

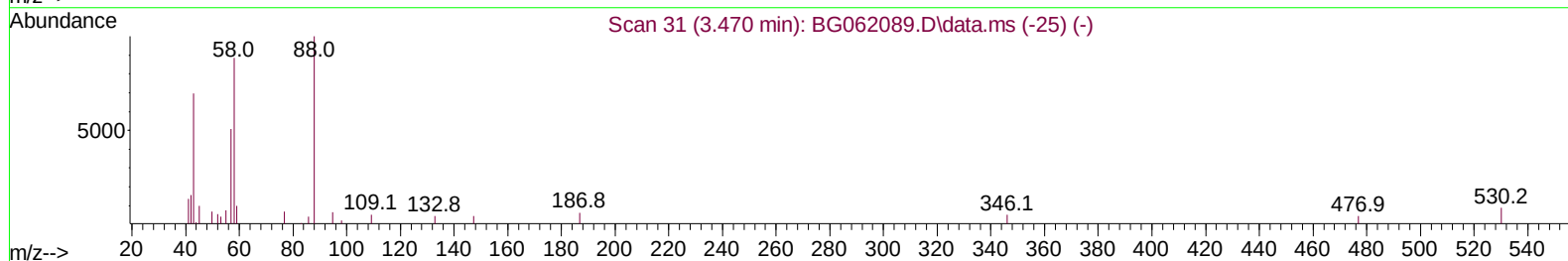
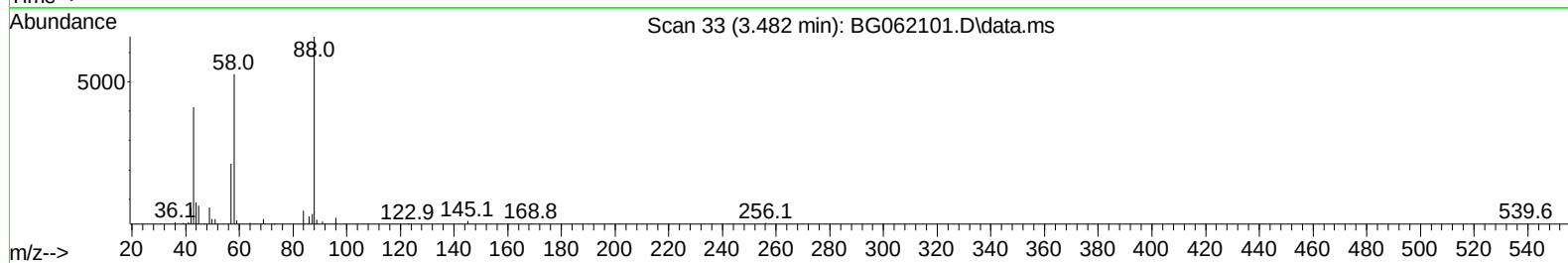
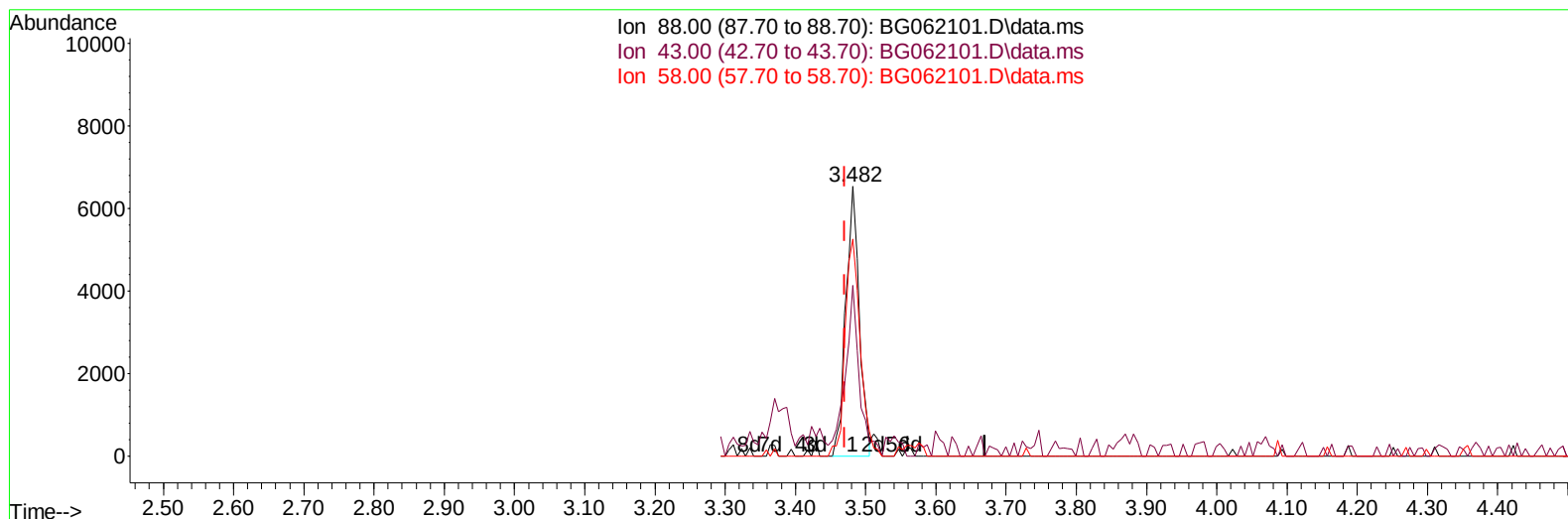
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 16 12:23:16 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/17/2024
Supervised By :mohammad ahmed 07/18/2024



TIC: BG062101.D\data.ms

(2) 1,4-Dioxane

3.482min (+ 0.011) 6.02 ng/uL

response 8730

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	49.30	63.25#
58.00	97.60	80.52
0.00	0.00	0.00

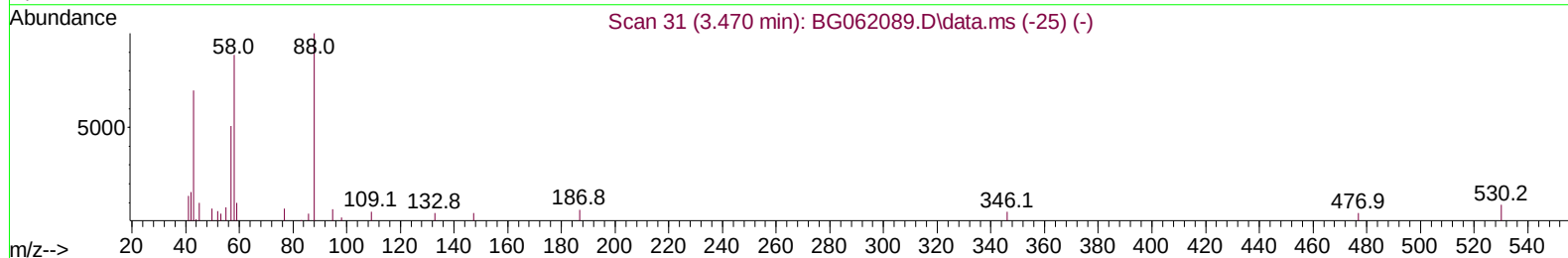
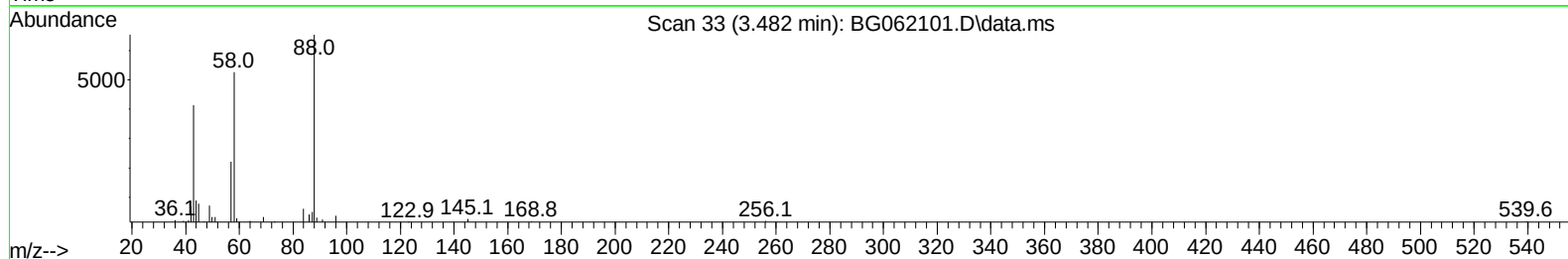
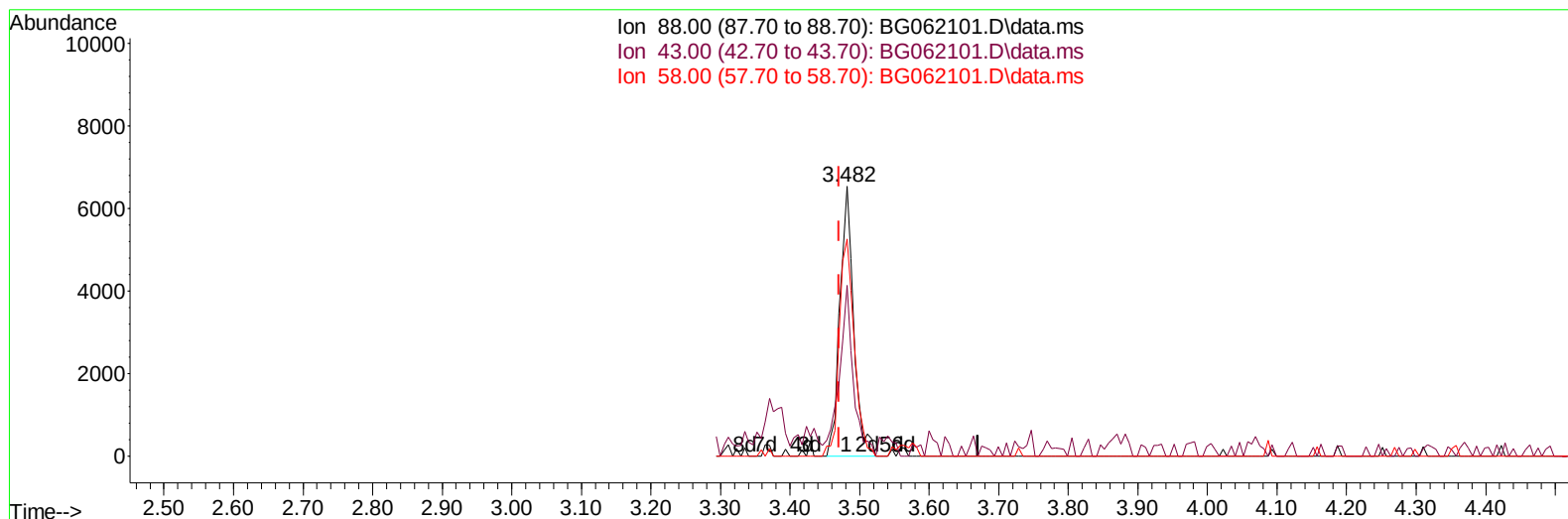
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
 Data File : BG062101.D
 Acq On : 16 Jul 2024 19:41
 Operator : MA/JU
 Sample : PB162022BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
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 Supervised By :mohammad ahmed 07/18/2024



TIC: BG062101.D\data.ms

(2) 1,4-Dioxane

3.482min (+ 0.011) 6.25 ng/uL m

response 9062

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	49.30	63.25#
58.00	97.60	80.52
0.00	0.00	0.00

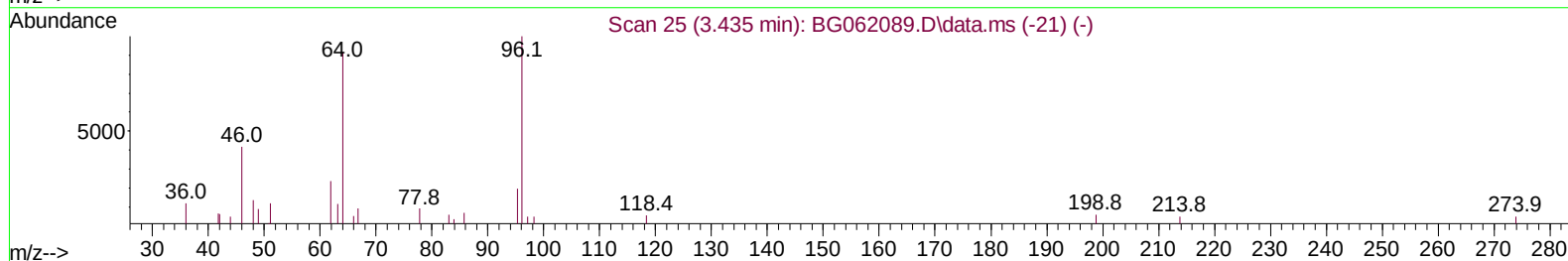
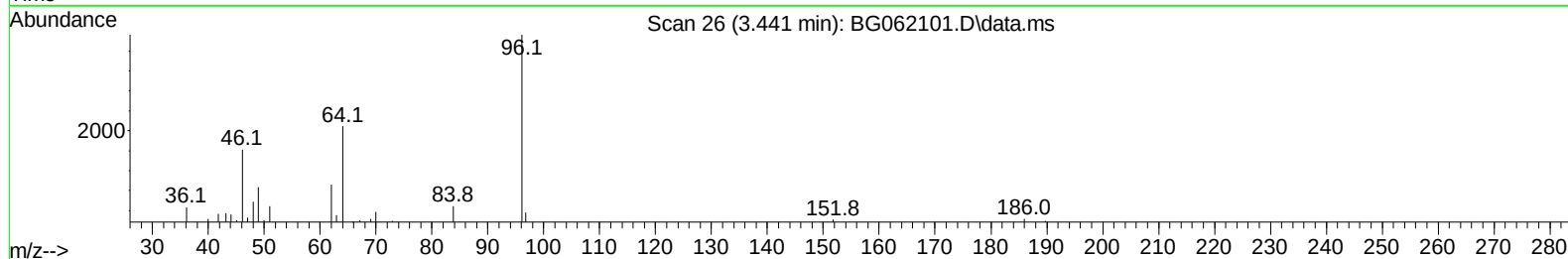
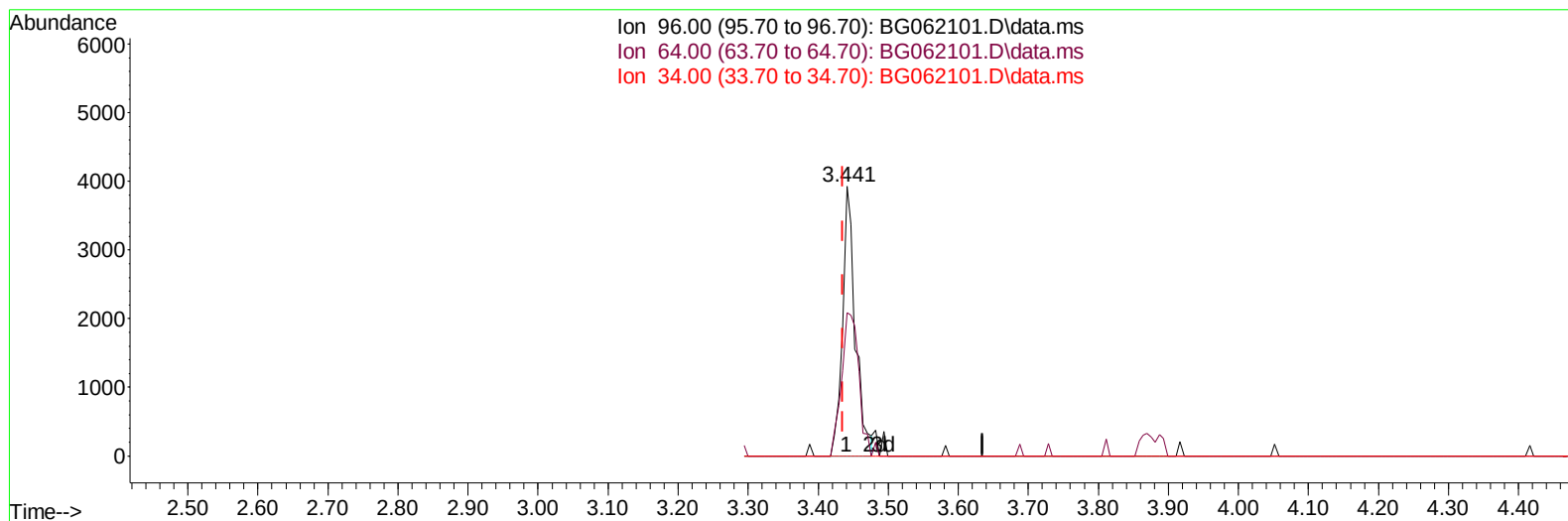
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
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QLast Update : Tue Jul 16 12:23:16 2024
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TIC: BG062101.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.441min (+ 0.006) 3.85 ng/uL

response 5063

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	104.80	53.22#
34.00	0.00	0.00
0.00	0.00	0.00

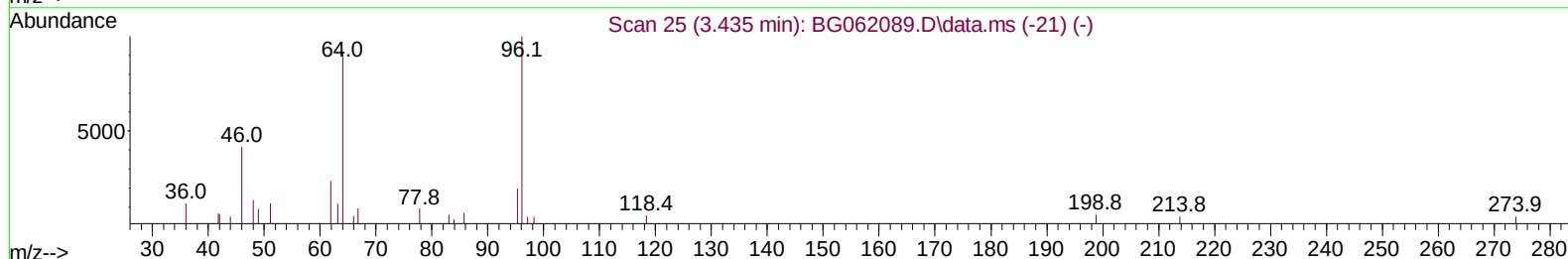
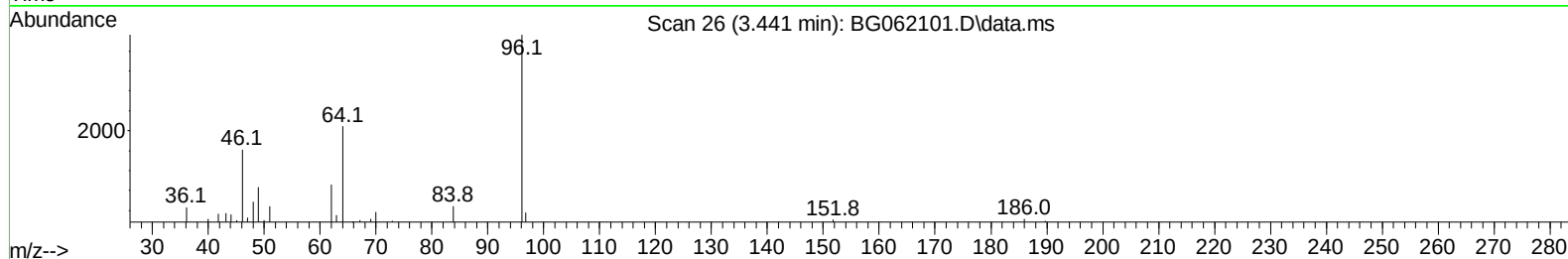
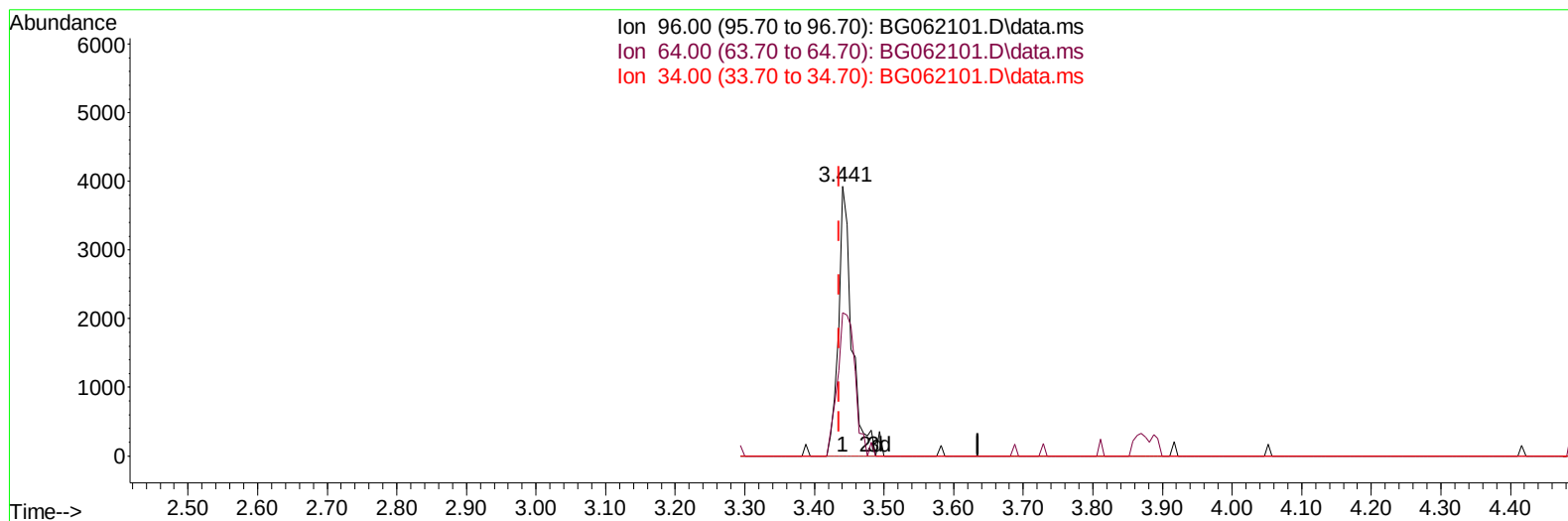
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
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QLast Update : Tue Jul 16 12:23:16 2024
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TIC: BG062101.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.441min (+ 0.006) 3.95 ng/uL m

response 5195

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	104.80	53.22#
34.00	0.00	0.00
0.00	0.00	0.00

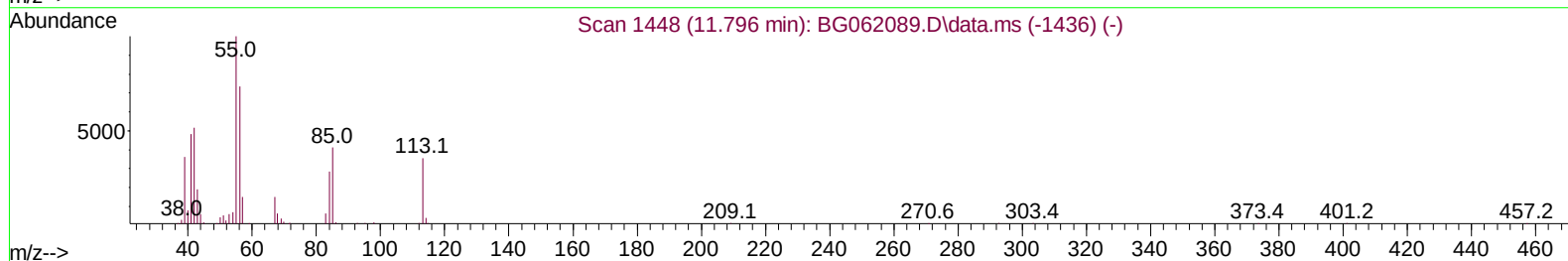
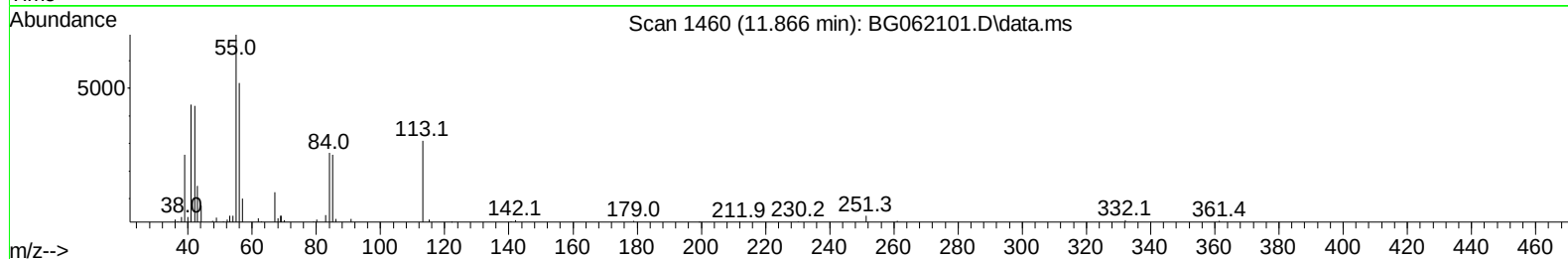
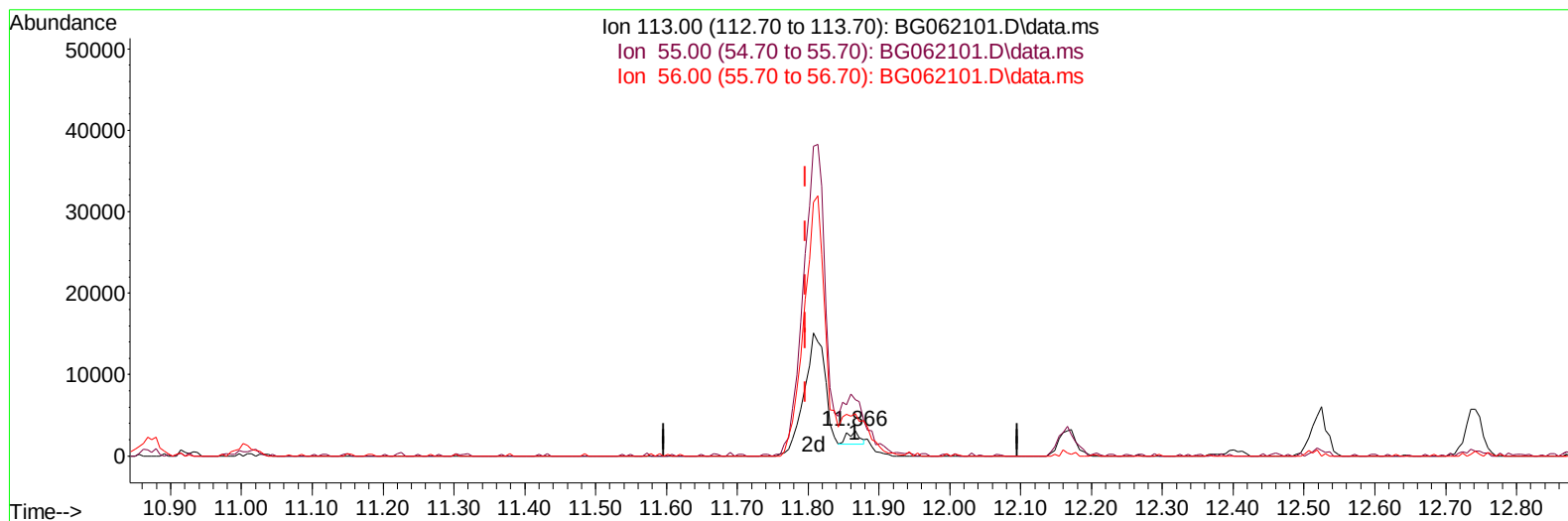
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
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TIC: BG062101.D\data.ms

(34) Caprolactam

11.866min (+ 0.070) 1.48 ng/ul

response 1980

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	230.70	223.37
56.00	168.80	167.29
0.00	0.00	0.00

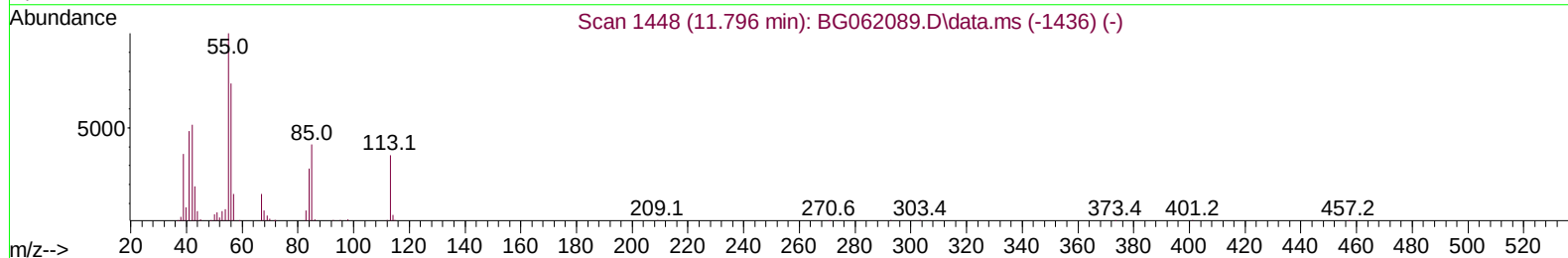
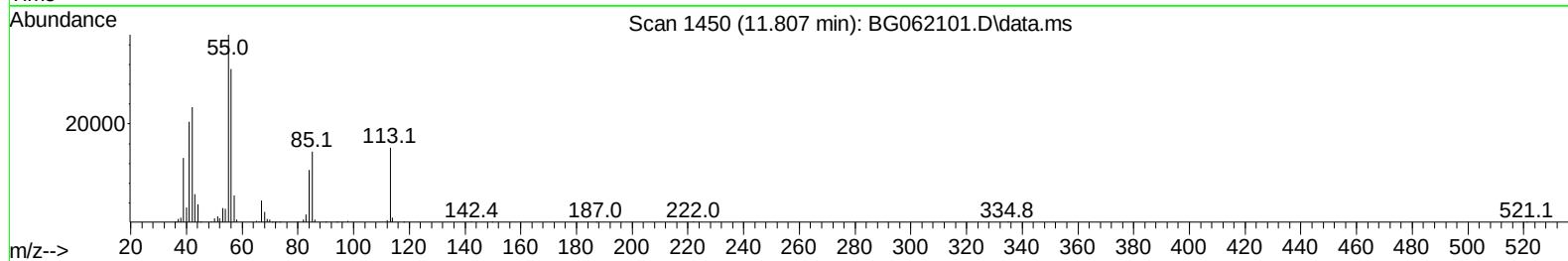
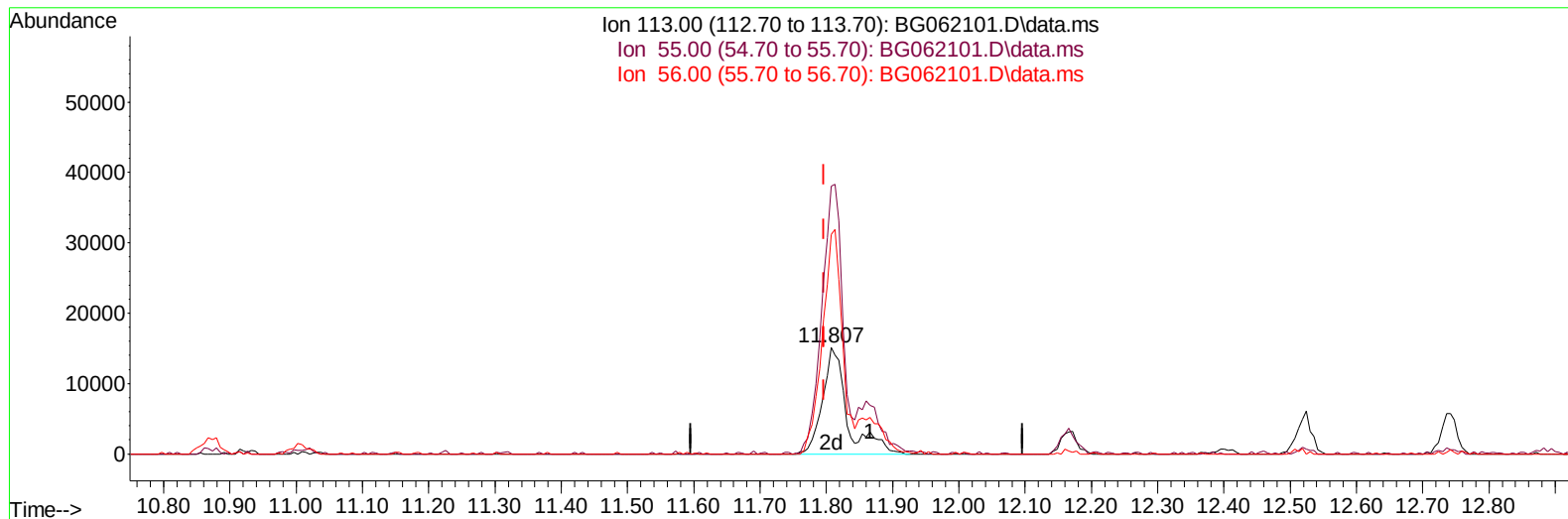
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
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QLast Update : Tue Jul 16 12:23:16 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/17/2024
Supervised By :mohammad ahmed 07/18/2024



TIC: BG062101.D\data.ms

(34) Caprolactam

11.807min (+ 0.011) 29.13 ng/ul m

response 39046

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	230.70	251.32
56.00	168.80	205.85#
0.00	0.00	0.00

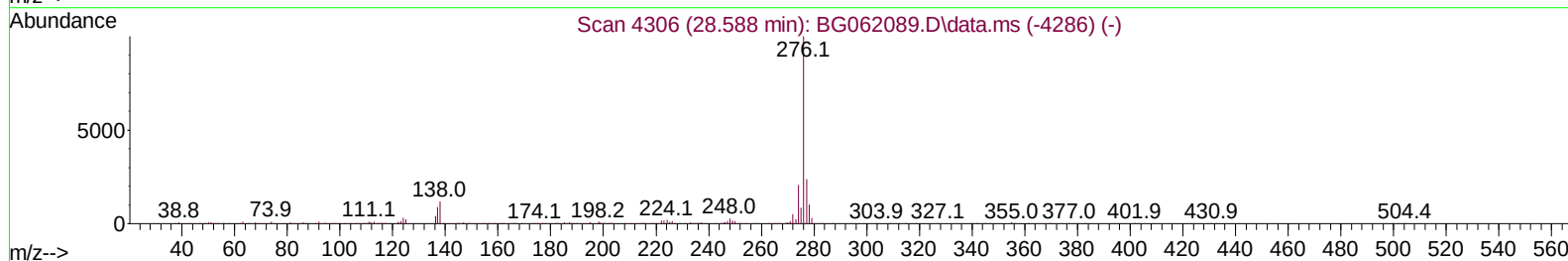
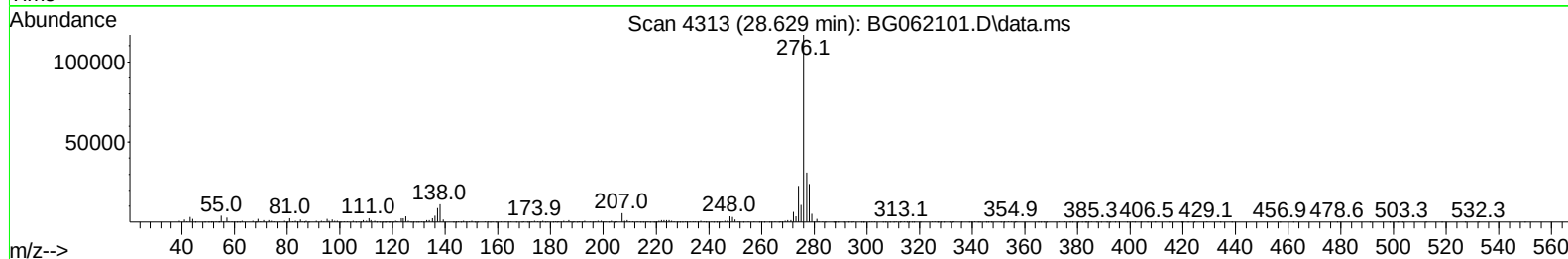
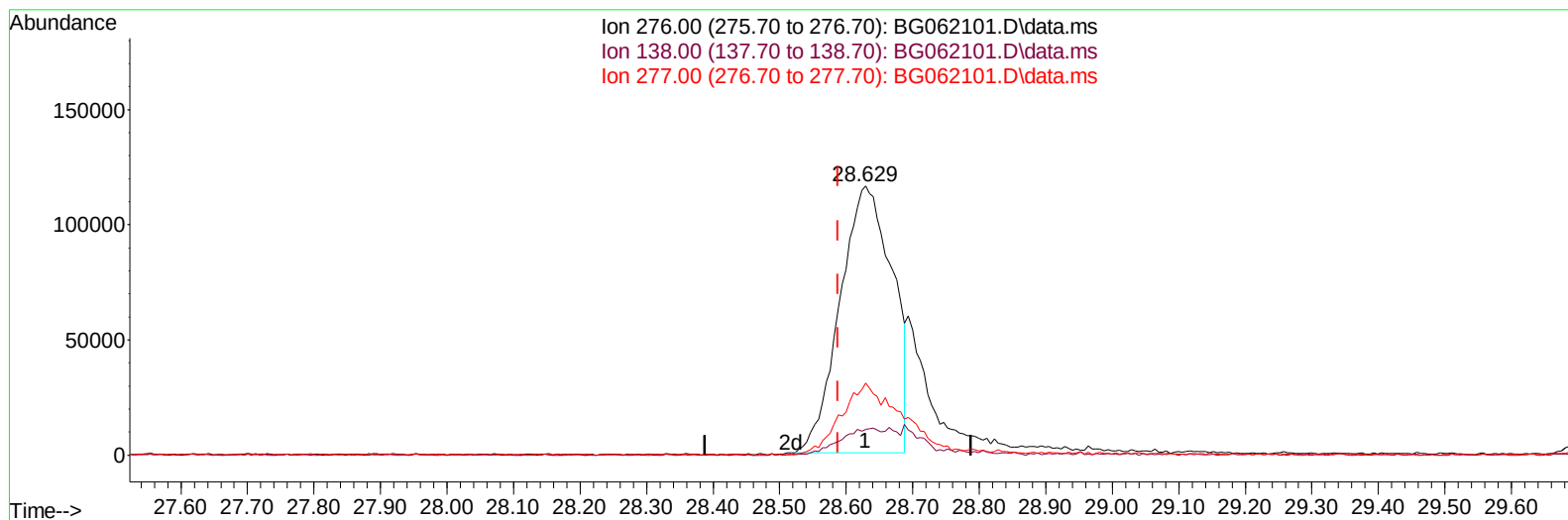
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

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TIC: BG062101.D\data.ms

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

28.629min (+ 0.041) 21.36 ng/ul

response 630962

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	14.90	9.60#
277.00	24.10	26.68
0.00	0.00	0.00

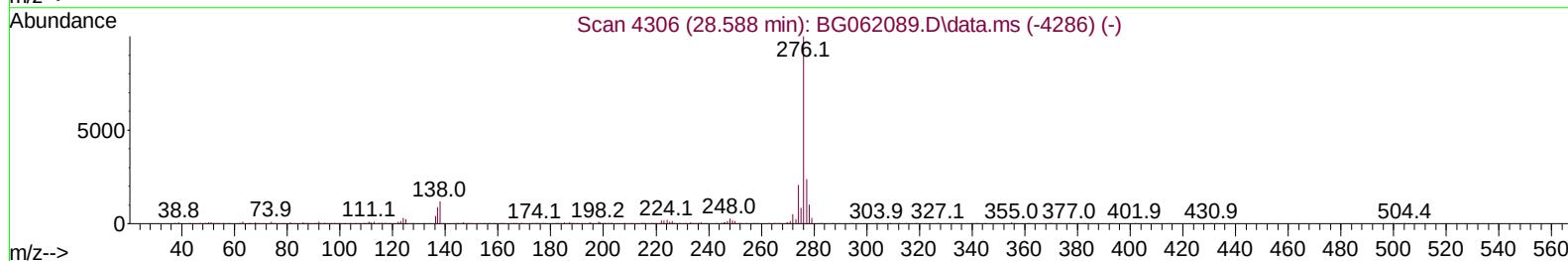
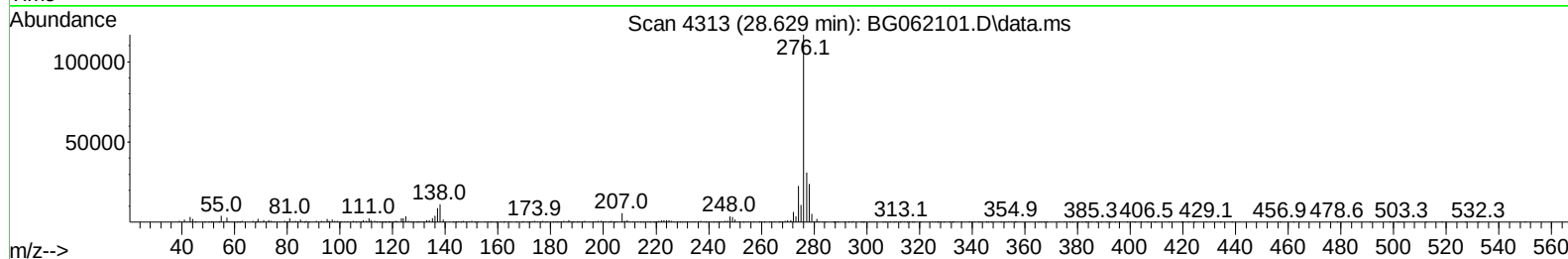
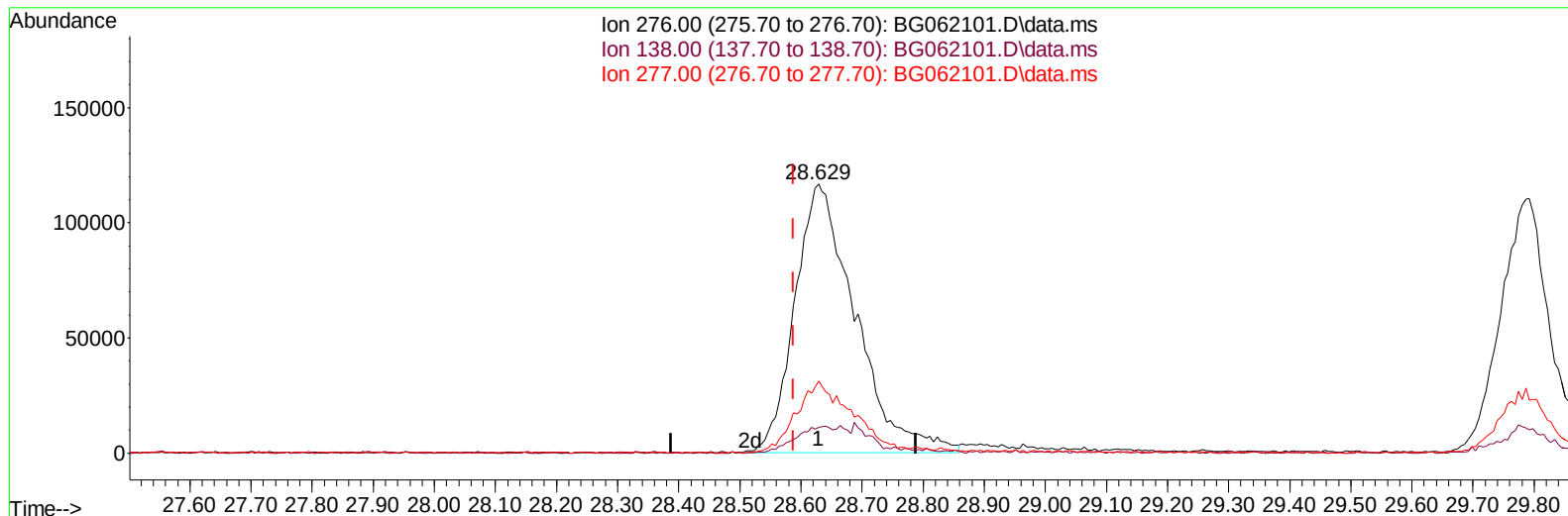
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 16 12:23:16 2024
Response via : Initial Calibration

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Supervised By :mohammad ahmed 07/18/2024



TIC: BG062101.D\data.ms

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

28.629min (+ 0.041) 27.29 ng/ul m

response 806231

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	14.90	9.60#
277.00	24.10	26.68
0.00	0.00	0.00

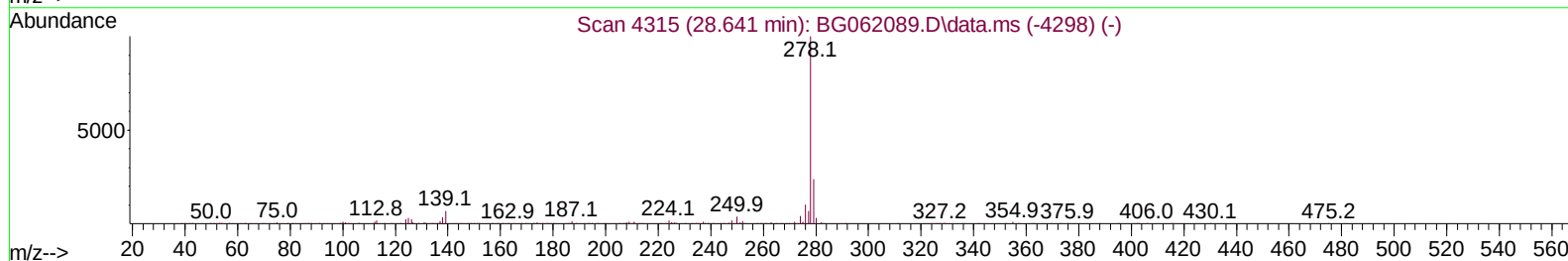
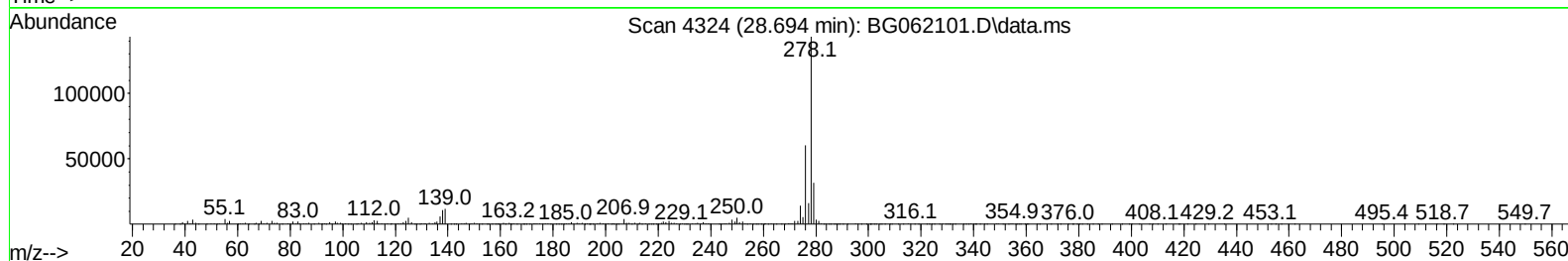
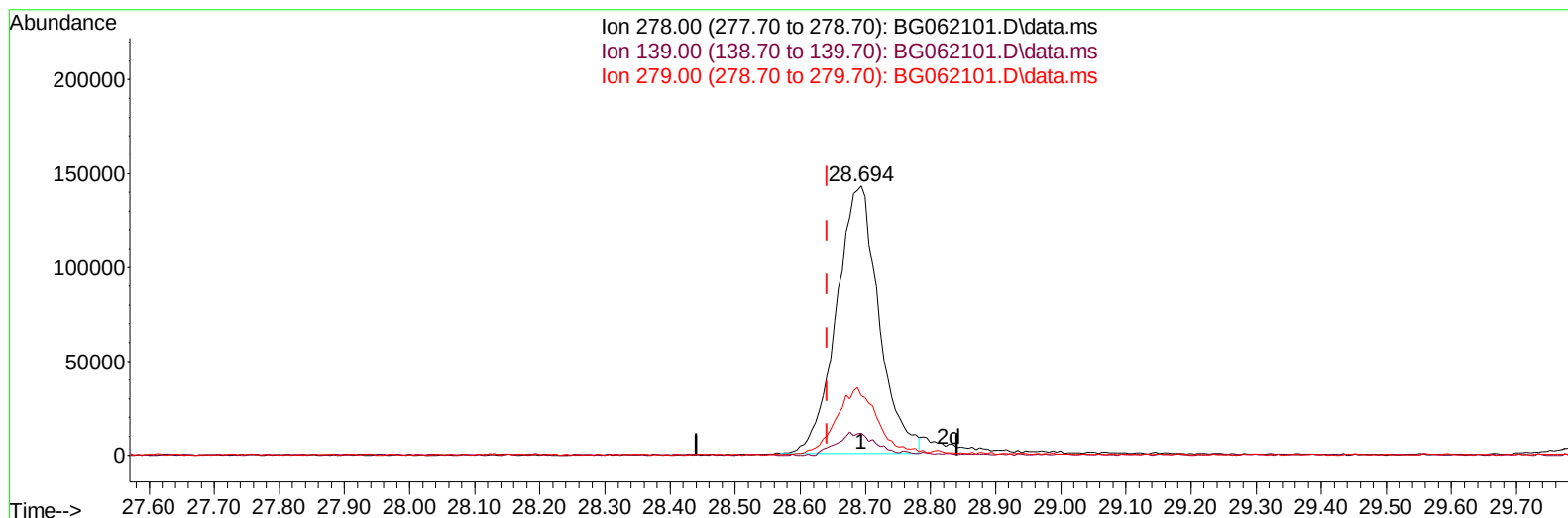
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
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TIC: BG062101.D\data.ms

(95) Dibenzo(a,h)anthracene

28.694min (+ 0.053) 26.42 ng/ul

response 644620

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	9.50	8.16
279.00	23.50	22.03
0.00	0.00	0.00

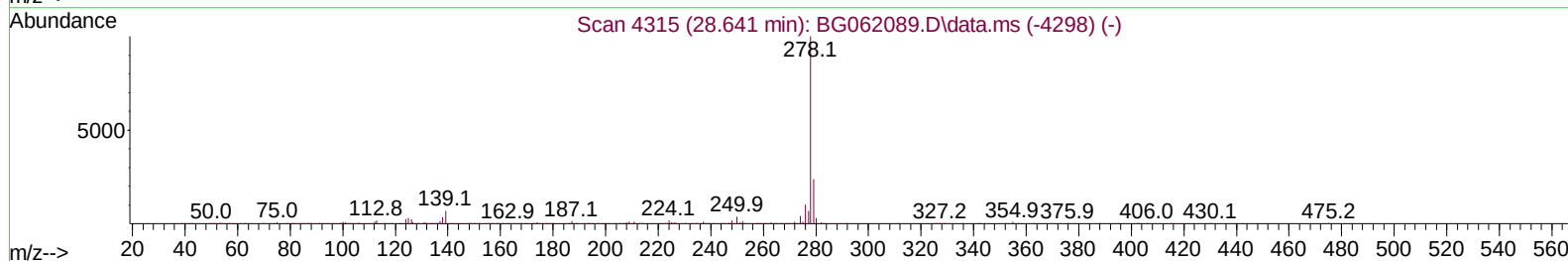
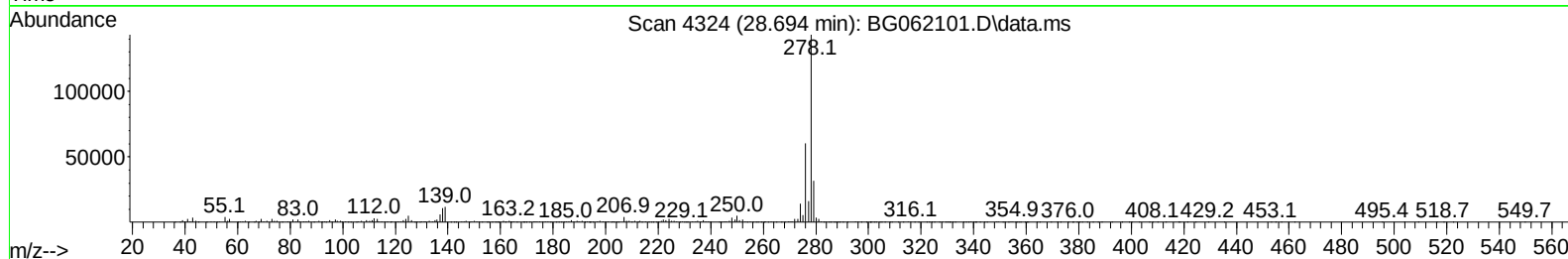
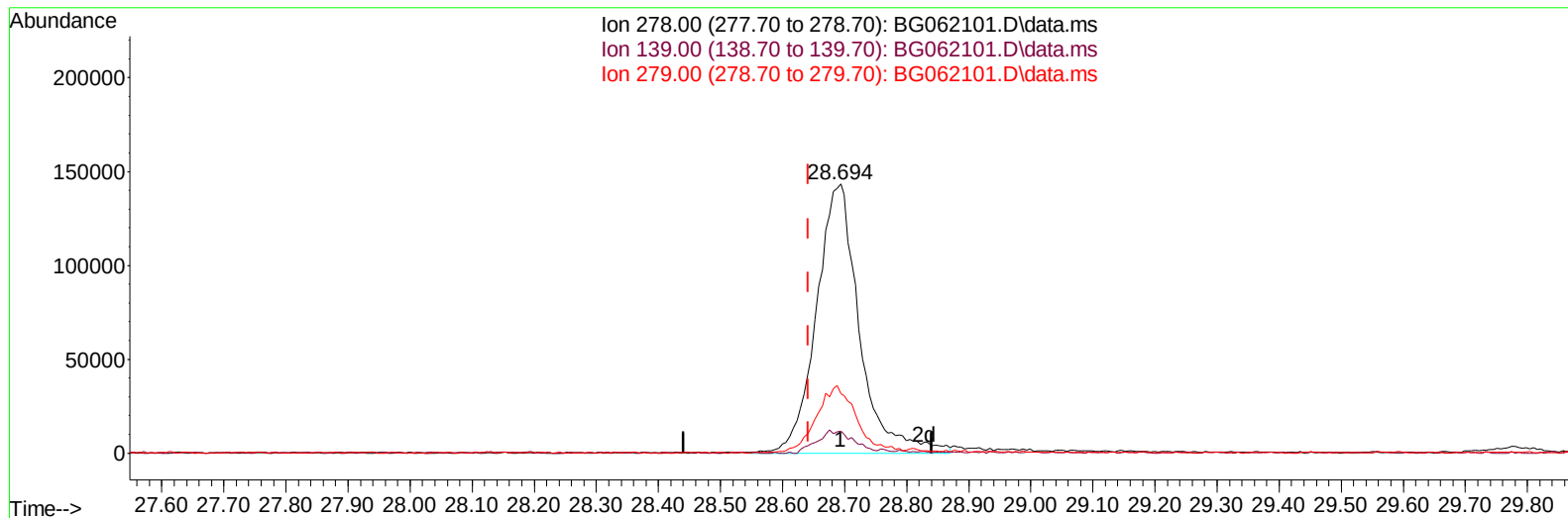
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
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TIC: BG062101.D\data.ms

(95) Dibenzo(a,h)anthracene

28.694min (+ 0.053) 28.21 ng/ul m

response 688255

Ion	Exp%	Act%
278.00	100.00	100.00
139.00	9.50	8.16
279.00	23.50	22.03
0.00	0.00	0.00

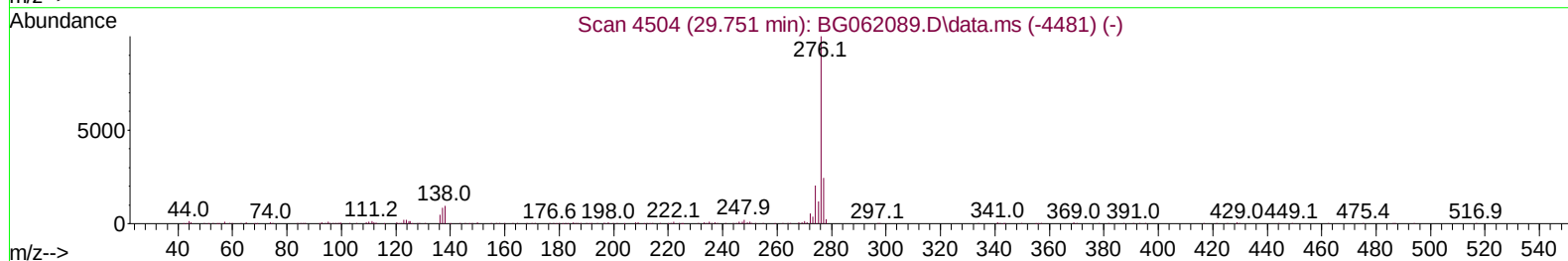
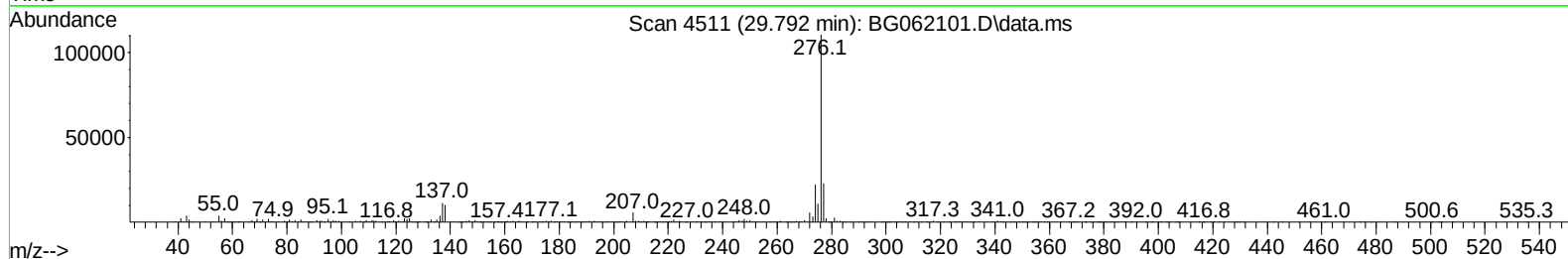
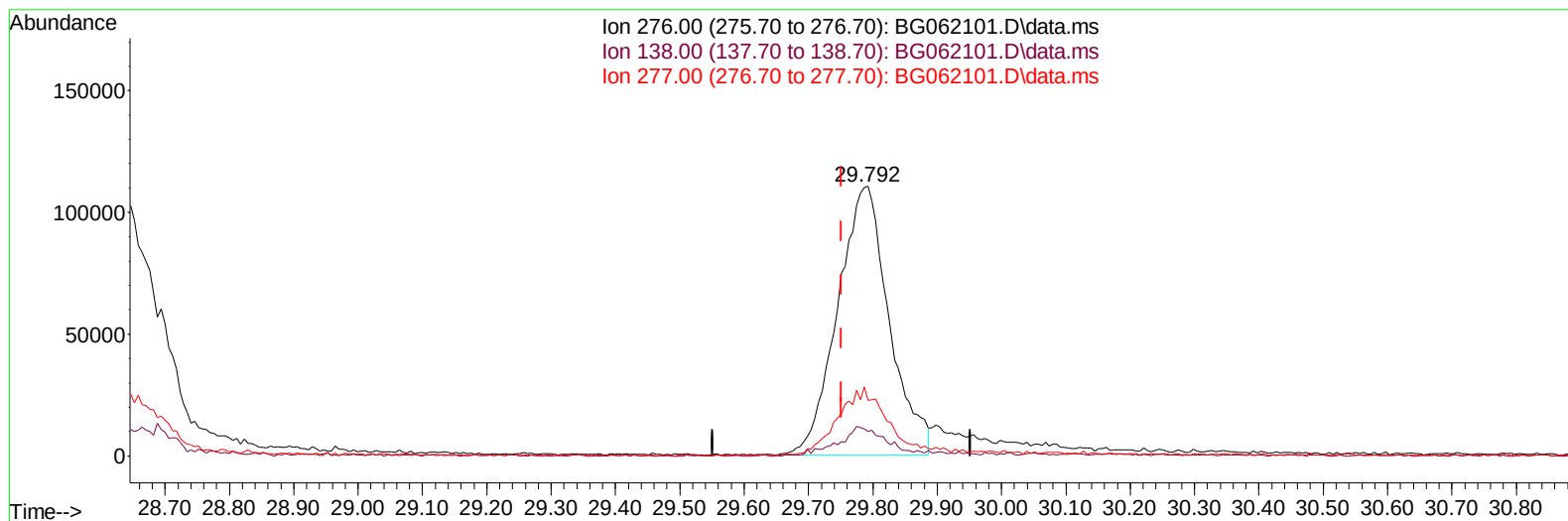
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
 Data File : BG062101.D
 Acq On : 16 Jul 2024 19:41
 Operator : MA/JU
 Sample : PB162022BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
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 QLast Update : Tue Jul 16 12:23:16 2024
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TIC: BG062101.D\data.ms

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

29.792min (+ 0.041) 26.07 ng/ul

response 612620

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	13.90	9.21#
277.00	23.50	20.76
0.00	0.00	0.00

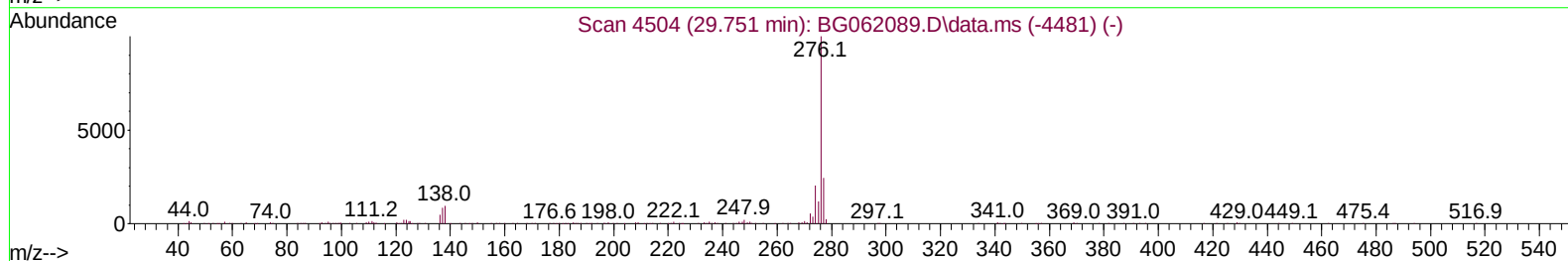
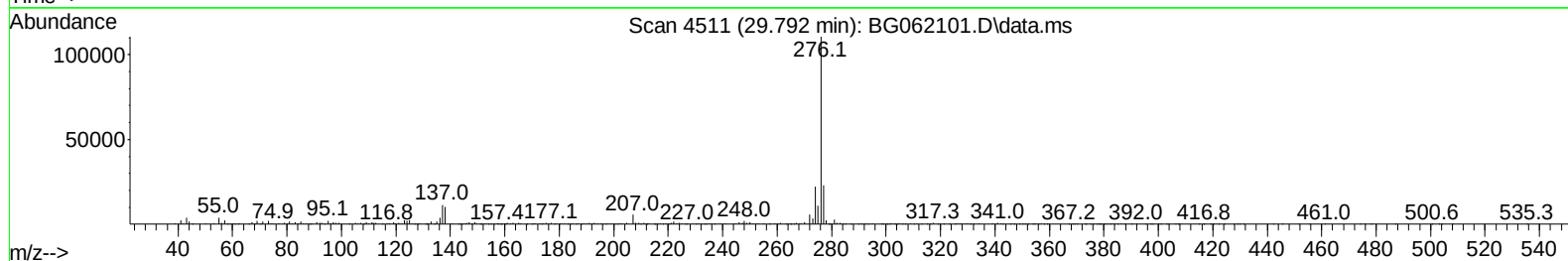
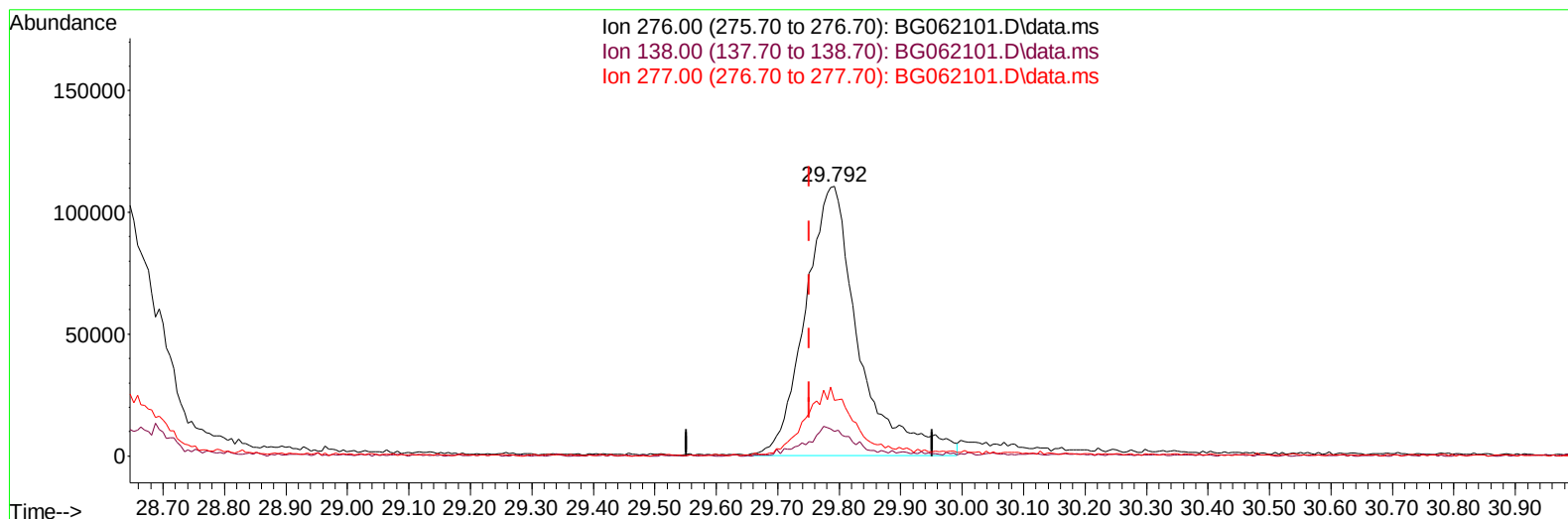
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
Data File : BG062101.D
Acq On : 16 Jul 2024 19:41
Operator : MA/JU
Sample : PB162022BS
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 00:47:24 2024
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Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 16 12:23:16 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/17/2024
Supervised By :mohammad ahmed 07/18/2024



TIC: BG062101.D\data.ms

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

29.792min (+ 0.041) 28.50 ng/ul m

response 669763

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	13.90	9.21#
277.00	23.50	20.76
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
 Data File : BG062101.D
 Acq On : 16 Jul 2024 19:41
 Operator : MA/JU
 Sample : PB162022BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 01:46:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 16 12:23:16 2024
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Reviewed By :Jagrut Upadhyay 07/17/2024
 Supervised By :mohammad ahmed 07/18/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.053	152	49365	20.000	ng/ul	0.00
20) Naphthalene-d8	10.873	136	217451	20.000	ng/ul	# 0.01
38) Acenaphthene-d10	14.692	164	164472	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.436	188	354943	20.000	ng/ul	# 0.01
79) Chrysene-d12	21.702	240	315596	20.000	ng/ul	# 0.01
88) Perylene-d12	24.933	264	387493	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.441	96	5195m	3.955	ng/uL	0.00
4) Pyridine-d5	3.864	84	57583	14.256	ng/ul	0.00
7) Phenol-d5	7.230	99	124089	22.970	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.377	67	69865	20.422	ng/ul	0.00
11) 2-Chlorophenol-d4	7.589	132	100497	27.119	ng/ul	0.00
15) 4-Methylphenol-d8	8.782	113	92665	21.785	ng/ul	0.02
21) Nitrobenzene-d5	9.228	128	52637	31.828	ng/ul	0.00
24) 2-Nitrophenol-d4	9.951	143	67346	35.659	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.503	165	127622	35.034	ng/ul	0.02
31) 4-Chloroaniline-d4	11.008	131	138063	26.105	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	414924	30.686	ng/ul	0.00
49) Acenaphthylene-d8	14.387	160	434920	30.760	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	56925	26.362	ng/ul	0.01
60) Fluorene-d10	15.679	176	309148	27.158	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.809	200	61622	33.057	ng/ul	0.01
73) Anthracene-d10	17.536	188	435889	26.303	ng/ul	0.01
81) Pyrene-d10	19.816	212	540722	29.129	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.710	264	612838	31.889	ng/ul	0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.482	88	9062m	6.250	ng/uL	
5) Pyridine	3.881	79	55768	13.463	ng/ul#	87
6) Benzaldehyde	7.195	77	74399	26.021	ng/ul	85
8) Phenol	7.260	94	124976	22.595	ng/ul#	81
10) Bis(2-Chloroethyl)ether	7.471	93	85173	20.045	ng/ul	97
12) 2-Chlorophenol	7.624	128	100681	26.668	ng/ul	92
13) 2-Methylphenol	8.511	108	94694	23.987	ng/ul	91
14) 2,2'-oxybis(1-Chloropr...	8.576	45	170952	18.416	ng/ul#	96
16) Acetophenone	8.881	105	156590	22.287	ng/ul	92
17) N-Nitroso-di-n-propyla...	8.864	70	80607	20.492	ng/ul#	83
18) 4-Methylphenol	8.846	108	97969	22.160	ng/ul	100
19) Hexachloroethane	9.134	117	42854	26.166	ng/ul#	73
22) Nitrobenzene	9.269	77	146754	31.261	ng/ul	97
23) Isophorone	9.792	82	272164	25.662	ng/ul#	95
25) 2-Nitrophenol	9.980	139	63635	33.040	ng/ul#	92
26) 2,4-Dimethylphenol	10.045	107	100099	21.716	ng/ul	95
27) Bis(2-Chloroethoxy)met...	10.268	93	139574	23.898	ng/ul	96
29) 2,4-Dichlorophenol	10.527	162	116222	33.965	ng/ul	93
30) Naphthalene	10.920	128	316929	26.533	ng/ul	97
32) 4-Chloroaniline	11.032	127	124207	23.935	ng/ul	99
33) Hexachlorobutadiene	11.191	225	81070	31.484	ng/ul	92
34) Caprolactam	11.807	113	39046m	29.125	ng/ul	
35) 4-Chloro-3-methylphenol	12.166	107	144774	31.767	ng/ul	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071624\
 Data File : BG062101.D
 Acq On : 16 Jul 2024 19:41
 Operator : MA/JU
 Sample : PB162022BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLCS022

Manual IntegrationsAPPROVED

Quant Time: Jul 17 01:46:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 16 12:23:16 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/17/2024
 Supervised By :mohammad ahmed 07/18/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.524	142	264067	30.515	ng/ul	96
37) 1-Methylnaphthalene	12.742	142	271695	30.212	ng/ul	94
39) 1,2,4,5-Tetrachloroben...	12.883	216	166198	33.649	ng/ul	95
40) Hexachlorocyclopentadiene	12.853	237	36401	11.071	ng/ul#	95
41) 2,4,6-Trichlorophenol	13.135	196	85804	26.155	ng/ul	99
42) 2,4,5-Trichlorophenol	13.212	196	92177	25.578	ng/ul	91
43) 1,1'-Biphenyl	13.523	154	331775	26.181	ng/ul#	96
44) 2-Chloronaphthalene	13.570	162	275938	29.013	ng/ul	93
45) 2-Nitroaniline	13.782	65	120883	31.739	ng/ul	93
47) Dimethylphthalate	14.140	163	366724	28.197	ng/ul	98
48) 2,6-Dinitrotoluene	14.269	165	79132	32.072	ng/ul	95
50) Acenaphthylene	14.416	152	437619	28.201	ng/ul	99
51) 3-Nitroaniline	14.598	138	75120	31.278	ng/ul	92
52) Acenaphthene	14.751	153	303331	28.308	ng/ul	94
53) 2,4-Dinitrophenol	14.810	184	34954	27.116	ng/ul	90
55) 4-Nitrophenol	14.933	109	80163	32.526	ng/ul#	84
56) Dibenzofuran	15.086	168	418517	27.635	ng/ul	91
57) 2,4-Dinitrotoluene	15.057	165	117856	32.571	ng/ul#	86
58) 2,3,4,6-Tetrachlorophenol	15.315	232	73119	22.470	ng/ul#	93
59) Diethylphthalate	15.491	149	340896	24.323	ng/ul	94
61) Fluorene	15.738	166	333288	25.878	ng/ul	98
62) 4-Chlorophenyl-phenyle...	15.721	204	174660	26.998	ng/ul	95
63) 4-Nitroaniline	15.768	138	73543	30.178	ng/ul#	71
66) 4,6-Dinitro-2-methylph...	15.820	198	64395	32.099	ng/ul#	85
67) N-Nitrosodiphenylamine	15.938	169	303772	31.339	ng/ul	100
68) 4-Bromophenyl-phenylether	16.619	248	139374	34.612	ng/ul	99
69) Hexachlorobenzene	16.737	284	159092	34.359	ng/ul	97
70) Atrazine	16.884	200	56526	14.772	ng/ul	89
71) Pentachlorophenol	17.090	266	51658	20.297	ng/ul	96
72) Phenanthrene	17.477	178	480751	25.687	ng/ul	100
74) Anthracene	17.571	178	480847	24.861	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.488	216	146639	33.264	ng/ul	97
76) Pentachlorobenzene	15.004	250	179522	35.704	ng/ul	95
77) Carbazole	17.842	167	449807	27.634	ng/ul	99
78) Di-n-butylphthalate	18.376	149	621144	29.722	ng/ul	97
80) Fluoranthene	19.481	202	555834	25.302	ng/ul#	92
82) Pyrene	19.845	202	660608	29.111	ng/ul#	92
83) Butylbenzylphthalate	20.715	149	224809	23.605	ng/ul	92
84) 3,3'-Dichlorobenzidine	21.590	252	238852	31.813	ng/ul	96
85) Benzo(a)anthracene	21.678	228	672736	29.441	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.561	149	375001	27.440	ng/ul	95
87) Chrysene	21.749	228	641743	30.000	ng/ul	98
89) Di-n-octyl phthalate	22.771	149	670527	28.324	ng/ul	100
90) Benzo(b)fluoranthene	23.905	252	722442	29.592	ng/ul#	95
91) Benzo(k)fluoranthene	23.970	252	727566	29.886	ng/ul#	92
93) Benzo(a)pyrene	24.786	252	628297	27.178	ng/ul#	96
94) Indeno(1,2,3-cd)pyrene	28.629	276	806231m	27.290	ng/ul	
95) Dibenzo(a,h)anthracene	28.694	278	688255m	28.208	ng/ul	
96) Benzo(g,h,i)perylene	29.792	276	669763m	28.497	ng/ul	

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

Instrument :

BNA_G

ClientSampleId :

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