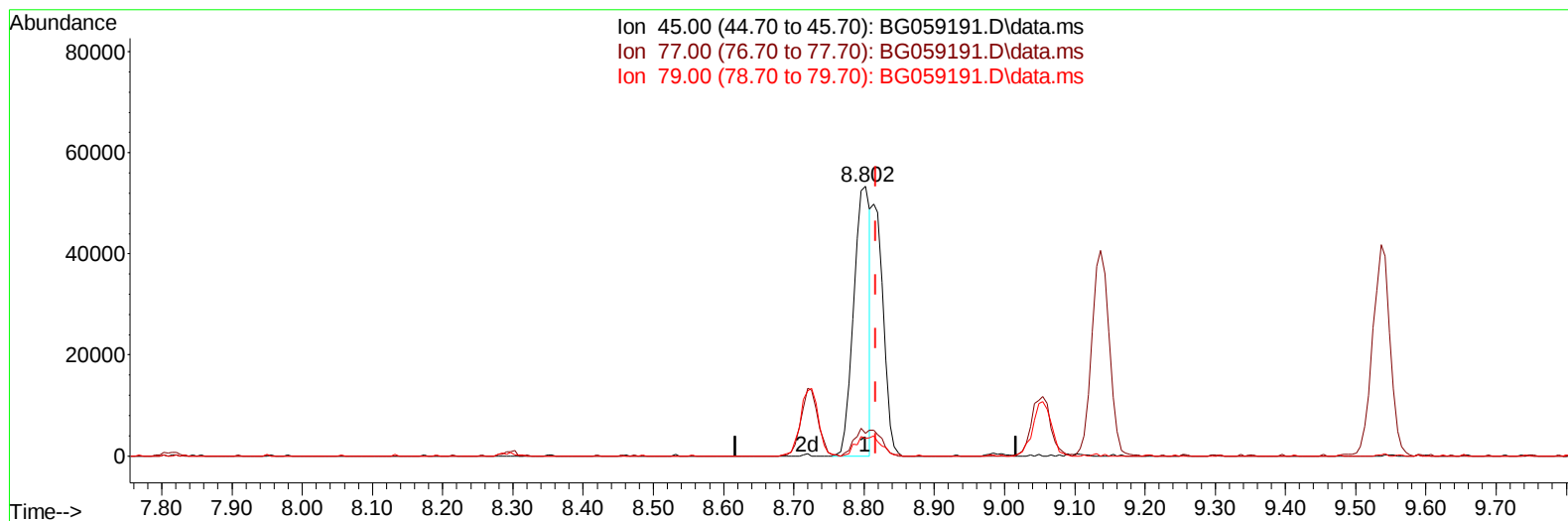
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092623\
 Data File : BG059191.D
 Acq On : 26 Sep 2023 10:15
 Operator : MA/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Sep 26 22:31:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 25 22:33:28 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059191.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.802min (-0.015) 14.10 ng/u1

response 86539

Ion	Exp%	Act%
45.00	100.00	100.00
77.00	9.70	8.35
79.00	7.00	6.60
0.00	0.00	0.00

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BNA_G

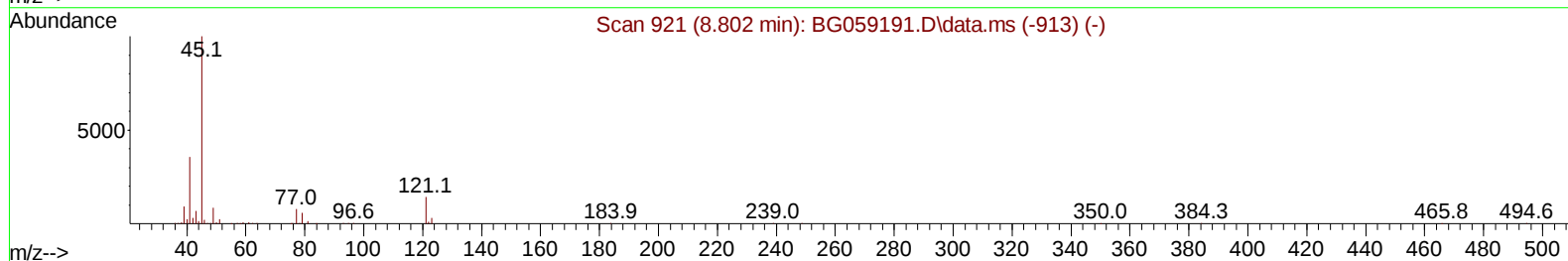
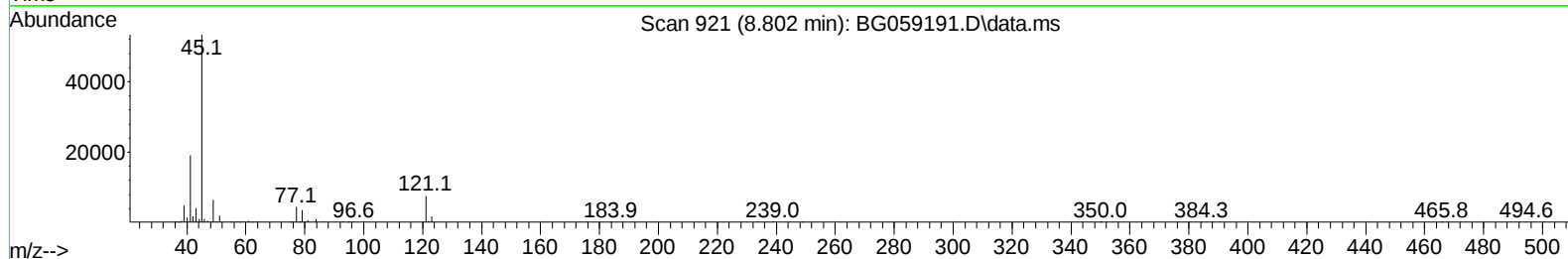
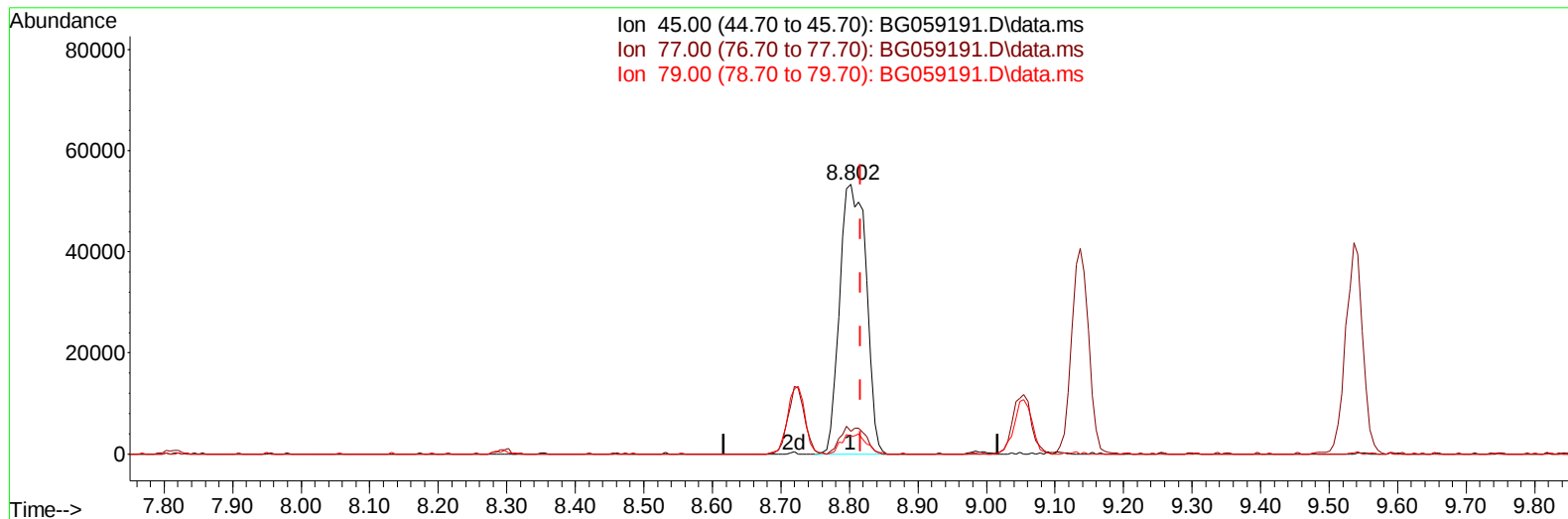
LabSampleId :

SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Sep 26 22:31:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
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 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059191.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.802min (-0.015) 23.28 ng/ul m

response 142902

Ion	Exp%	Act%
45.00	100.00	100.00
77.00	9.70	8.35
79.00	7.00	6.60
0.00	0.00	0.00

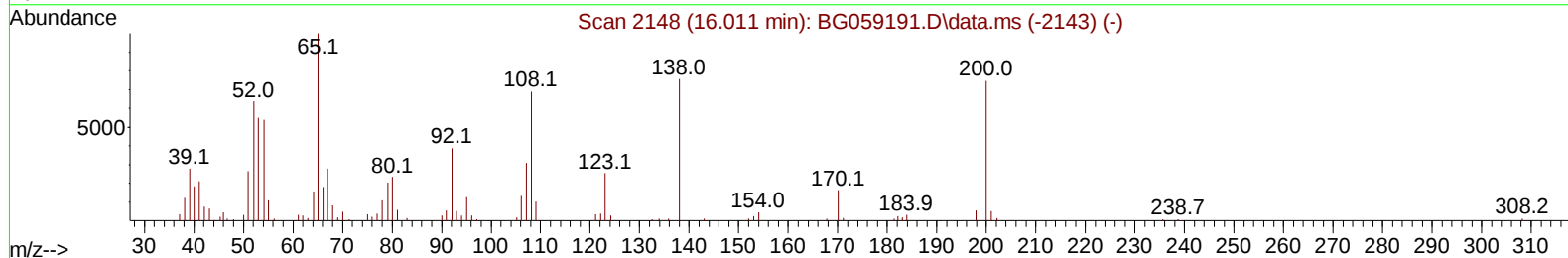
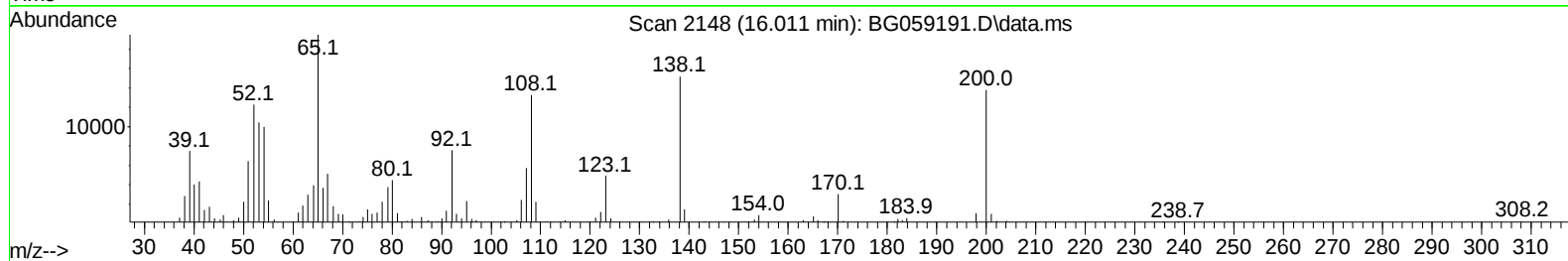
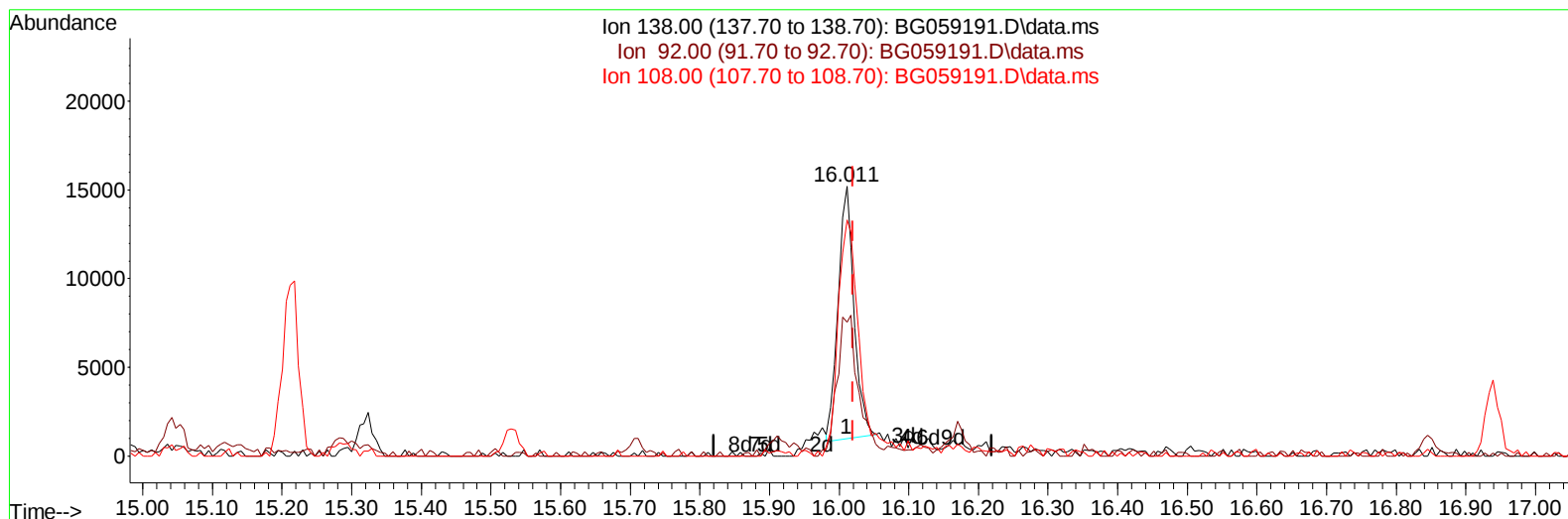
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Instrument :
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Manual IntegrationsAPPROVED

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TIC: BG059191.D\data.ms

(63) 4-Nitroaniline

16.011min (-0.009) 11.83 ng/ul

response 22685

Ion	Exp%	Act%
138.00	100.00	100.00
92.00	48.80	49.70
108.00	66.90	87.50#
0.00	0.00	0.00

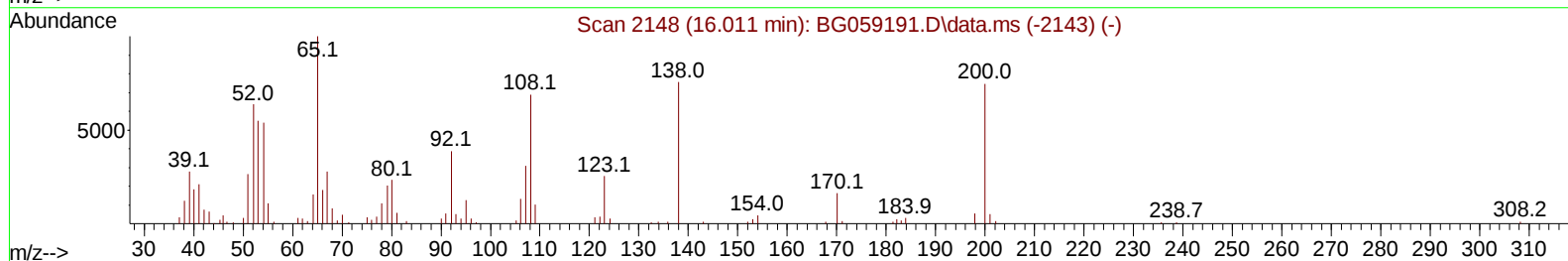
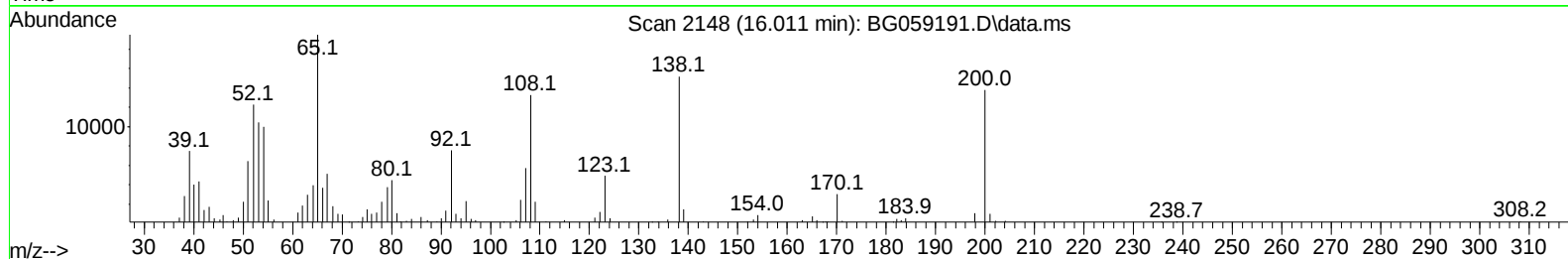
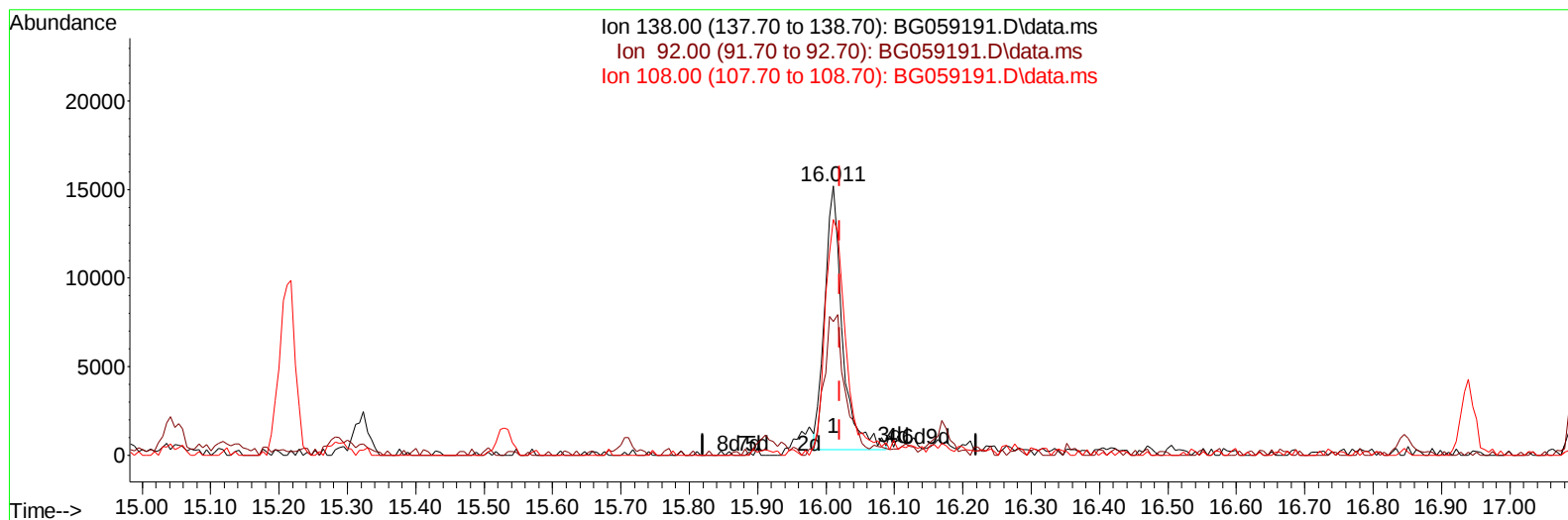
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092623\
 Data File : BG059191.D
 Acq On : 26 Sep 2023 10:15
 Operator : MA/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
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Manual IntegrationsAPPROVED

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Reviewed By :Yogesh Patel 09/27/2023
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059191.D\data.ms

(63) 4-Nitroaniline

16.011min (-0.009) 14.02 ng/ul m

response 26882

Ion	Exp%	Act%
138.00	100.00	100.00
92.00	48.80	49.70
108.00	66.90	87.50#
0.00	0.00	0.00

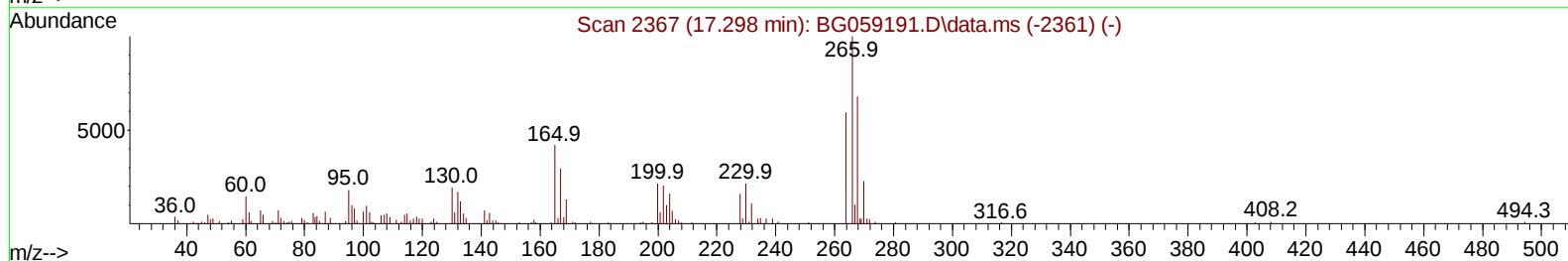
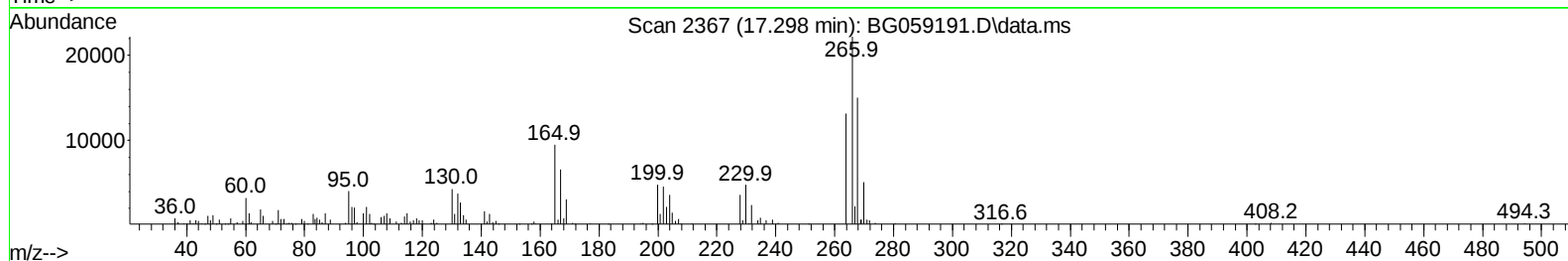
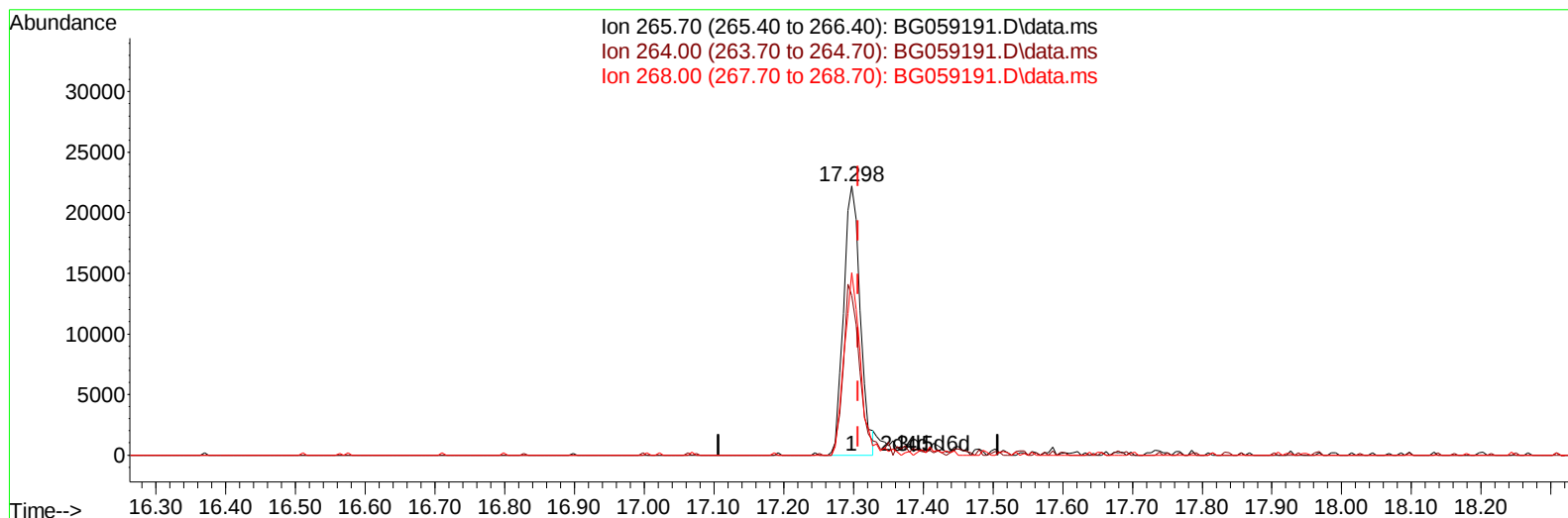
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Data File : BG059191.D
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TIC: BG059191.D\data.ms

(71) Pentachlorophenol (C)

17.298min (-0.009) 15.89 ng/u1

response 36527

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	63.30	59.26
268.00	70.50	70.76
0.00	0.00	0.00

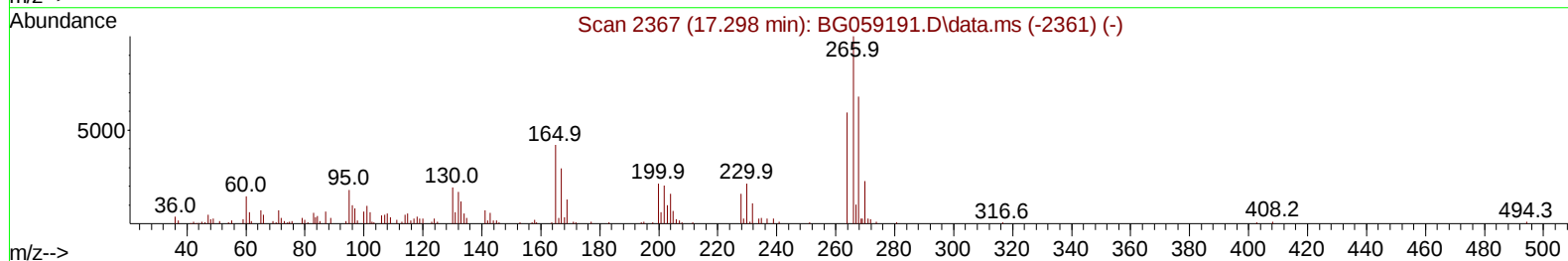
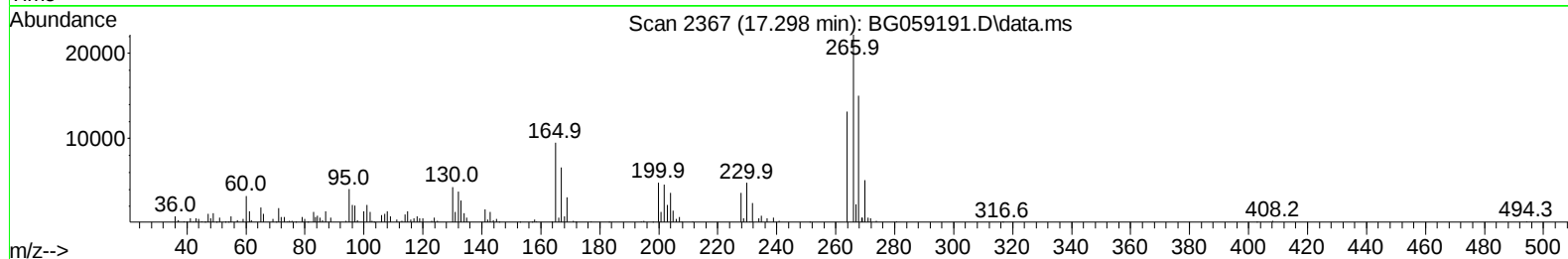
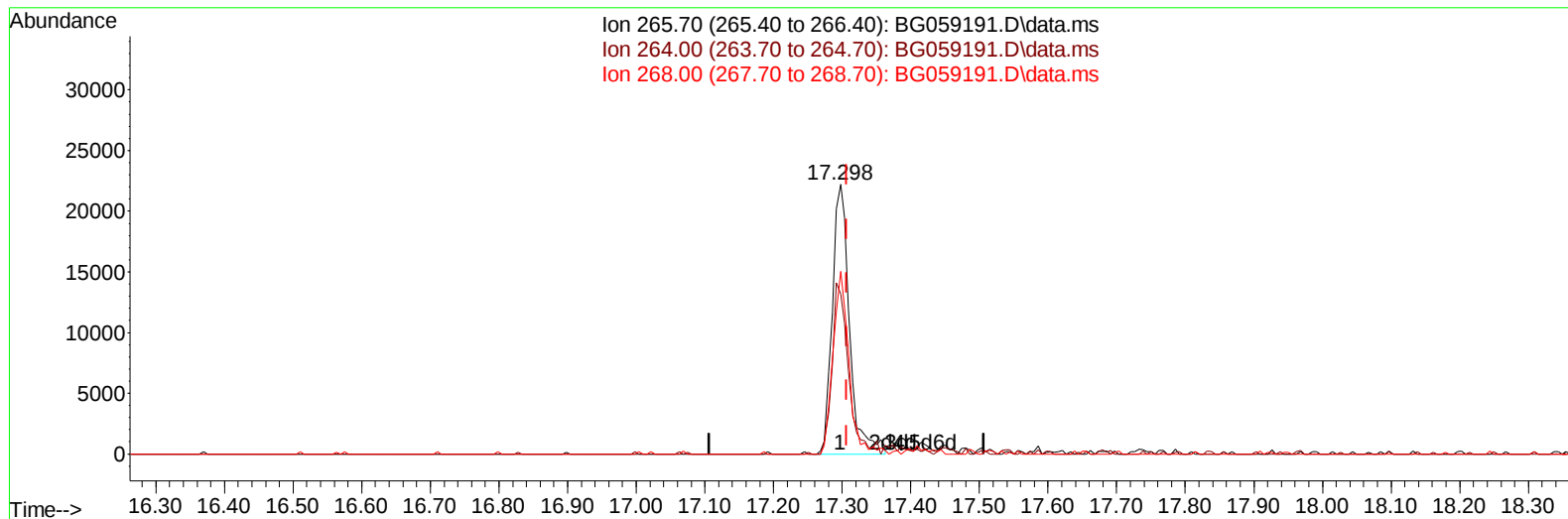
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092623\
 Data File : BG059191.D
 Acq On : 26 Sep 2023 10:15
 Operator : MA/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Sep 26 22:31:47 2023
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Reviewed By :Yogesh Patel 09/27/2023
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059191.D\data.ms

(71) Pentachlorophenol (C)

17.298min (-0.009) 16.86 ng/ul m

response 38767

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	63.30	59.26
268.00	70.50	67.82
0.00	0.00	0.00

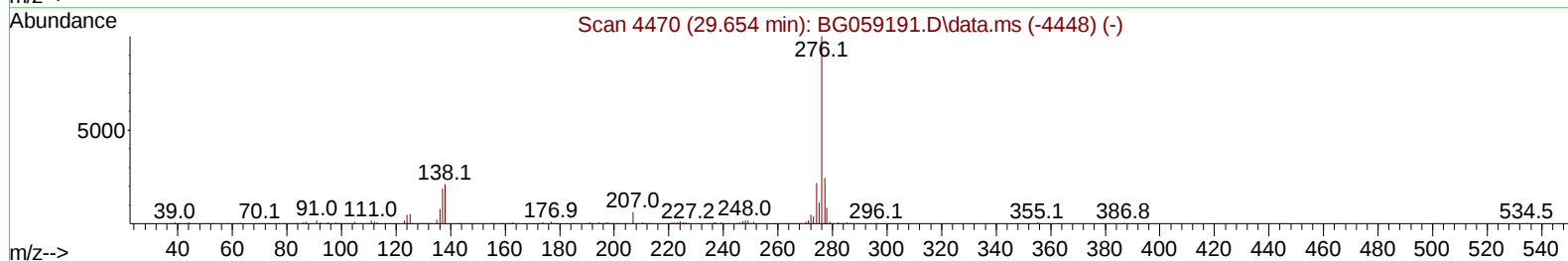
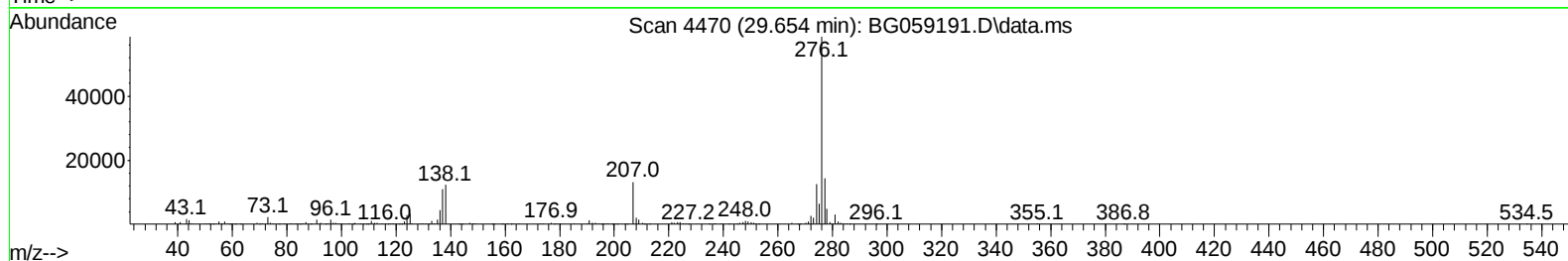
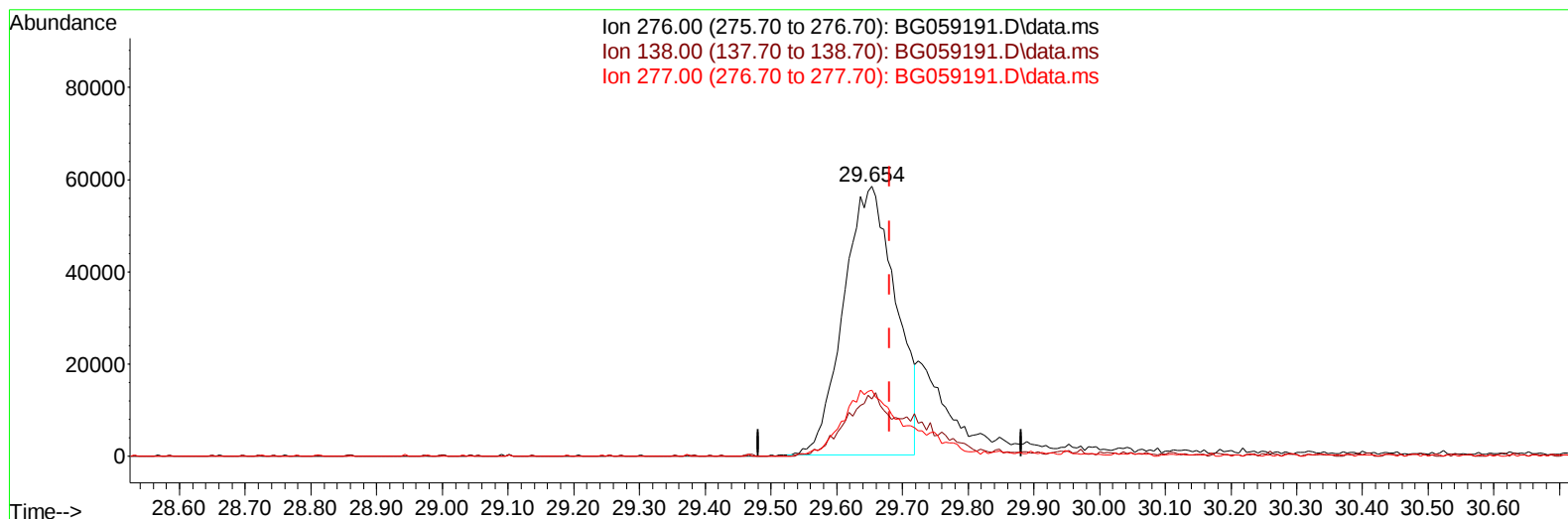
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092623\
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Instrument :
 BNA_G
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Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059191.D\data.ms

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

29.654min (-0.027) 15.18 ng/u1

response 321164

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	21.40	21.14
277.00	25.50	24.47
0.00	0.00	0.00

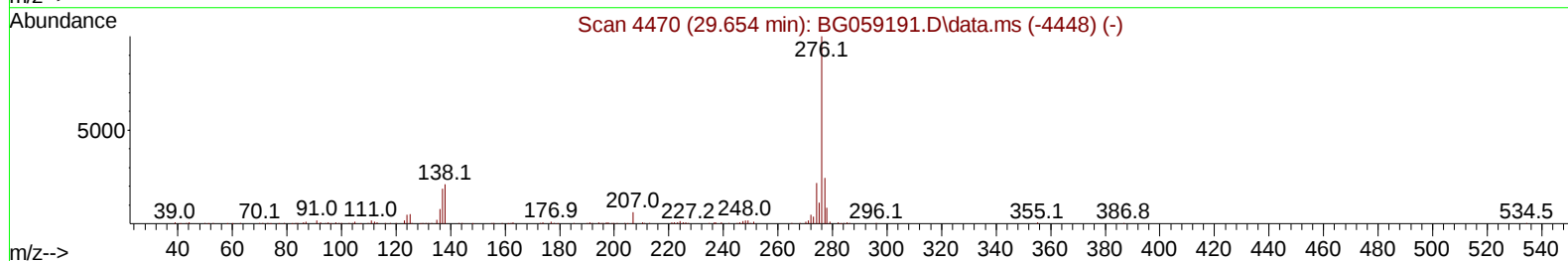
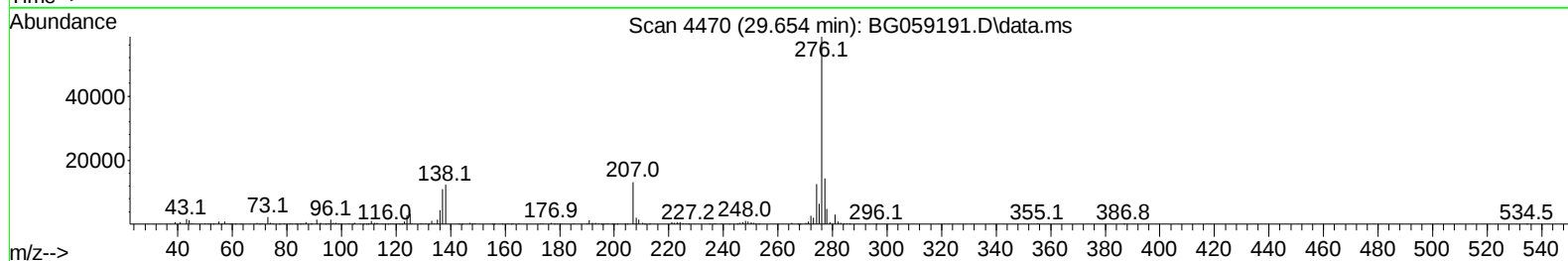
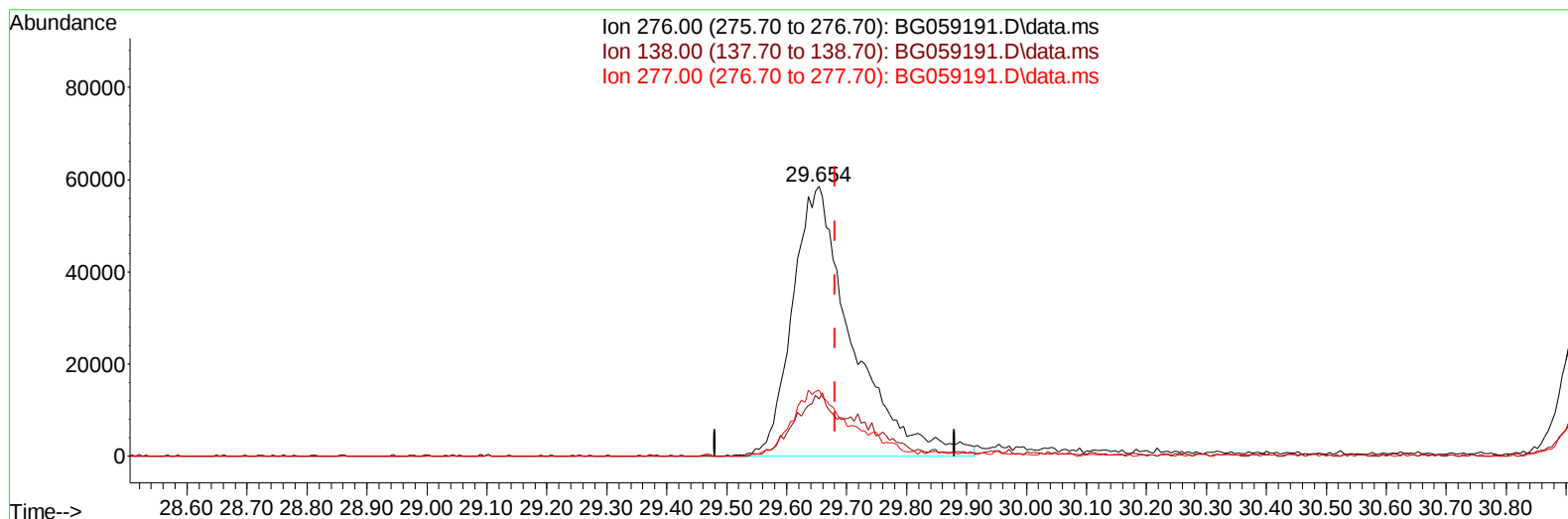
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092623\
Data File : BG059191.D
Acq On : 26 Sep 2023 10:15
Operator : MA/JU
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Sep 26 22:31:47 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Sep 25 22:33:28 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023
Supervised By :mohammad ahmed 09/29/2023



TIC: BG059191.D\data.ms

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

29.654min (-0.027) 19.15 ng/ul m

response 405264

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	21.40	21.14
277.00	25.50	24.47
0.00	0.00	0.00

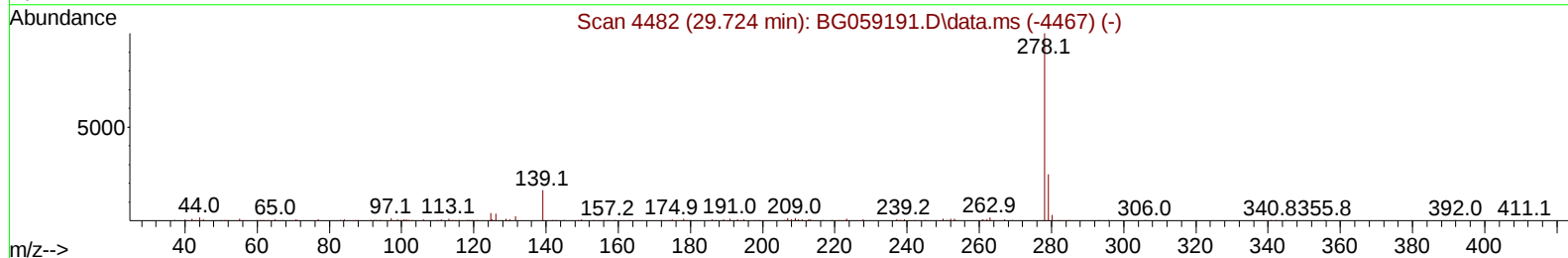
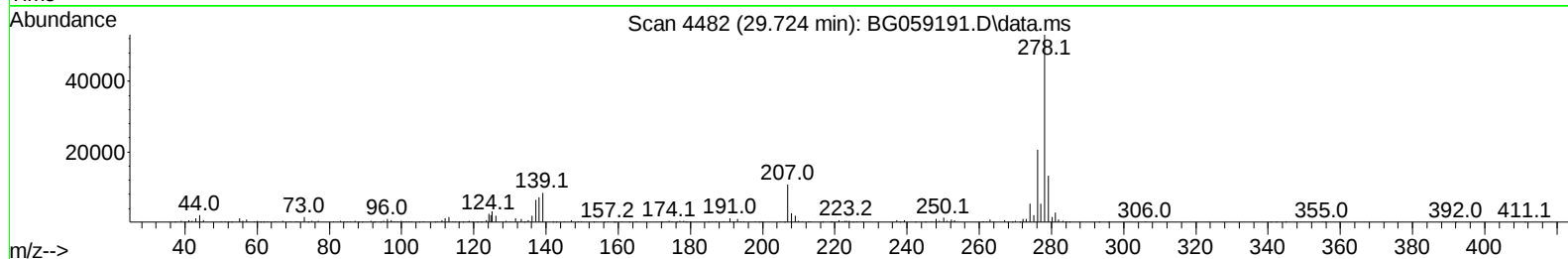
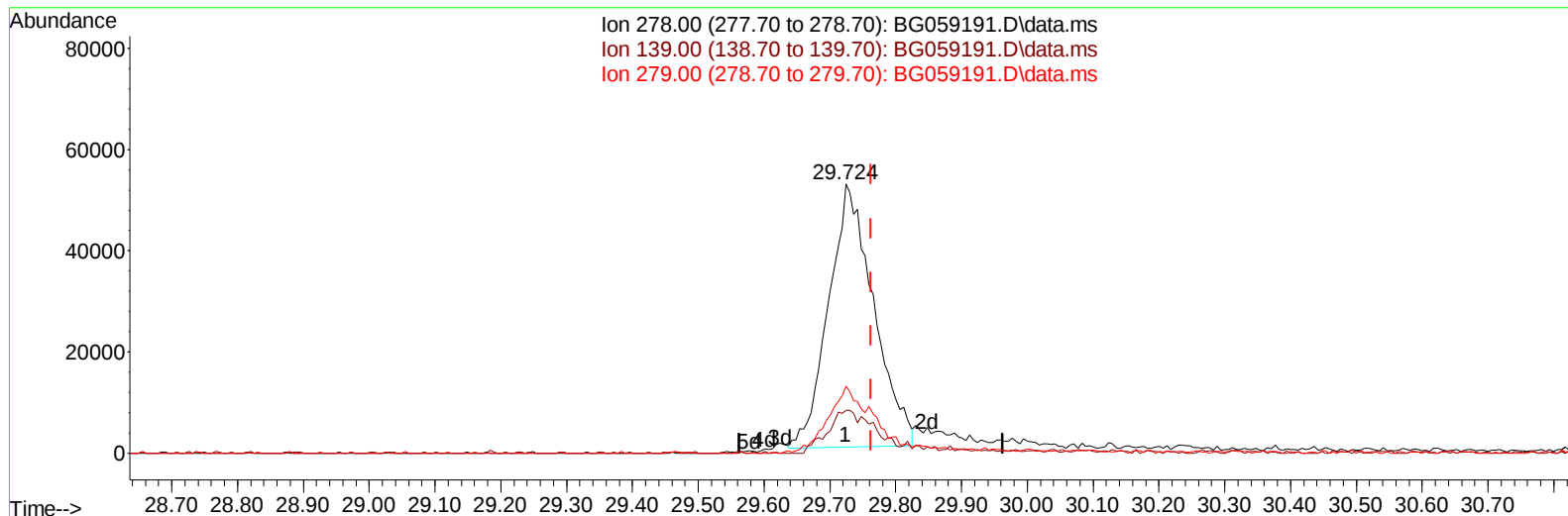
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 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
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TIC: BG059191.D\data.ms

(95) Dibenzo(a,h)anthracene

29.724min (-0.039) 14.47 ng/u1

response 246101

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	17.20	15.91
279.00	23.70	24.88
0.00	0.00	0.00

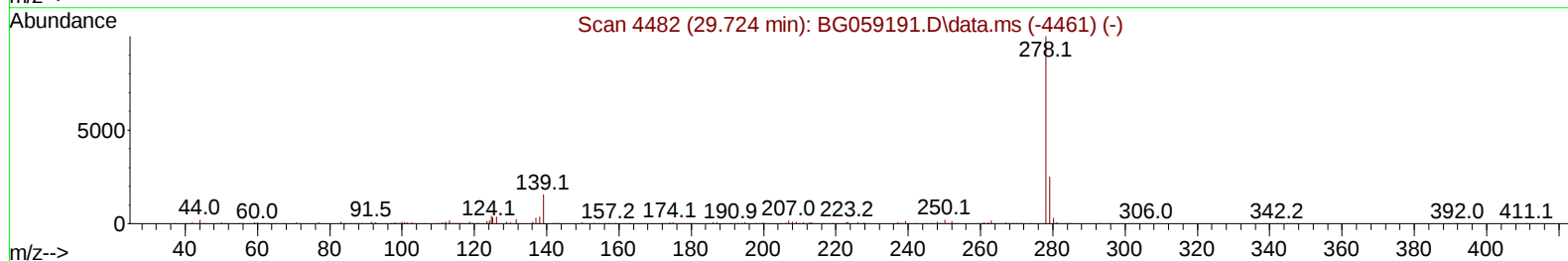
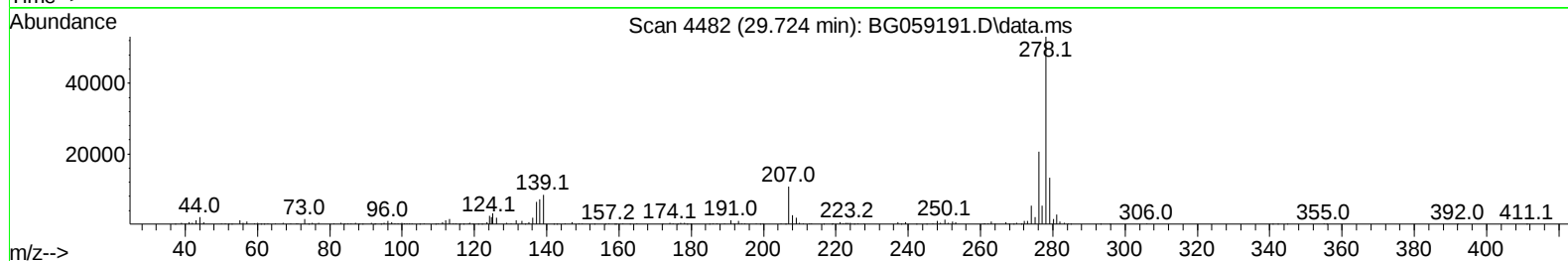
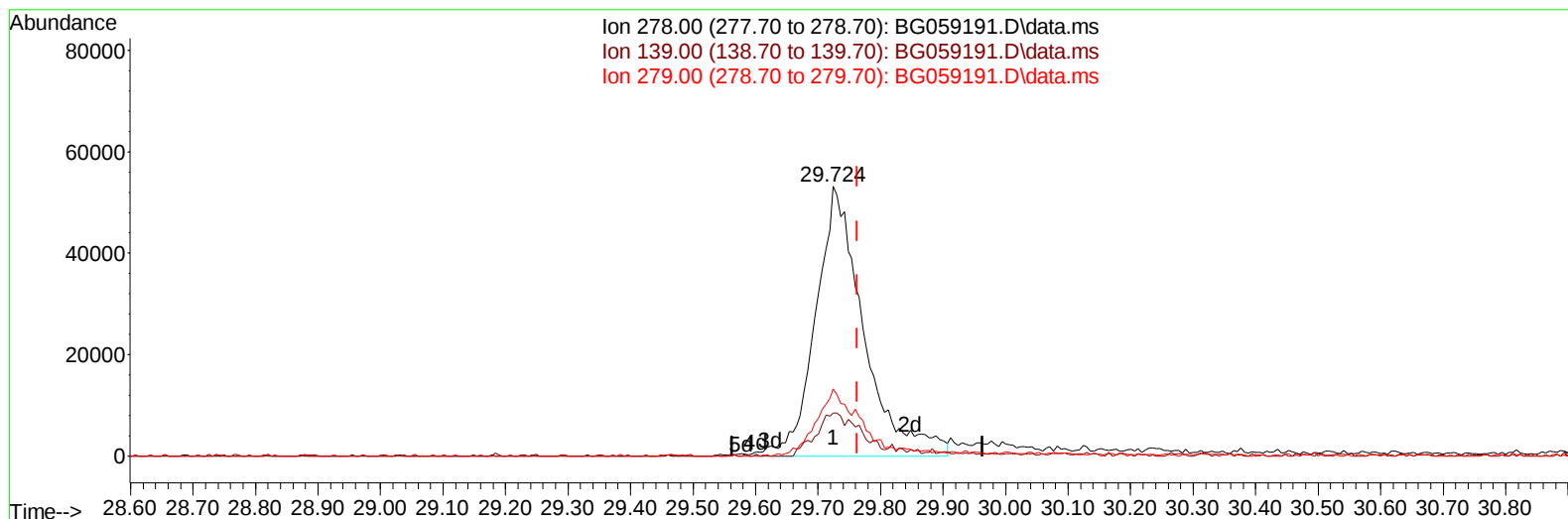
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 Operator : MA/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Sep 26 22:38:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
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 QLast Update : Mon Sep 25 22:33:28 2023
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Reviewed By :Yogesh Patel 09/27/2023
 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059191.D\data.ms

(95) Dibenzo(a,h)anthracene

29.724min (-0.039) 16.64 ng/ul m

response 282848

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	17.20	15.91
279.00	23.70	24.88
0.00	0.00	0.00

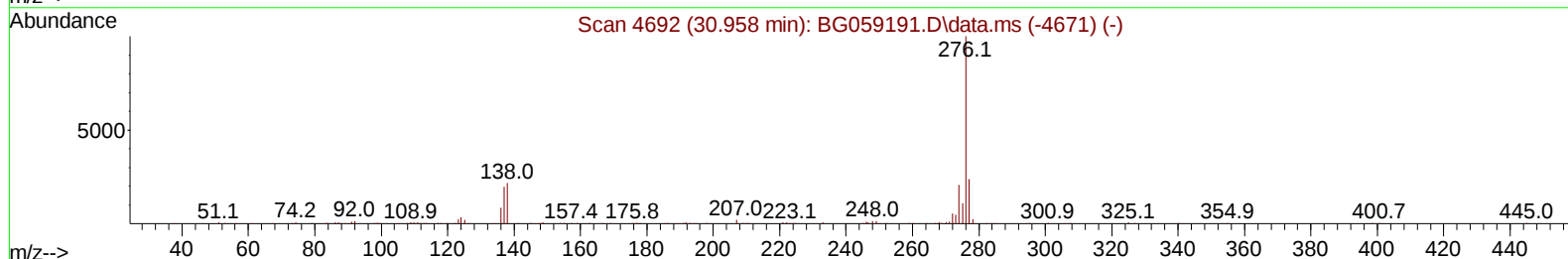
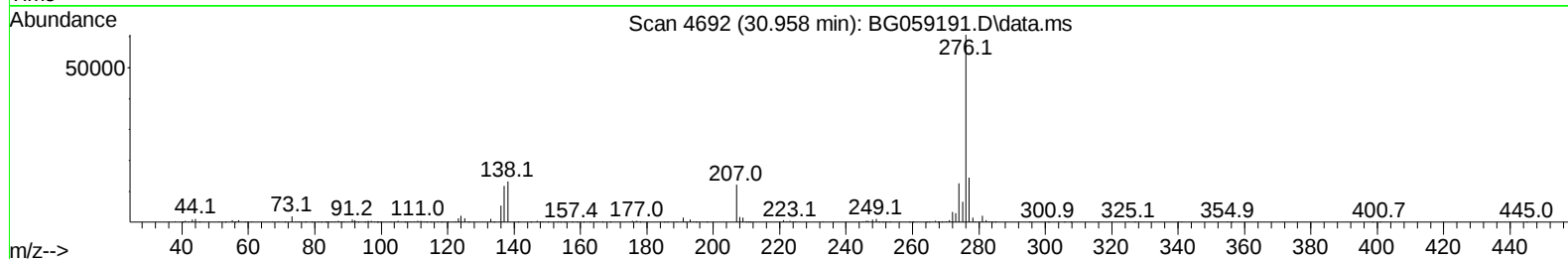
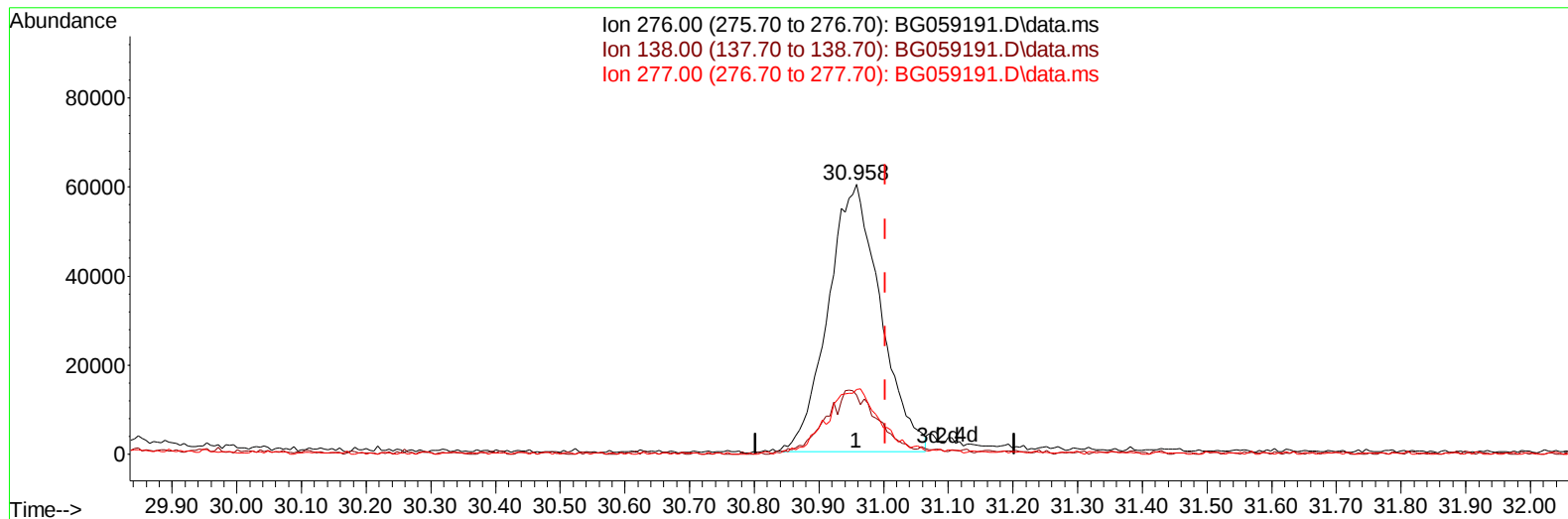
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TIC: BG059191.D\data.ms

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

30.958min (-0.044) 19.81 ng/u1

response 337670

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	21.10	21.91
277.00	24.10	24.06
0.00	0.00	0.00

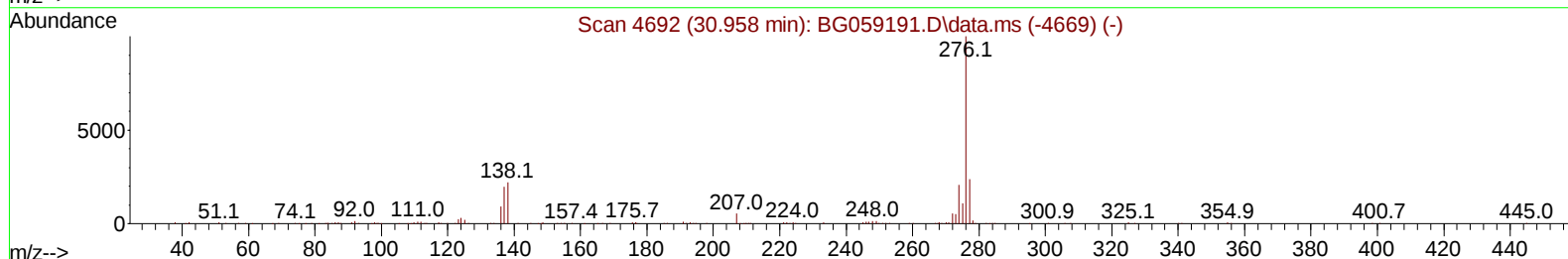
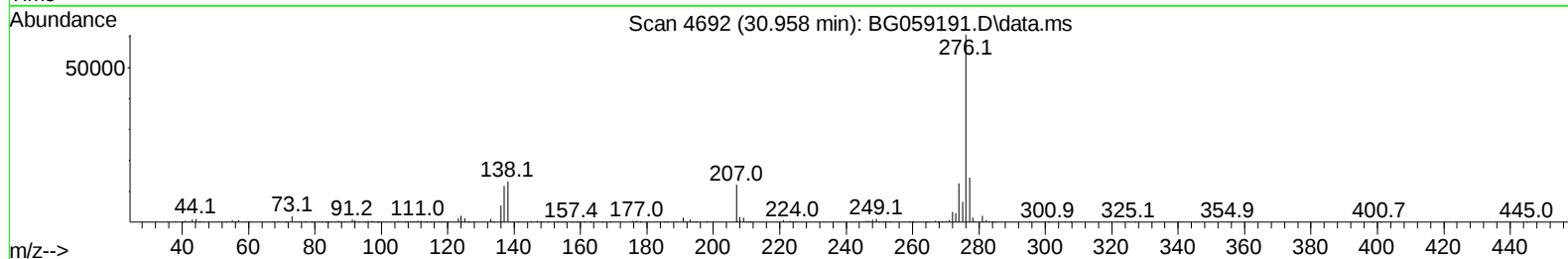
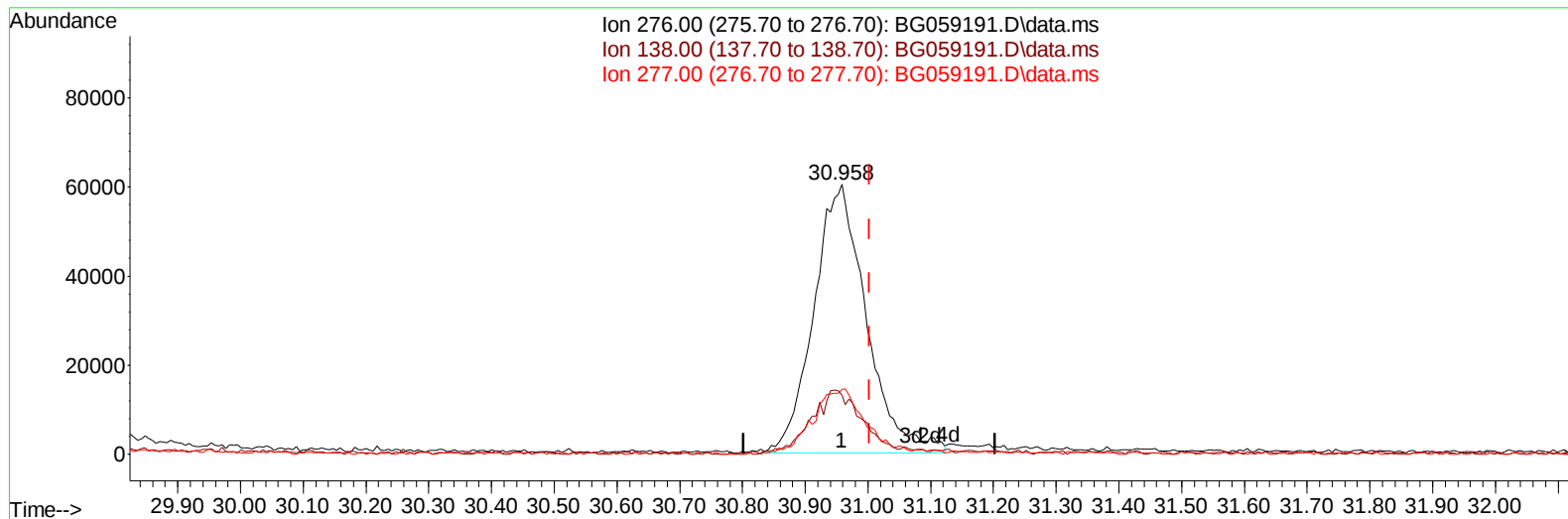
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 ALS Vial : 2 Sample Multiplier: 1

Instrument :
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 LabSampleId :
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Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 09/29/2023



TIC: BG059191.D\data.ms

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

30.958min (-0.044) 20.66 ng/ul m

response 352140

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	21.10	21.91
277.00	24.10	24.06
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	8.296	152	36433	20.000	ng/ul	# 0.00
20) Naphthalene-d8	11.140	136	185144	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.924	164	129162	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.674	188	317631	20.000	ng/ul	0.00
79) Chrysene-d12	21.998	240	272574	20.000	ng/ul	0.00
88) Perylene-d12	25.541	264	317659	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.608	96	8994	9.663	ng/uL	-0.02
4) Pyridine-d5	4.043	84	64487	23.616	ng/ul	-0.02
7) Phenol-d5	7.421	99	78677	21.962	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.615	67	49830	22.666	ng/ul	0.00
11) 2-Chlorophenol-d4	7.815	132	59613	22.549	ng/ul	0.00
15) 4-Methylphenol-d8	8.990	113	61641	21.605	ng/ul	0.00
21) Nitrobenzene-d5	9.495	128	33035	23.123	ng/ul	0.00
24) 2-Nitrophenol-d4	10.212	143	32901	22.797	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.741	165	61937	22.605	ng/ul	0.00
31) 4-Chloroaniline-d4	11.281	131	91991	21.511	ng/ul	-0.02
46) Dimethylphthalate-d6	14.319	166	232361	23.392	ng/ul	0.00
49) Acenaphthylene-d8	14.624	160	263798	21.810	ng/ul	0.00
54) 4-Nitrophenol-d4	15.112	143	27425	15.132	ng/ul	0.00
60) Fluorene-d10	15.911	176	202664	22.289	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.017	200	36778	20.366	ng/ul	-0.02
73) Anthracene-d10	17.774	188	311651	21.564	ng/ul	0.00
81) Pyrene-d10	20.047	212	360695	22.640	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.294	264	327313	21.102	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.649	88	9473	9.001	ng/uL	96
5) Pyridine	4.066	79	61296	22.029	ng/ul#	90
6) Benzaldehyde	7.439	77	47379	25.506	ng/ul	97
8) Phenol	7.450	94	79991	21.724	ng/ul	96
10) Bis(2-Chloroethyl)ether	7.709	93	66669	22.552	ng/ul	92
12) 2-Chlorophenol	7.844	128	62946	23.337	ng/ul	93
13) 2-Methylphenol	8.720	108	59202	21.665	ng/ul	93
14) 2,2'-oxybis(1-Chloropr...	8.802	45	142902m	23.282	ng/ul	
16) Acetophenone	9.137	105	99111	22.564	ng/ul	98
17) N-Nitroso-di-n-propyla...	9.096	70	52615	23.695	ng/ul#	96
18) 4-Methylphenol	9.054	108	65476	22.359	ng/ul	81
19) Hexachloroethane	9.378	117	23104	21.786	ng/ul	97
22) Nitrobenzene	9.536	77	72426	22.387	ng/ul	96
23) Isophorone	10.041	82	155687	22.441	ng/ul#	94
25) 2-Nitrophenol	10.247	139	35900	22.259	ng/ul	98
26) 2,4-Dimethylphenol	10.271	107	66695	22.118	ng/ul	93
27) Bis(2-Chloroethoxy)met...	10.523	93	93963	21.771	ng/ul#	96
29) 2,4-Dichlorophenol	10.770	162	61377	22.160	ng/ul	93
30) Naphthalene	11.193	128	218506	22.055	ng/ul	98
32) 4-Chloroaniline	11.305	127	87436	20.770	ng/ul	94
33) Hexachlorobutadiene	11.416	225	39215	21.997	ng/ul#	72
34) Caprolactam	12.086	113	26182	25.812	ng/ul#	42
35) 4-Chloro-3-methylphenol	12.386	107	70739	24.247	ng/ul	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092623\
 Data File : BG059191.D
 Acq On : 26 Sep 2023 10:15
 Operator : MA/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Sep 26 22:38:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 25 22:33:28 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/27/2023
 Supervised By :mohammad ahmed 09/29/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.768	142	142941	21.603	ng/ul	96
37) 1-Methylnaphthalene	12.985	142	152542	22.822	ng/ul	96
39) 1,2,4,5-Tetrachloroben...	13.114	216	77211	20.391	ng/ul	97
40) Hexachlorocyclopentadiene	13.067	237	41094	17.023	ng/ul#	93
41) 2,4,6-Trichlorophenol	13.355	196	50407	20.386	ng/ul	86
42) 2,4,5-Trichlorophenol	13.432	196	55865	20.887	ng/ul	95
43) 1,1'-Biphenyl	13.761	154	217209	21.069	ng/ul	97
44) 2-Chloronaphthalene	13.814	162	161581	20.526	ng/ul	96
45) 2-Nitroaniline	14.031	65	53490	21.436	ng/ul	91
47) Dimethylphthalate	14.366	163	228712	22.542	ng/ul	99
48) 2,6-Dinitrotoluene	14.507	165	42761	22.807	ng/ul#	90
50) Acenaphthylene	14.654	152	279177	21.610	ng/ul	98
51) 3-Nitroaniline	14.854	138	39360	19.026	ng/ul	96
52) Acenaphthene	14.989	153	186671	21.307	ng/ul	98
53) 2,4-Dinitrophenol	15.047	184	16047	14.205	ng/ul#	80
55) 4-Nitrophenol	15.130	109	17531	14.636	ng/ul#	75
56) Dibenzofuran	15.318	168	252725	21.748	ng/ul	97
57) 2,4-Dinitrotoluene	15.288	165	62243	23.027	ng/ul#	99
58) 2,3,4,6-Tetrachlorophenol	15.529	232	55783	22.532	ng/ul#	88
59) Diethylphthalate	15.711	149	228911	22.897	ng/ul	98
61) Fluorene	15.970	166	217641	21.935	ng/ul	97
62) 4-Chlorophenyl-phenyle...	15.946	204	110861	22.512	ng/ul	94
63) 4-Nitroaniline	16.011	138	26882m	14.018	ng/ul	
66) 4,6-Dinitro-2-methylph...	16.034	198	38830	19.950	ng/ul#	98
67) N-Nitrosodiphenylamine	16.170	169	180595	21.356	ng/ul	94
68) 4-Bromophenyl-phenylether	16.851	248	69747	22.569	ng/ul	93
69) Hexachlorobenzene	16.939	284	79519	22.465	ng/ul	95
70) Atrazine	17.104	200	81857	24.327	ng/ul	97
71) Pentachlorophenol	17.298	266	38767m	16.861	ng/ul	
72) Phenanthrene	17.715	178	390289	21.957	ng/ul	99
74) Anthracene	17.809	178	396925	21.854	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.720	216	83011	19.603	ng/uL	92
76) Pentachlorobenzene	15.212	250	80994	20.678	ng/uL	97
77) Carbazole	18.085	167	303531	20.102	ng/ul	97
78) Di-n-butylphthalate	18.590	149	434919	23.381	ng/ul	99
80) Fluoranthene	19.713	202	454627	22.899	ng/ul	99
82) Pyrene	20.077	202	464767	22.443	ng/ul	98
83) Butylbenzylphthalate	20.929	149	181634	24.282	ng/ul	96
84) 3,3'-Dichlorobenzidine	21.886	252	108971	16.544	ng/ul#	94
85) Benzo(a)anthracene	21.975	228	421528	21.301	ng/ul	96
86) Bis(2-ethylhexyl)phtha...	21.804	149	282272	25.221	ng/ul	98
87) Chrysene	22.045	228	401438	21.752	ng/ul	99
89) Di-n-octyl phthalate	23.120	149	465355	23.962	ng/ul	100
90) Benzo(b)fluoranthene	24.389	252	432976	22.368	ng/ul	99
91) Benzo(k)fluoranthene	24.466	252	418121	20.834	ng/ul	97
93) Benzo(a)pyrene	25.371	252	388132	20.864	ng/ul	98
94) Indeno(1,2,3-cd)pyrene	29.654	276	405264m	19.155	ng/ul	
95) Dibenzo(a,h)anthracene	29.724	278	282848m	16.636	ng/ul	
96) Benzo(g,h,i)perylene	30.958	276	352140m	20.662	ng/ul	

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

Instrument :
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LabSampleId :
SSTDCCC020

Manual IntegrationsAPPROVED

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Manual IntegrationsAPPROVED

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