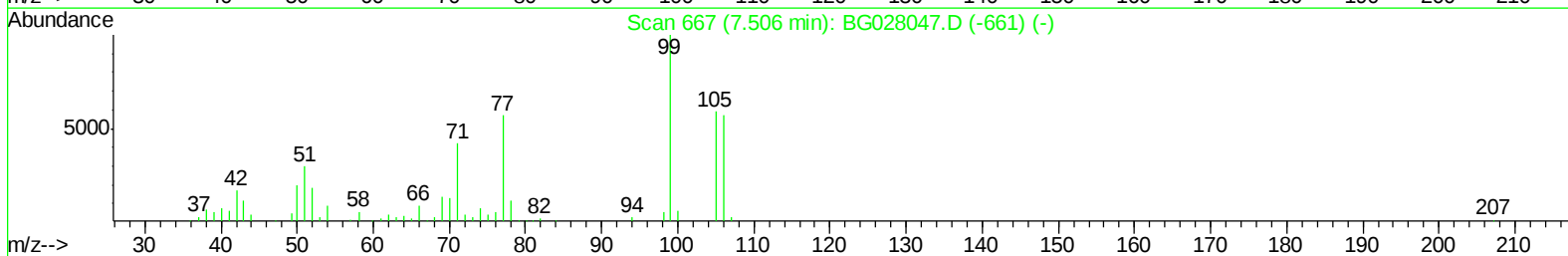
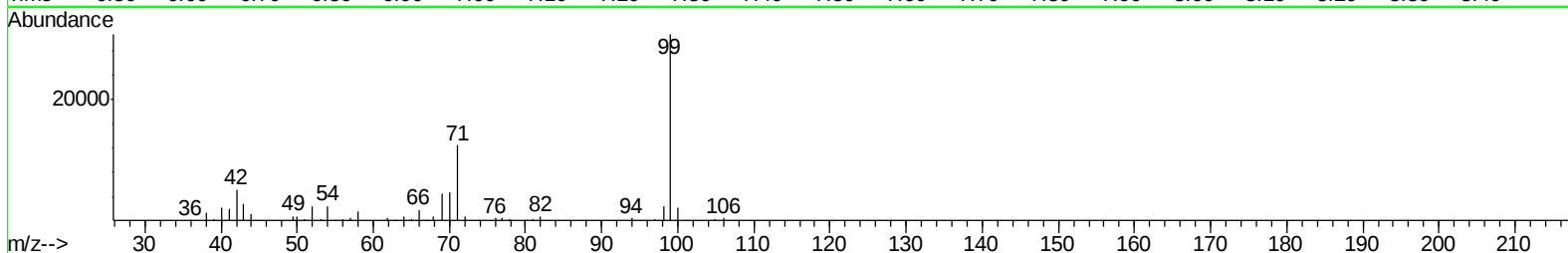
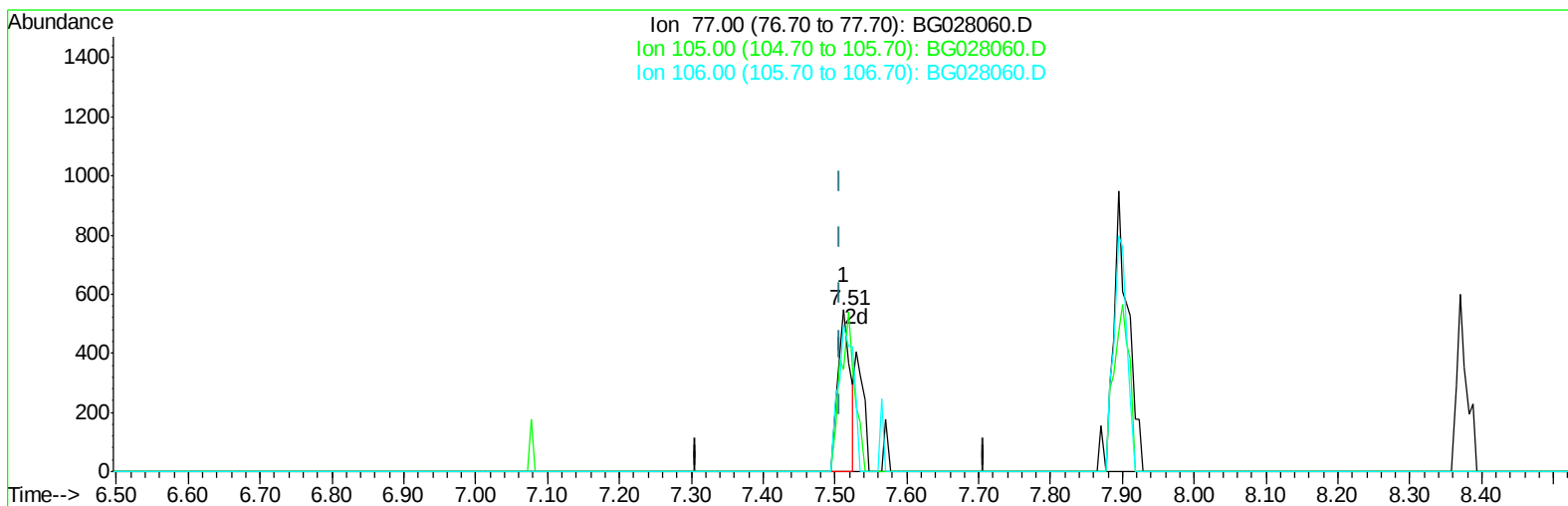


Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(4) Benzaldehyde

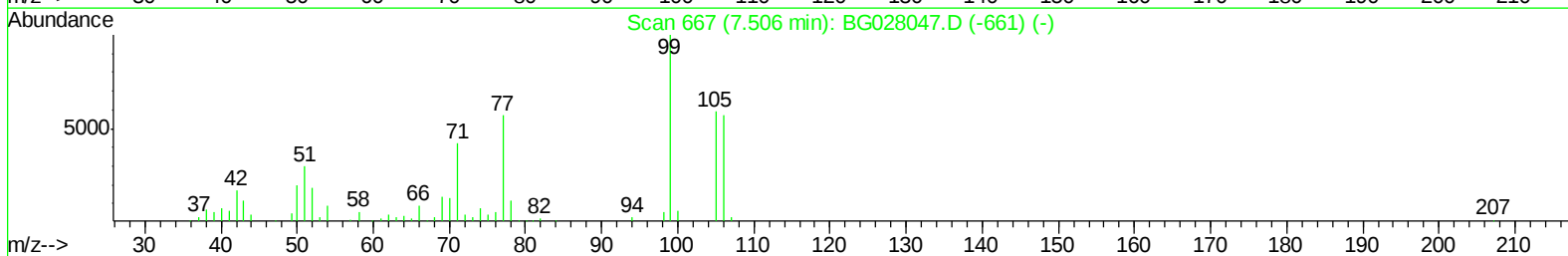
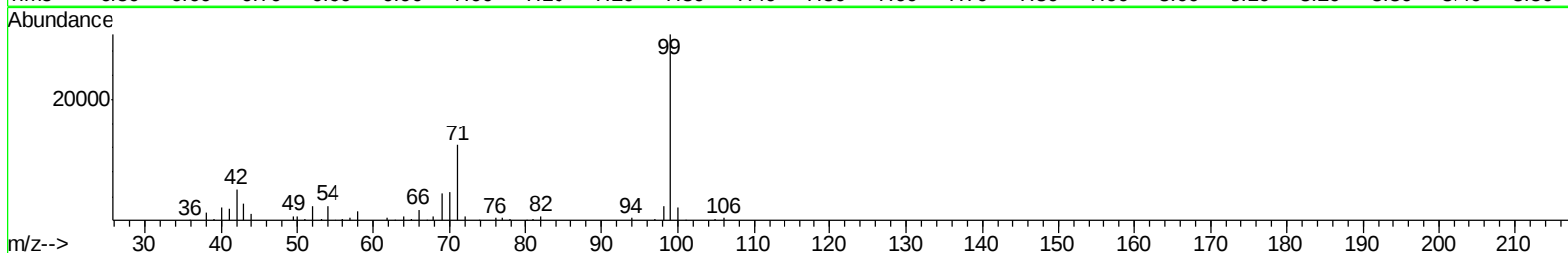
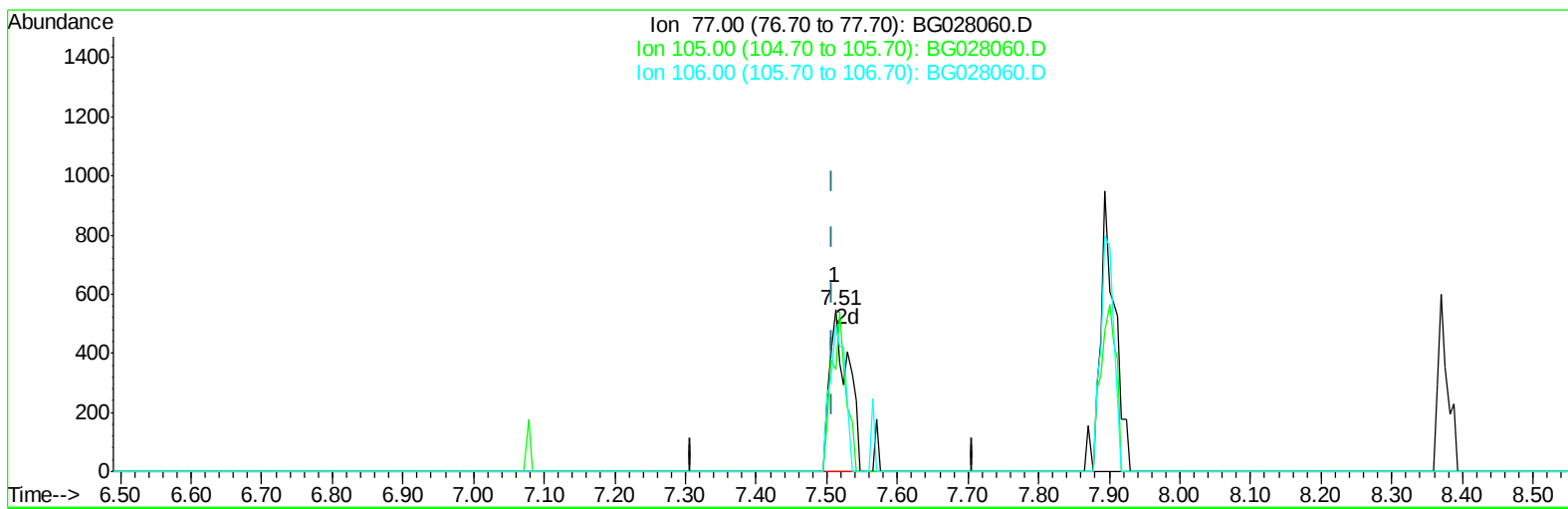
7.512min (+0.005) 0.58ng/ul

response 650

Ion	Exp%	Act%
77.00	100	100
105.00	88.80	63.00#
106.00	86.90	90.66
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(4) Benzaldehyde

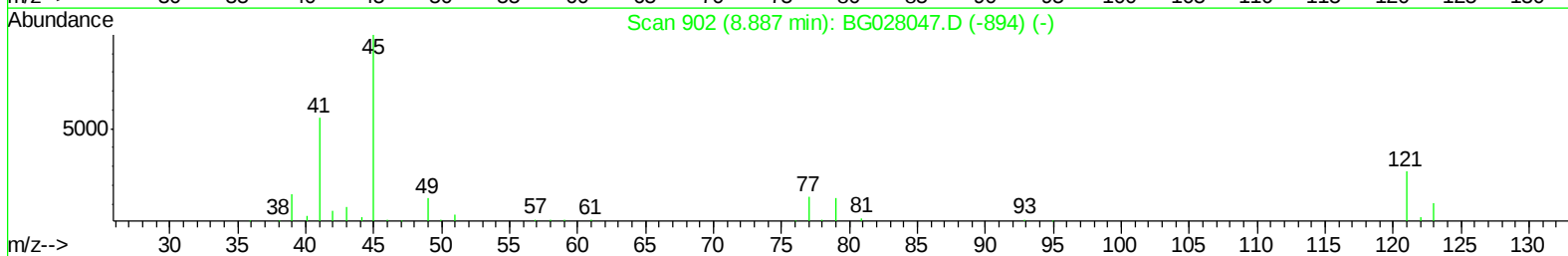
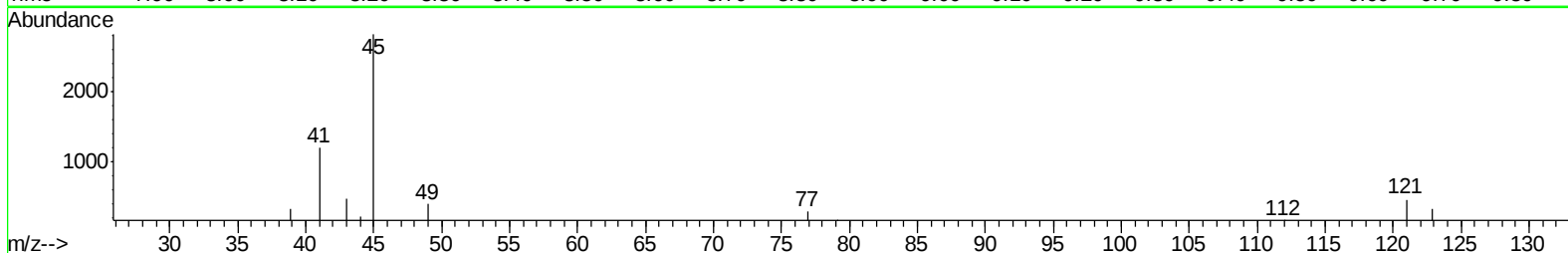
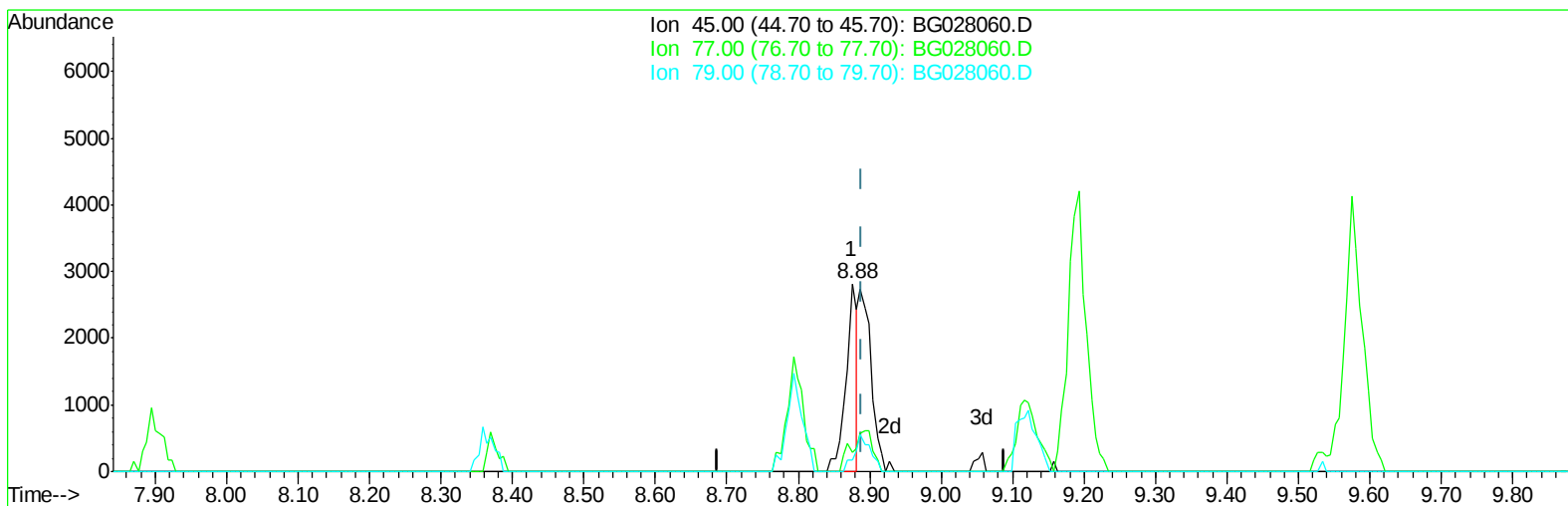
7.512min (+0.005) 0.88ng/ul m

response 994

Ion	Exp%	Act%
77.00	100	100
105.00	88.80	63.00#
106.00	86.90	90.66
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(12) 2,2'-oxybis(1-Chloropropane)

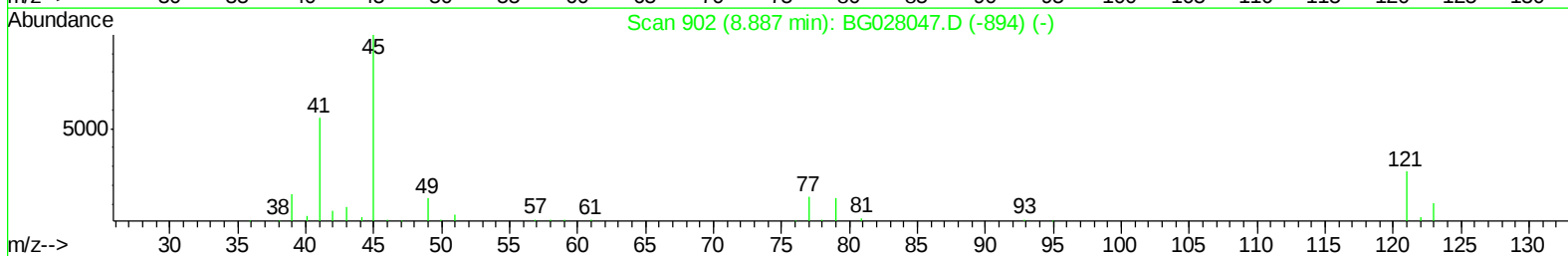
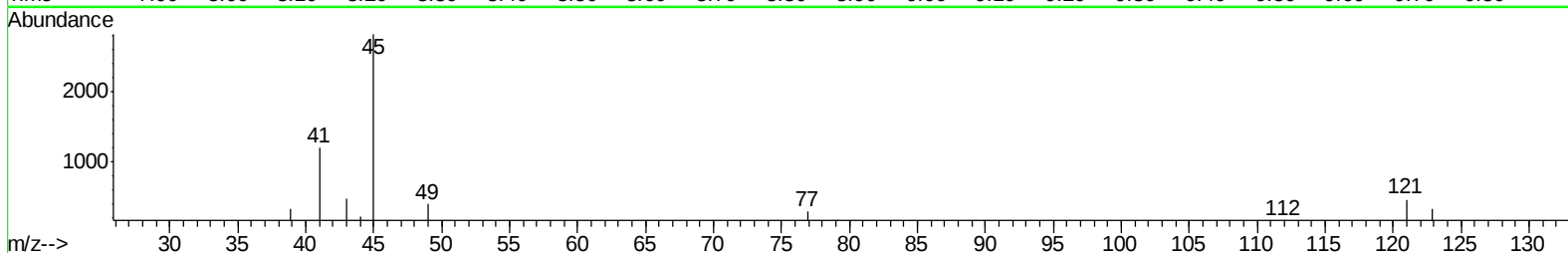
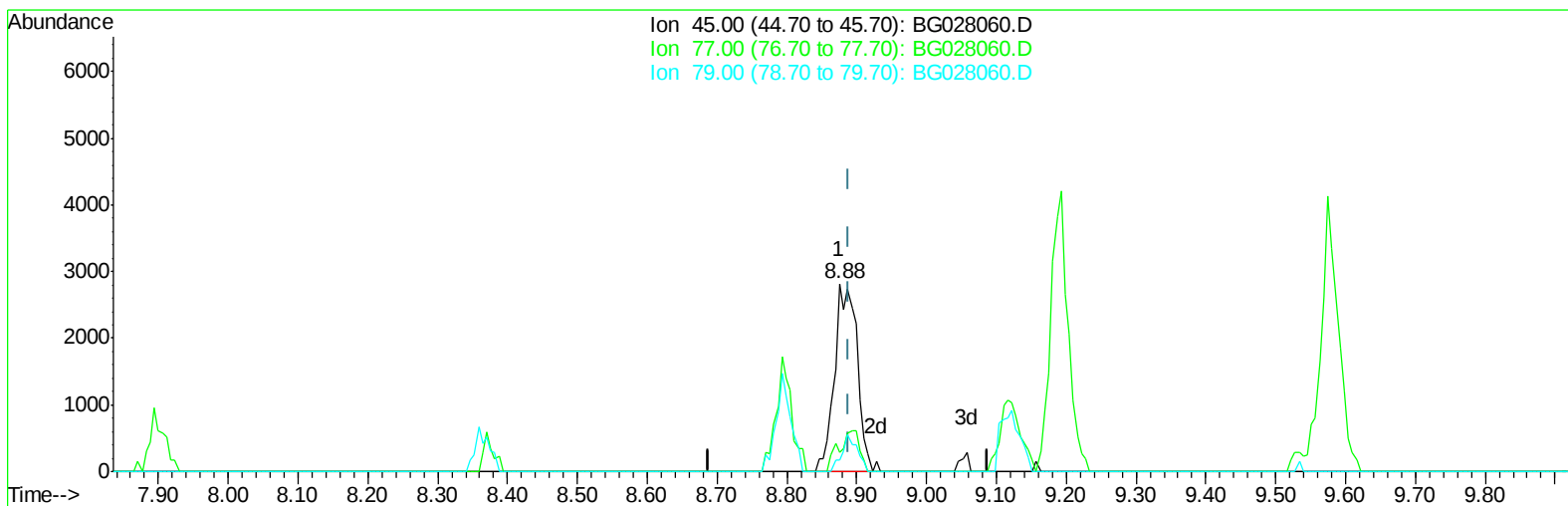
8.875min (-0.013) 1.47ng/ul

response 3032

Ion	Exp%	Act%
45.00	100	100
77.00	15.60	10.27#
79.00	11.80	6.01#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(12) 2,2'-oxybis(1-Chloropropane)

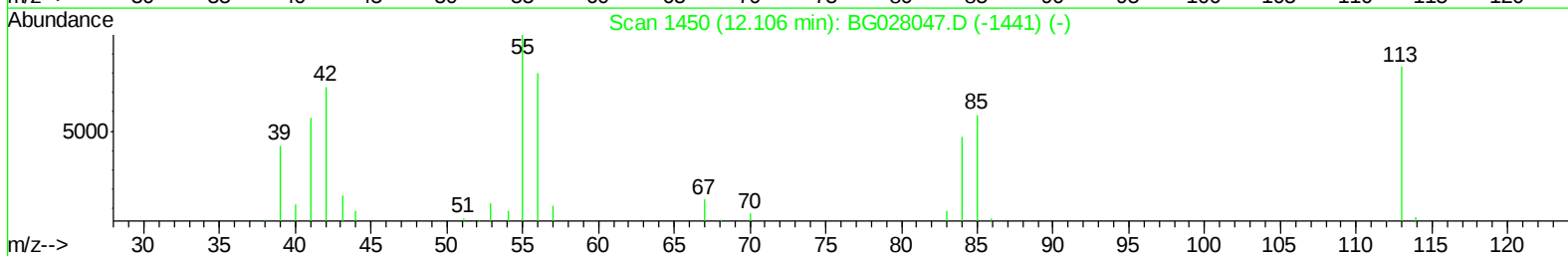
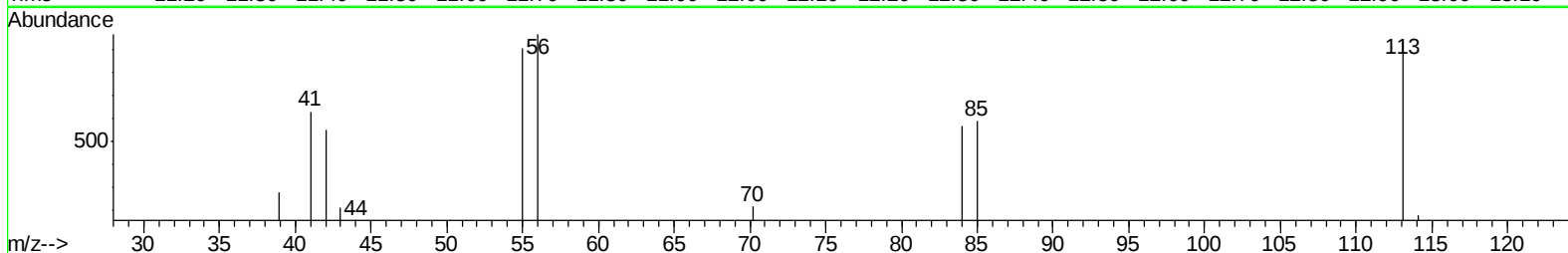
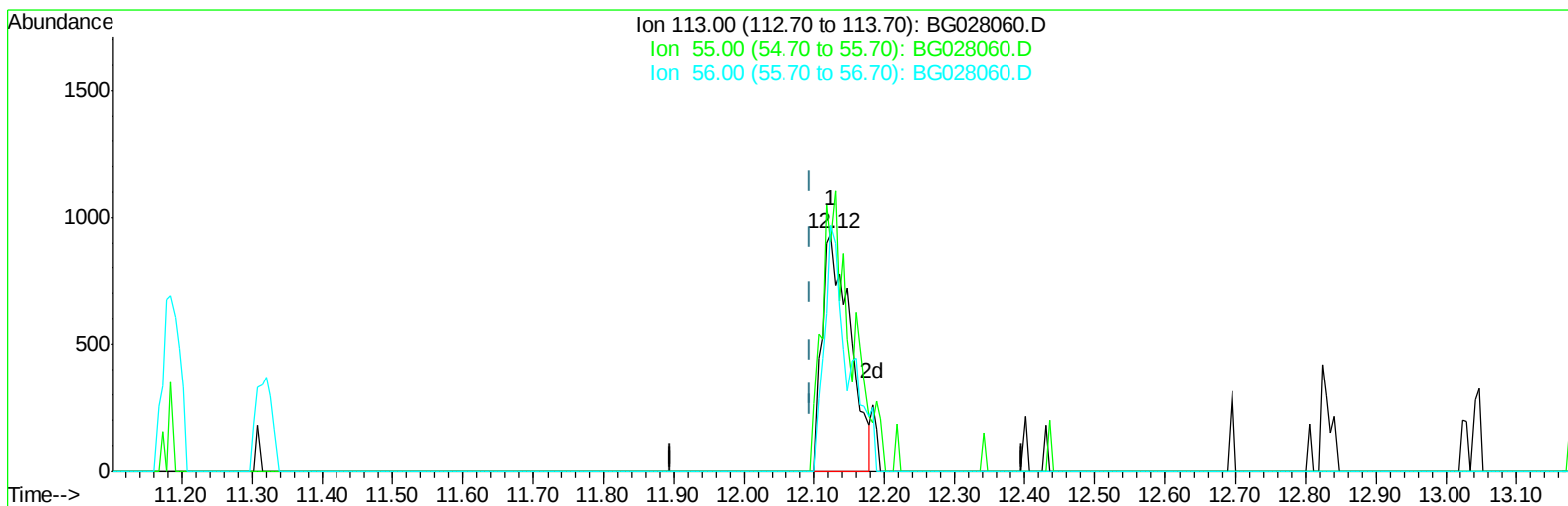
8.875min (-0.013) 3.05ng/ul m

response 6284

Ion	Exp%	Act%
45.00	100	100
77.00	15.60	10.27#
79.00	11.80	6.01#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(32) Caprolactam

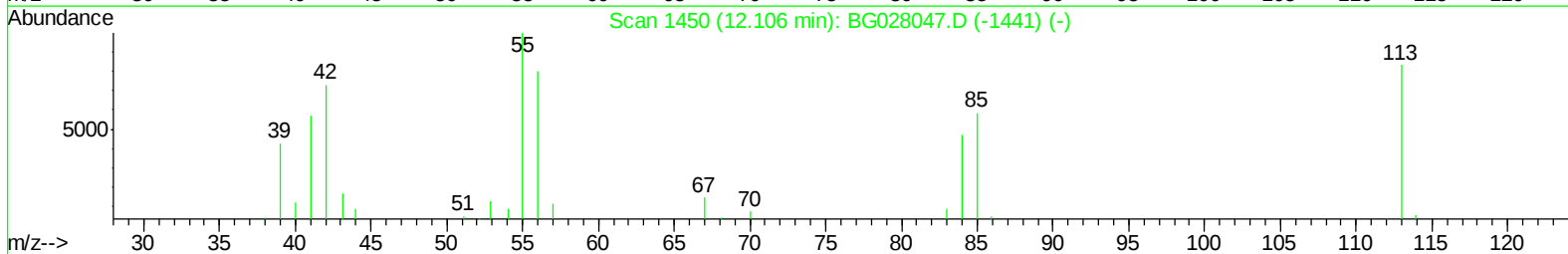
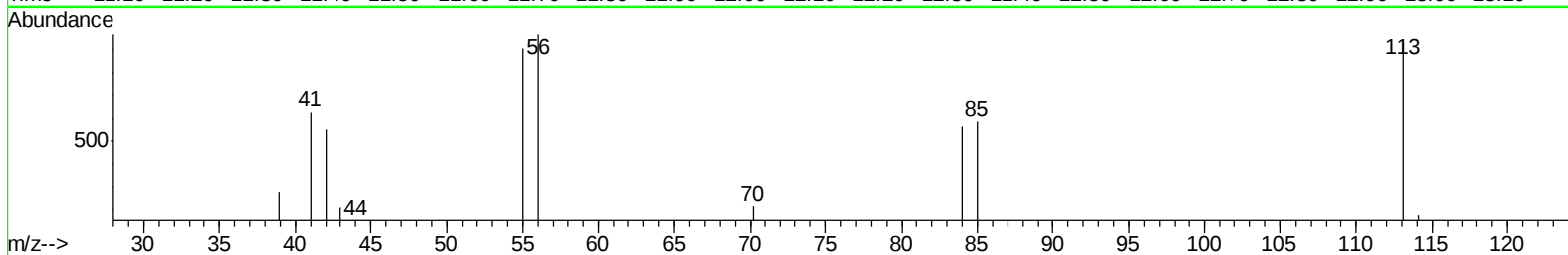
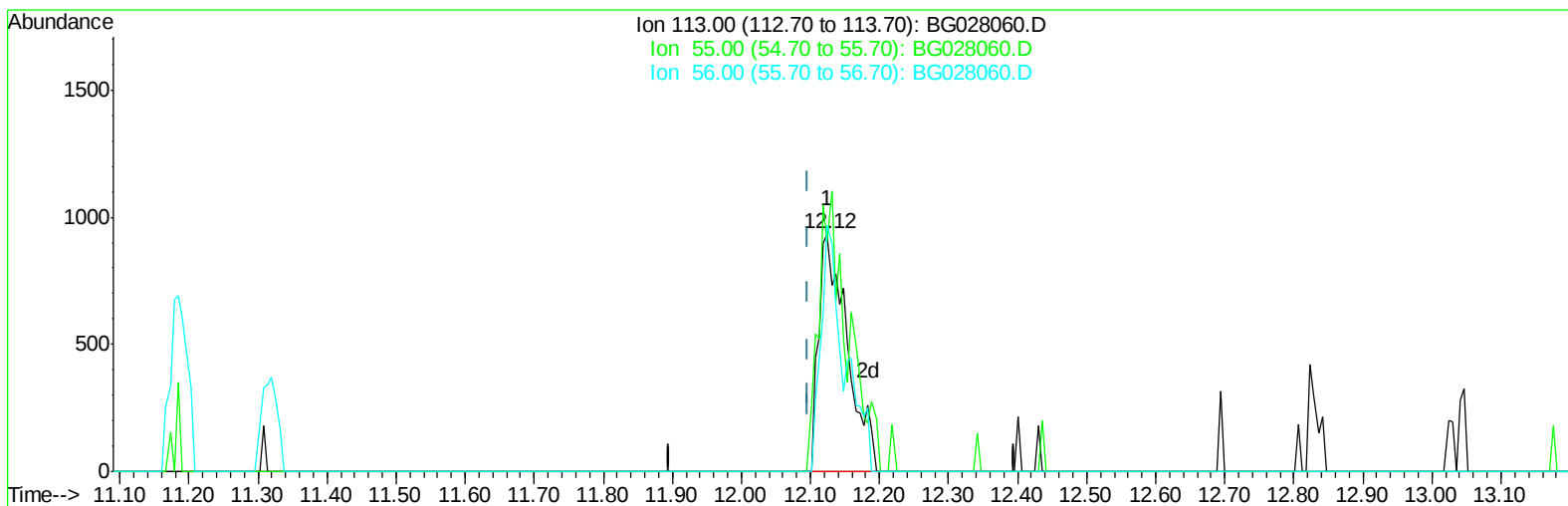
12.124min (+0.028) 3.62ng/ul

response 2542

Ion	Exp%	Act%
113.00	100	100
55.00	164.60	96.79#
56.00	115.90	103.53
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(32) Caprolactam

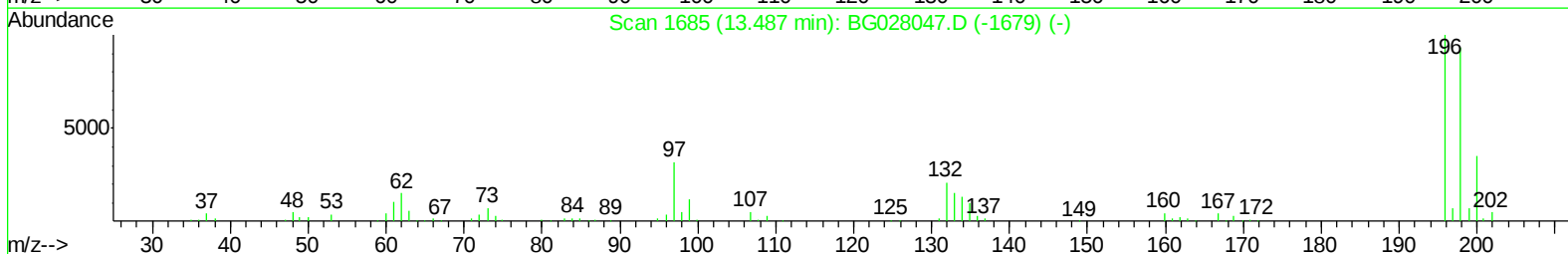
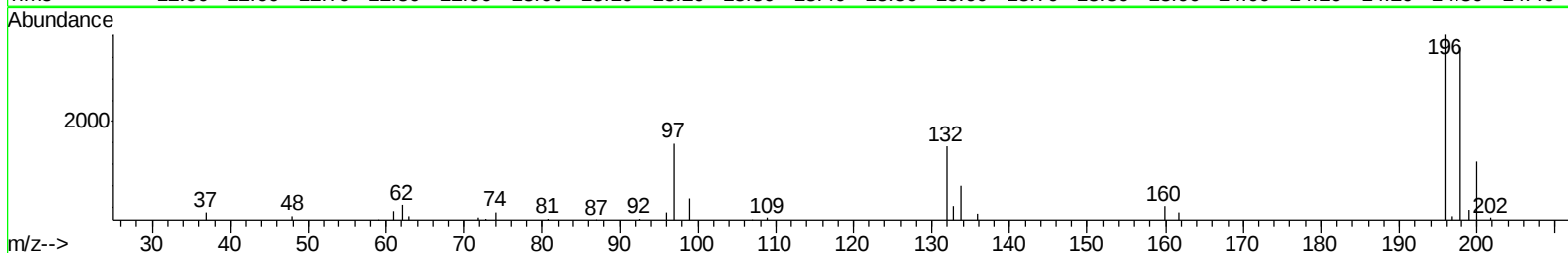
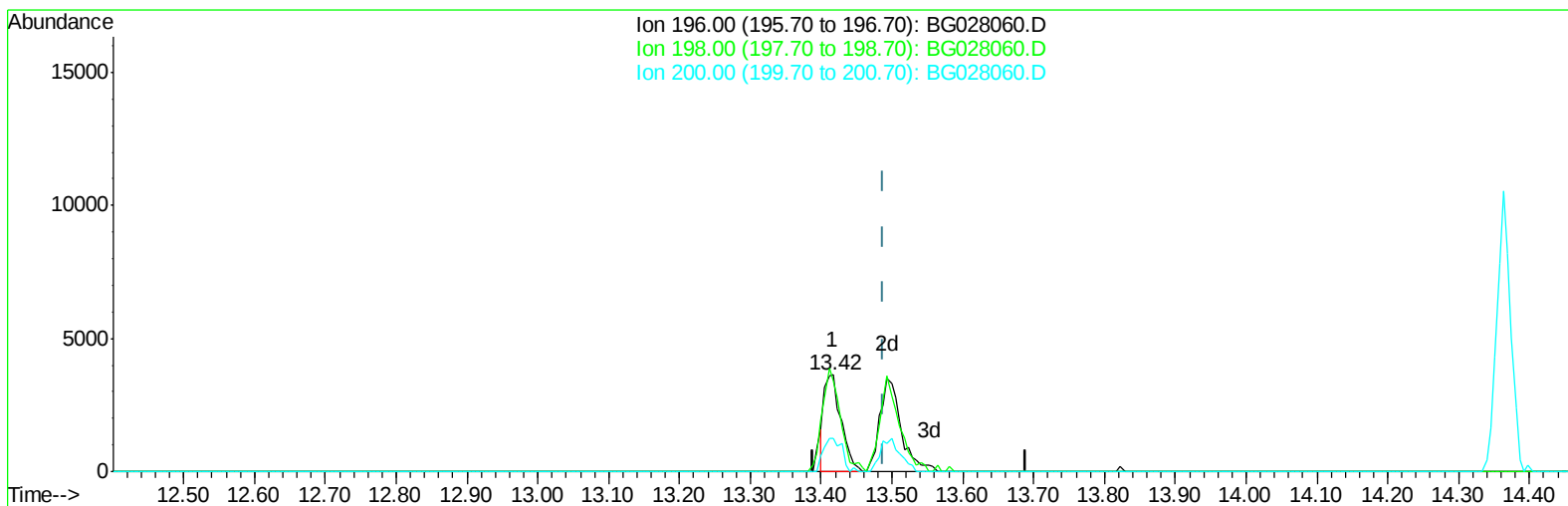
12.124min (+0.028) 3.84ng/ul m

response 2694

Ion	Exp%	Act%
113.00	100	100
55.00	164.60	96.79#
56.00	115.90	103.53
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(40) 2,4,5-Trichlorophenol

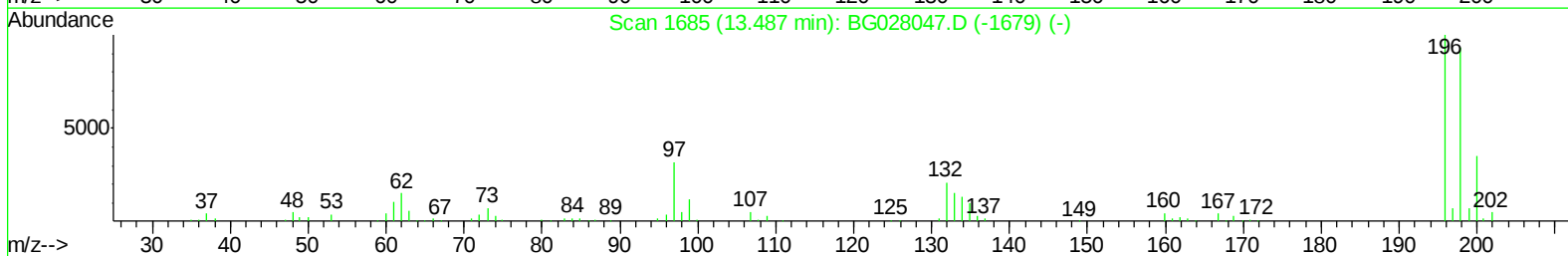
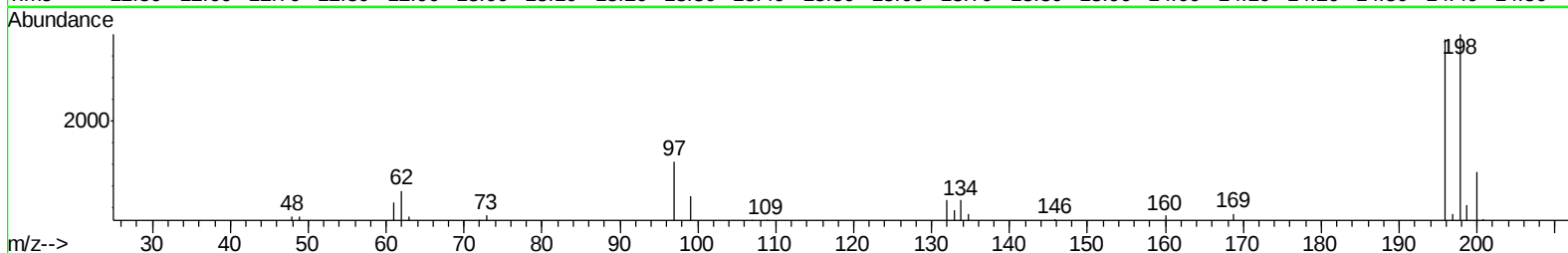
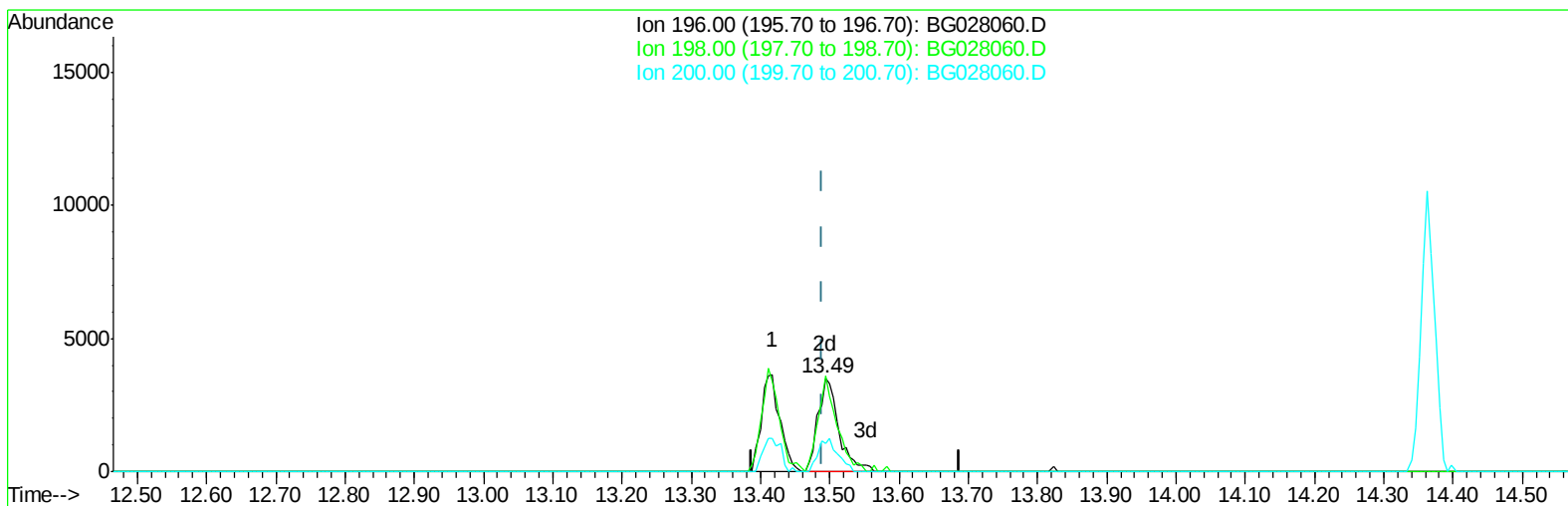
13.417min (-0.072) 2.35ng/ul

response 5966

Ion	Exp%	Act%
196.00	100	100
198.00	93.40	88.12
200.00	31.70	32.64
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(40) 2,4,5-Trichlorophenol

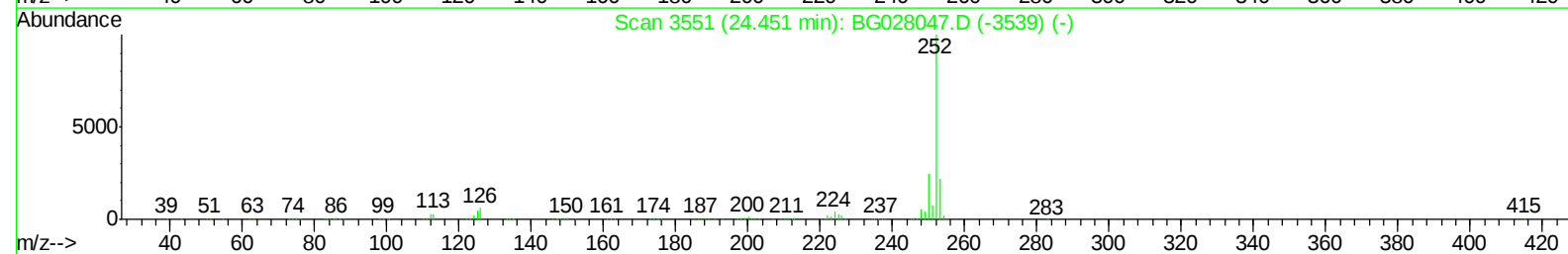
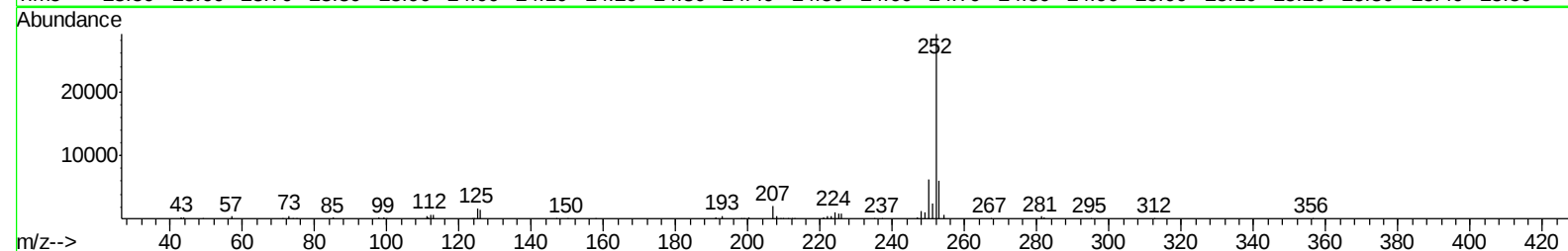
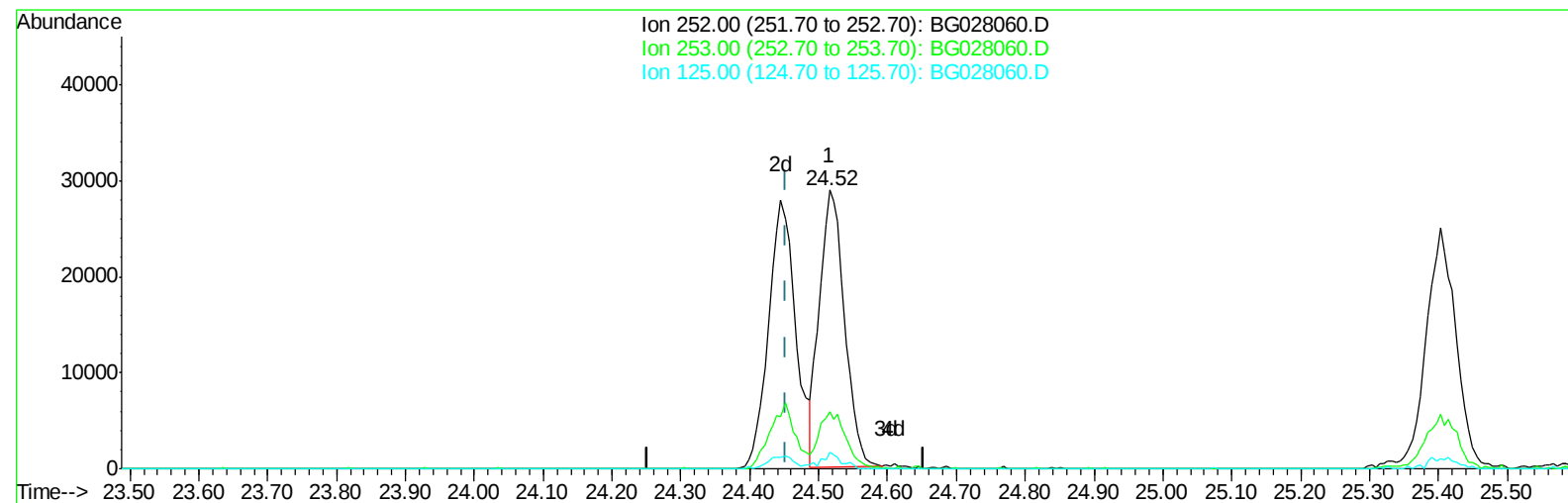
13.493min (+0.004) 2.89ng/ul m

response 7350

Ion	Exp%	Act%
196.00	100	100
198.00	93.40	102.39
200.00	31.70	29.72
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(88) Benzo(b)fluoranthene

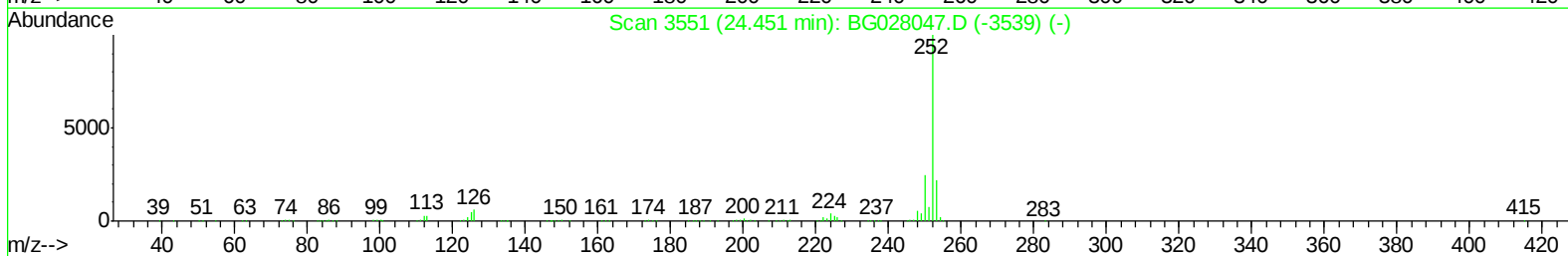
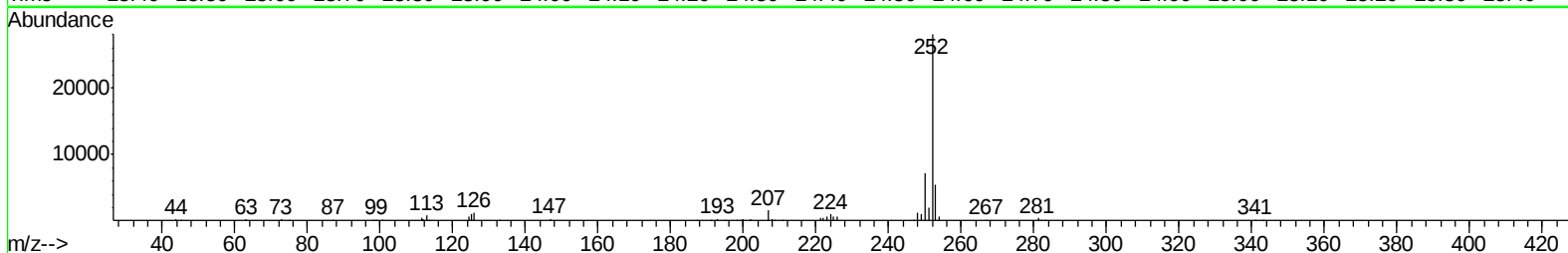
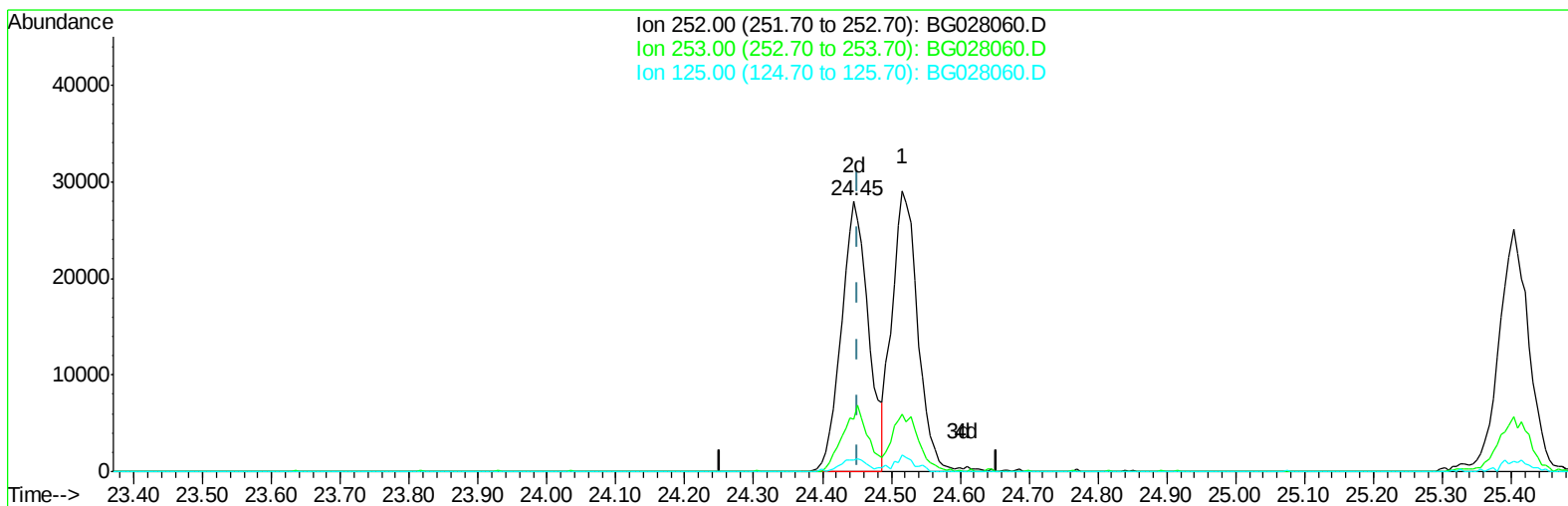
24.516min (+0.063) 3.25ng/ul

response 73092

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	20.66
125.00	6.10	6.15
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(88) Benzo(b)fluoranthene

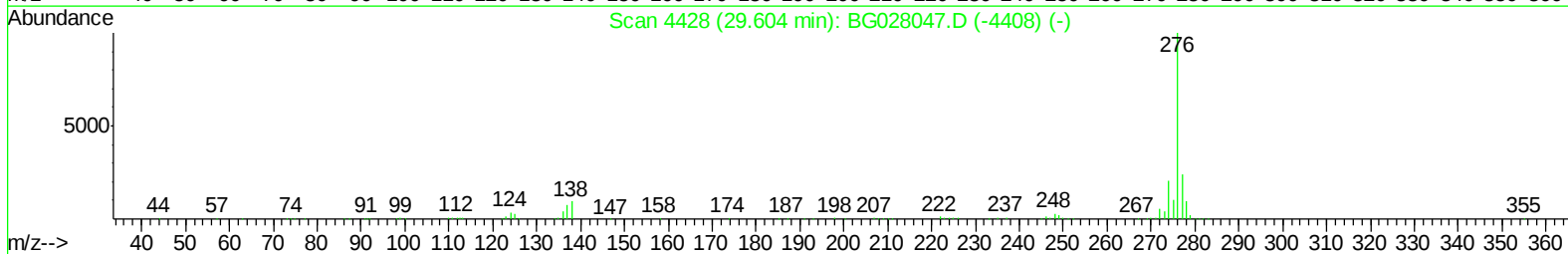
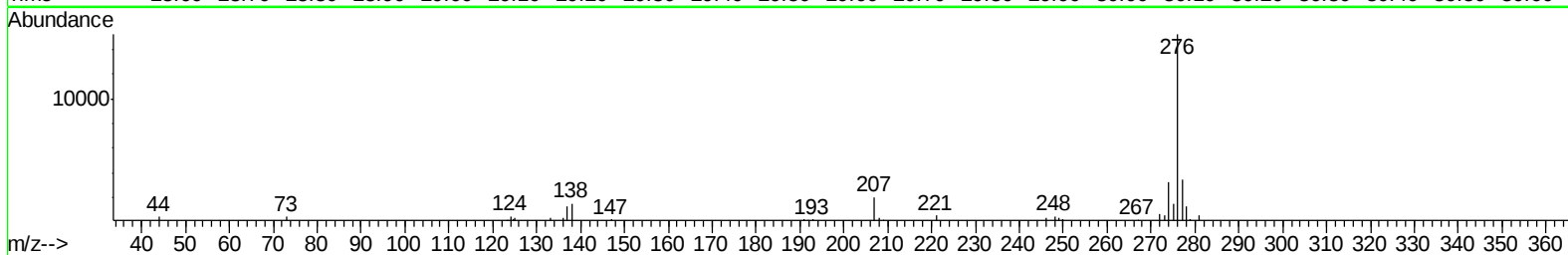
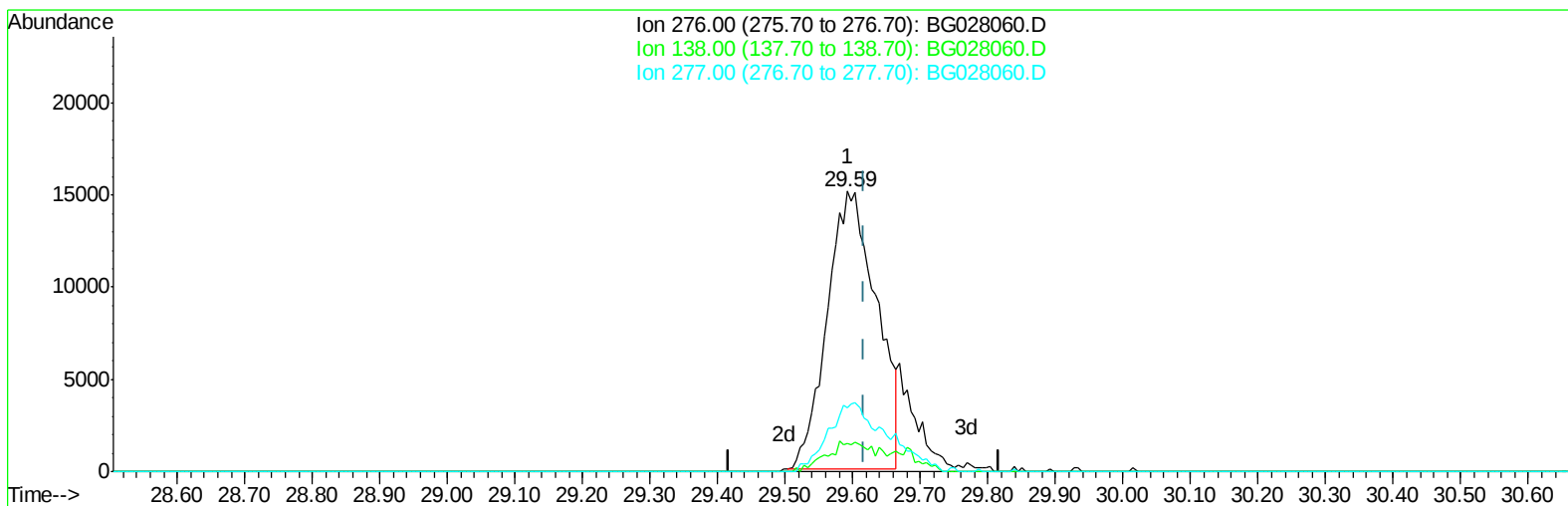
24.445min (-0.007) 3.42ng/ul m

response 76710

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	19.38
125.00	6.10	4.10#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

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

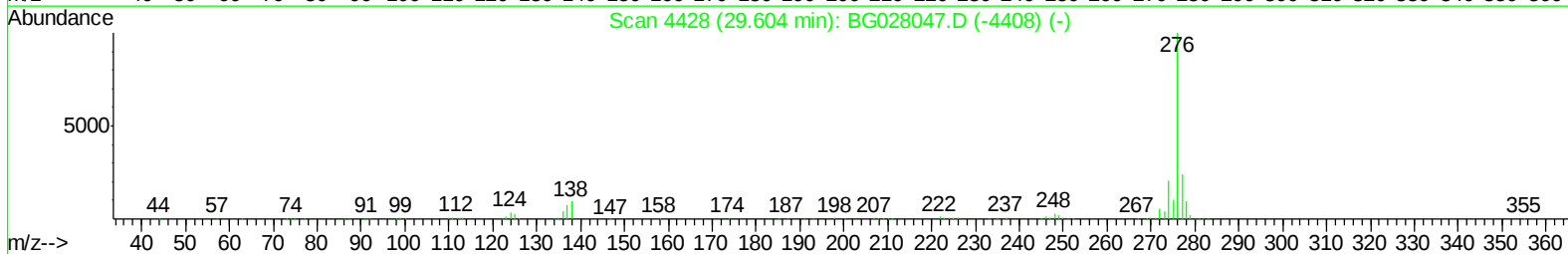
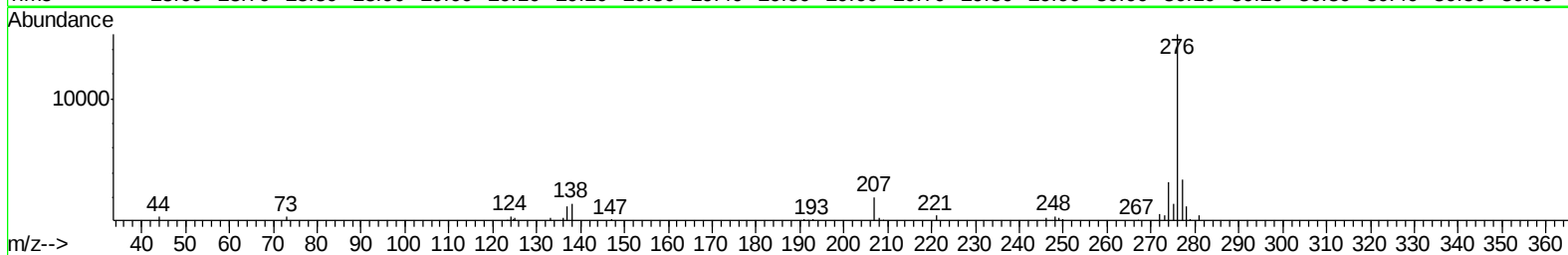
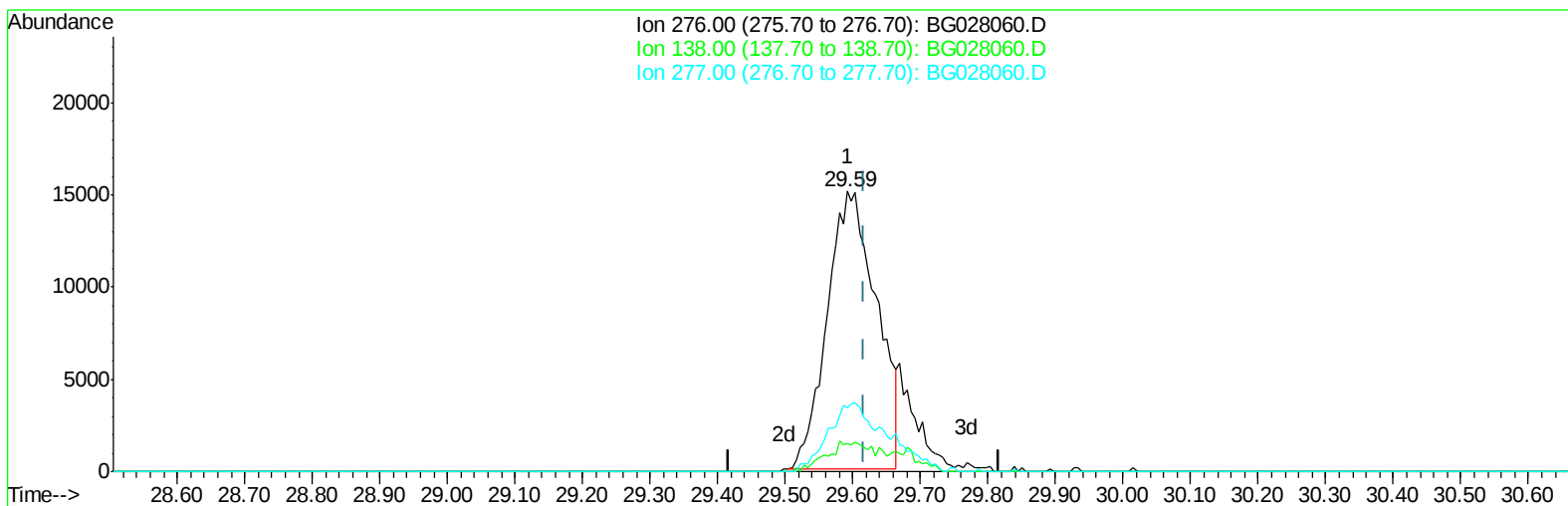
29.592min (-0.025) 2.91ng/ul

response 76264

Ion	Exp%	Act%
276.00	100	100
138.00	13.60	9.94#
277.00	23.60	22.56
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

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

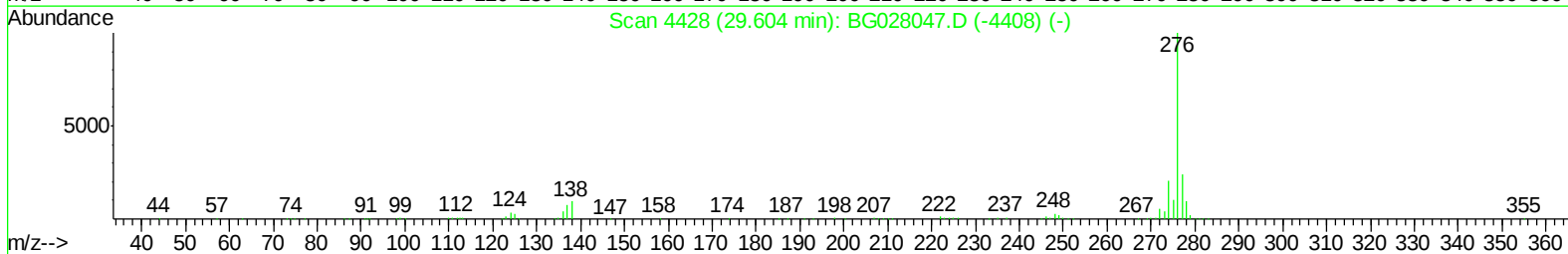
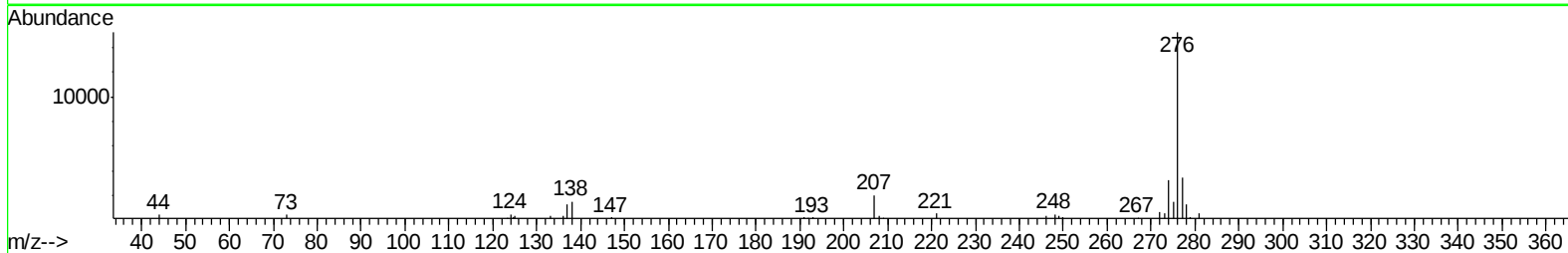
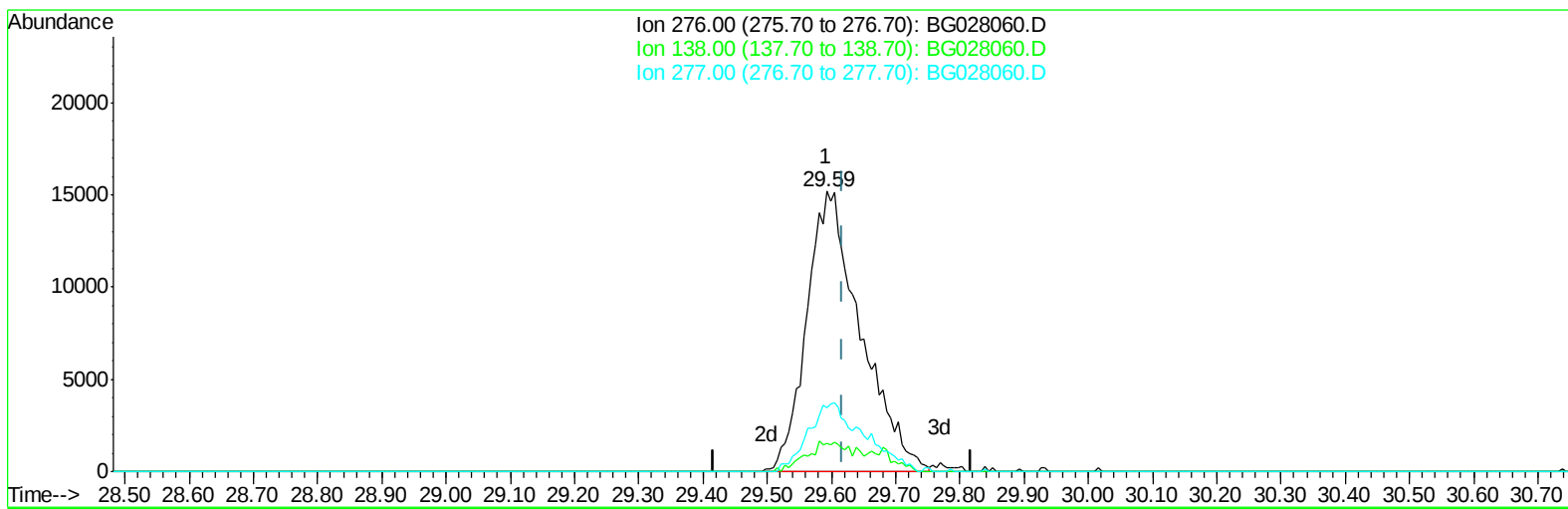
29.592min (-0.025) 2.91ng/ul

response 76264

Ion	Exp%	Act%
276.00	100	100
138.00	13.60	9.94#
277.00	23.60	22.56
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

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

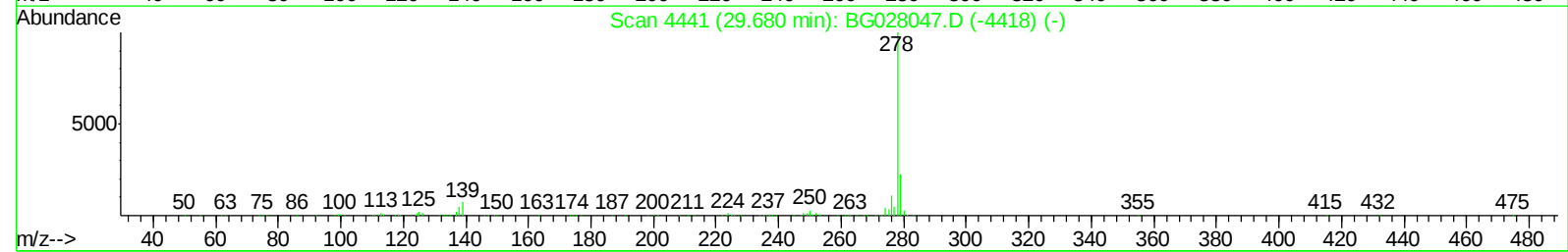
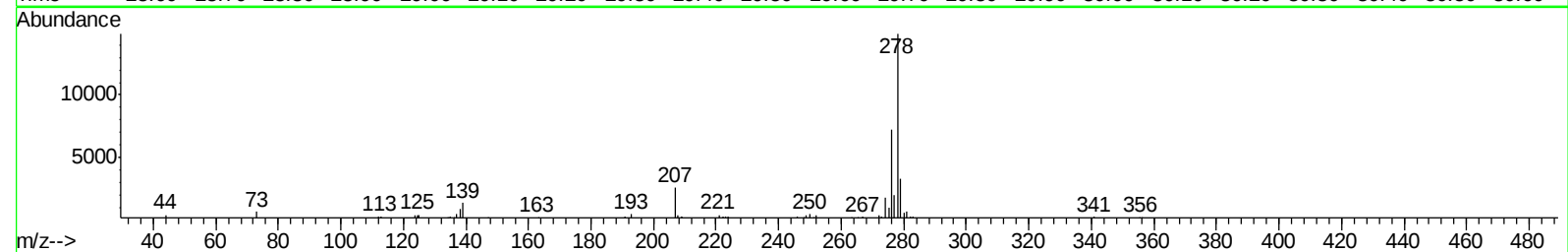
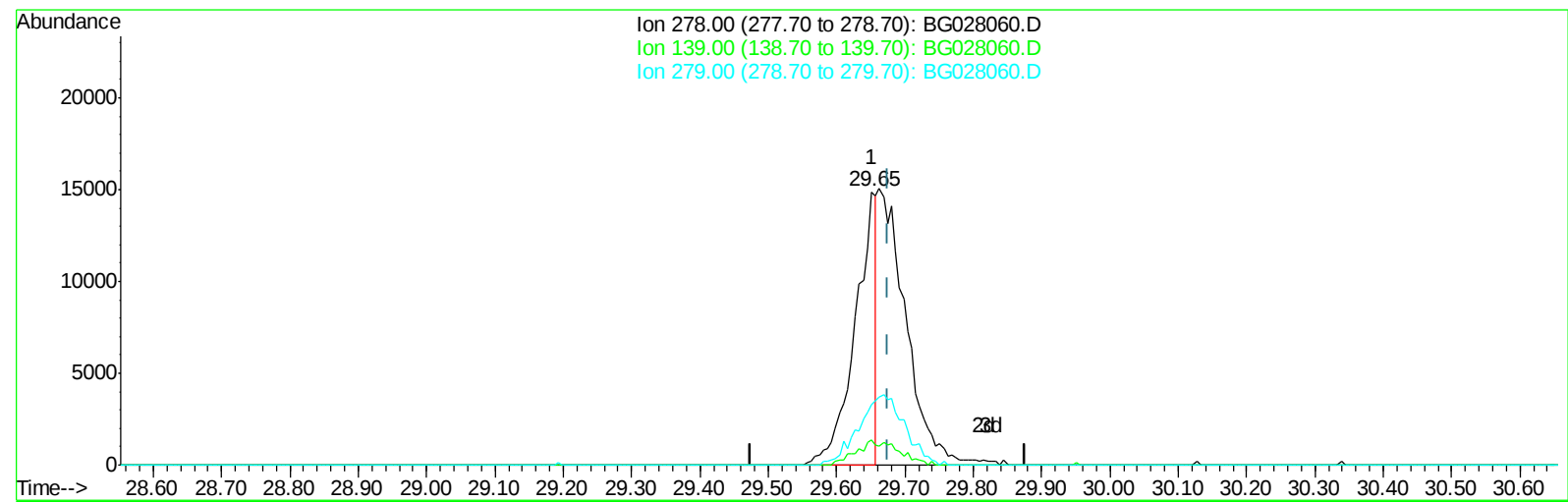
29.592min (-0.025) 3.39ng/ul m

response 89037

Ion	Exp%	Act%
276.00	100	100
138.00	13.60	9.94#
277.00	23.60	22.56
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(93) Dibenzo(a,h)anthracene

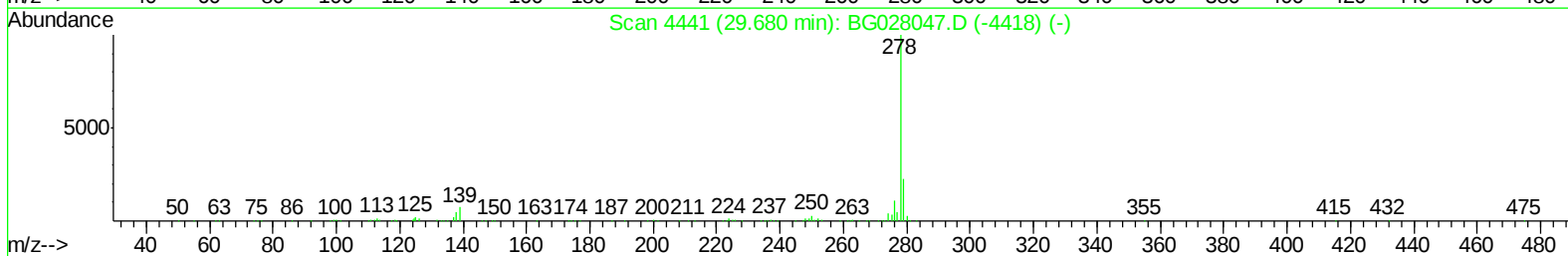
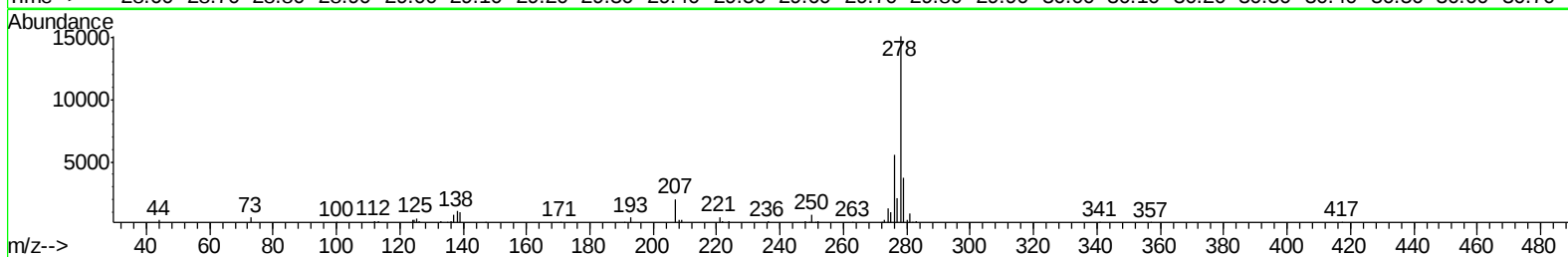
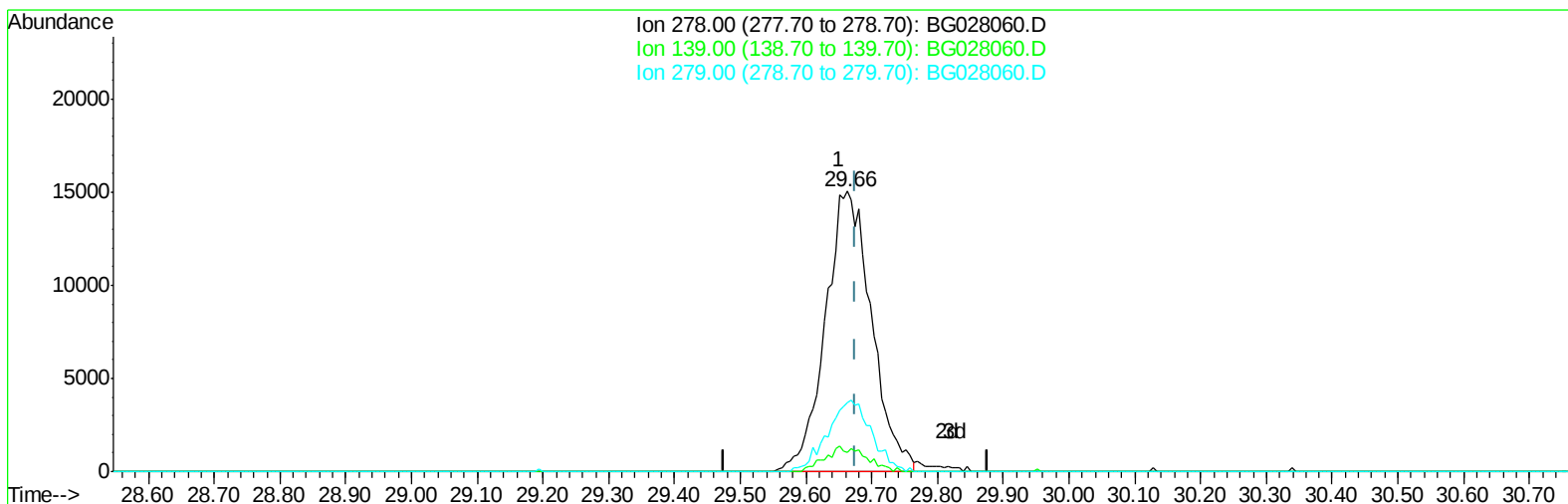
29.651min (-0.025) 1.44ng/ul

response 32408

Ion	Exp%	Act%
278.00	100	100
139.00	11.20	9.02
279.00	23.90	22.02
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



TIC: BG028060.D

(93) Dibenzo(a,h)anthracene

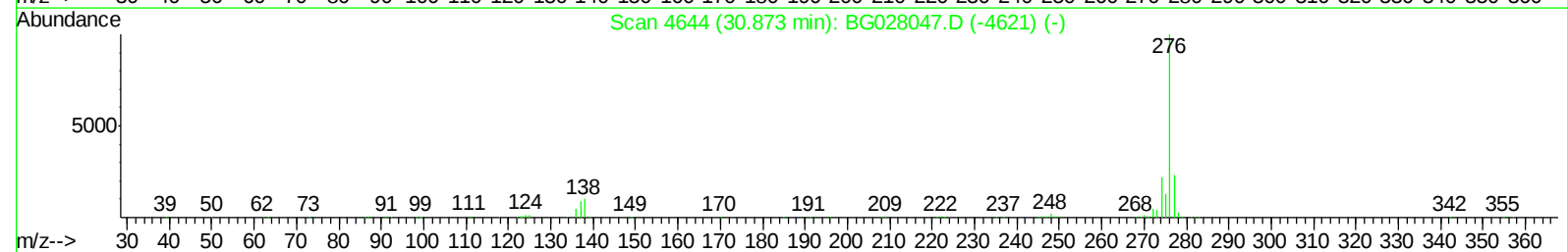
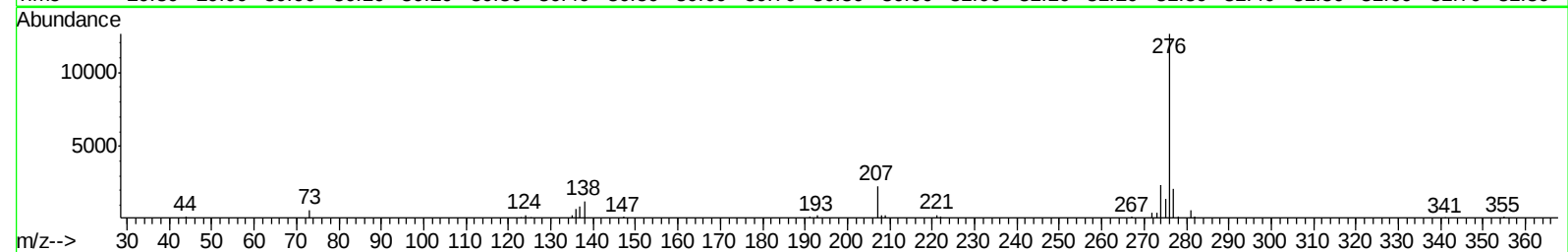
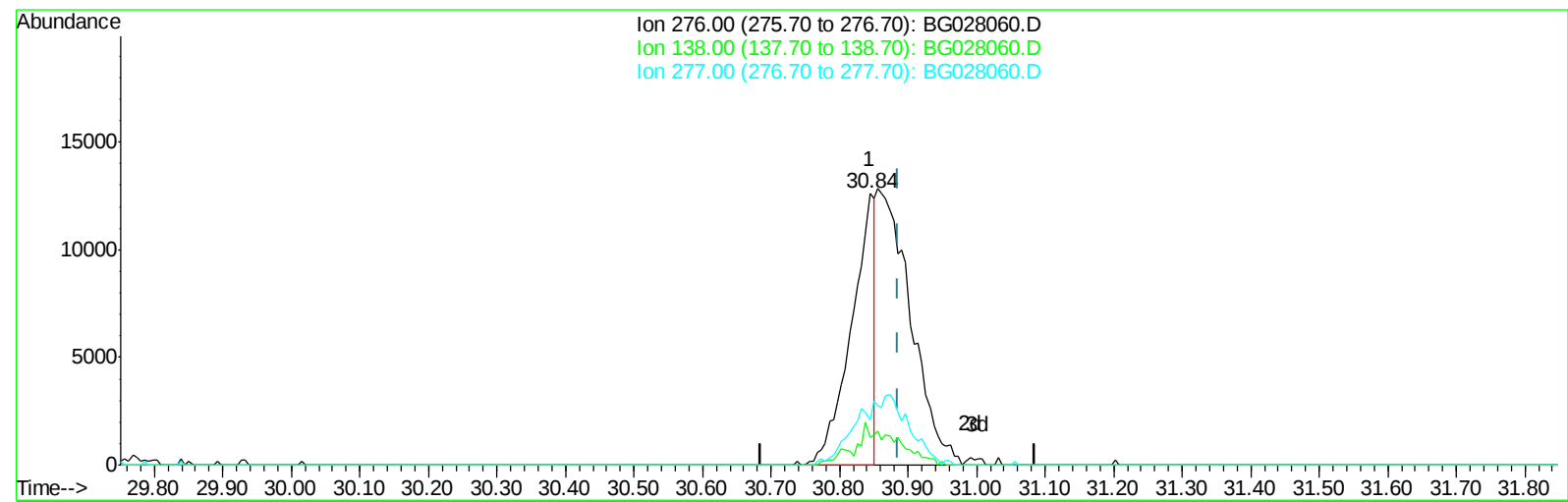
29.663min (-0.013) 3.29ng/ul m

response 73927

Ion	Exp%	Act%
278.00	100	100
139.00	11.20	6.66#
279.00	23.90	24.57
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072817\
Data File : BG028060.D
Acq On : 28 Jul 2017 17:27
Operator : SJ/JU
Sample : MDL-S-ML-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:39:43 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 28 13:35:26 2017
Response via : Initial Calibration



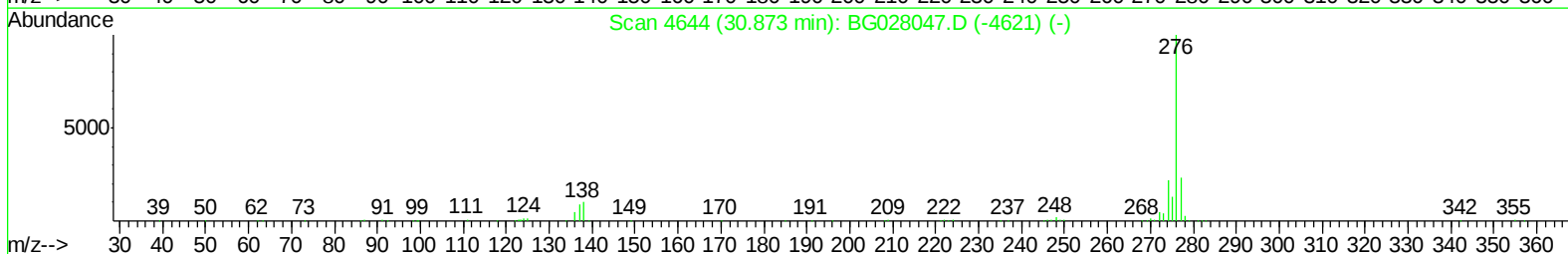
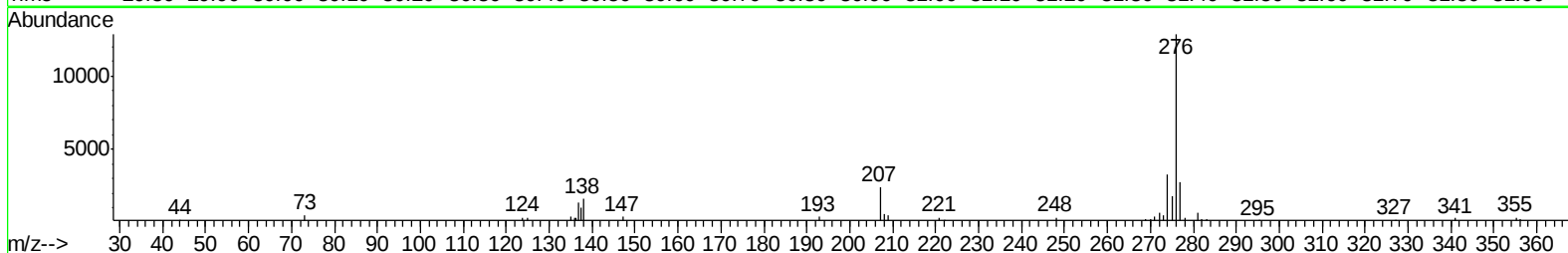
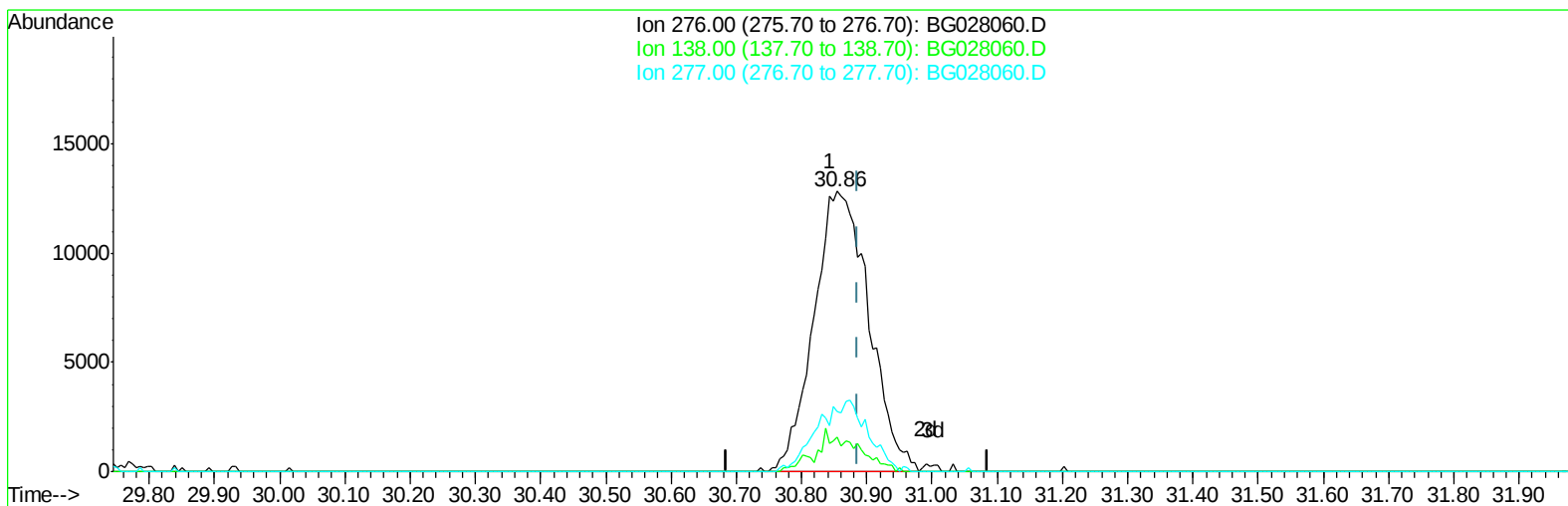
TIC: BG028060.D

(94) Benzo(g,h,i)perylene
30.844min (-0.043) 1.37ng/ul
response 29817

Ion	Exp%	Act%
276.00	100	100
138.00	14.00	10.24#
277.00	23.80	16.67#
0.00	0.00	0.00

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TIC: BG028060.D

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

30.855min (-0.031) 3.41ng/ul m

response 74045

Ion	Exp%	Act%
276.00	100	100
138.00	14.00	12.40
277.00	23.80	21.55
0.00	0.00	0.00

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 Acq On : 28 Jul 2017 17:27
 Operator : SJ/JU
 Sample : MDL-S-ML-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jul 28 18:52:52 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 28 13:35:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	30222	20.00	ng/ul	0.00
18) Naphthalene-d8	11.18	136	130906	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.97	164	103932	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.72	188	302541	20.00	ng/ul	0.00
78) Chrysene-d12	22.06	240	399989	20.00	ng/ul	0.00
86) Perylene-d12	25.57	264	403453	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.85	96	3168	6.79	ng/uL	0.00
5) Phenol-d5	7.51	99	59109	25.89	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.68	67	32090	28.11	ng/ul	0.00
9) 2-Chlorophenol-d4	7.90	132	57003	29.62	ng/ul	0.00
13) 4-Methylphenol-d8	9.06	113	56307	27.63	ng/ul	0.00
19) Nitrobenzene-d5	9.53	128	28140	30.76	ng/ul	0.00
22) 2-Nitrophenol-d4	10.26	143	36699	31.49	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.80	165	65660	28.63	ng/ul	0.00
29) 4-Chloroaniline-d4	11.32	131	78838	34.08	ng/ul	0.00
44) Dimethylphthalate-d6	14.36	166	295453	31.41	ng/ul	0.00
47) Acenaphthylene-d8	14.67	160	305468	30.02	ng/ul	0.00
52) 4-Nitrophenol-d4	15.16	143	36410	25.77	ng/ul	0.00
58) Fluorene-d10	15.96	176	262548	30.30	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	16.07	200	56593	26.27	ng/ul	0.00
71) Anthracene-d10	17.82	188	449706	31.04	ng/ul	0.00
79) Pyrene-d10	20.10	212	565656	31.84	ng/ul	0.00
90) Benzo(a)pyrene-d12	25.33	264	576466	31.06	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.89	88	782	1.564 ng/uL#	88
4) Benzaldehyde	7.51	77	994m	0.880 ng/ul	
6) Phenol	7.53	94	7578	3.423 ng/ul	89
8) Bis(2-Chloroethyl)ether	7.78	93	5370	3.427 ng/ul	97
10) 2-Chlorophenol	7.93	128	5946	3.354 ng/ul#	81
11) 2-Methylphenol	8.79	108	5178	3.000 ng/ul#	79
12) 2,2'-oxybis(1-Chloropropan	8.88	45	6284m	3.055 ng/ul	
14) Acetophenone	9.19	105	9598	3.242 ng/ul#	86
15) N-Nitroso-di-n-propylamine	9.16	70	4177	3.005 ng/ul	93
16) 4-Methylphenol	9.12	108	6399	3.341 ng/ul	80
17) Hexachloroethane	9.46	117	2442	3.443 ng/ul#	70
20) Nitrobenzene	9.57	77	7067	3.597 ng/ul#	83
21) Isophorone	10.09	82	13423	3.575 ng/ul#	92
23) 2-Nitrophenol	10.29	139	4318	3.845 ng/ul#	66
24) 2,4-Dimethylphenol	10.34	107	7345	3.270 ng/ul#	92
25) Bis(2-Chloroethoxy)methane	10.57	93	7657	3.511 ng/ul	99
27) 2,4-Dichlorophenol	10.83	162	8092	3.702 ng/ul#	93
28) Naphthalene	11.24	128	21914	3.709 ng/ul	95
30) 4-Chloroaniline	11.34	127	5592	2.540 ng/ul	92
31) Hexachlorobutadiene	11.52	225	7422	3.925 ng/ul	99
32) Caprolactam	12.12	113	2694m	3.841 ng/ul	
33) 4-Chloro-3-methylphenol	12.44	107	6577	3.088 ng/ul	95
34) 2-Methylnaphthalene	12.82	142	17918	3.514 ng/ul	95

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Quant Time: Jul 28 18:52:52 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072717MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 28 13:35:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	13.04	142	18573	3.756	ng/ul#	90
37) 1,2,4,5-Tetrachlorobenzene	13.18	216	14160	3.536	ng/ul	91
38) Hexachlorocyclopentadiene	13.15	237	5416	2.307	ng/ul	90
39) 2,4,6-Trichlorophenol	13.42	196	6827	2.857	ng/ul	95
40) 2,4,5-Trichlorophenol	13.49	196	7350m	2.895	ng/ul	
41) 1,1'-Biphenyl	13.81	154	26432	3.499	ng/ul	99
42) 2-Chloronaphthalene	13.86	162	21576	3.541	ng/ul	97
43) 2-Nitroaniline	14.06	65	4440	3.092	ng/ul#	74
45) Dimethylphthalate	14.41	163	31309	3.633	ng/ul	98
46) 2,6-Dinitrotoluene	14.54	165	5631	3.255	ng/ul#	95
48) Acenaphthylene	14.70	152	35792	3.712	ng/ul	99
49) 3-Nitroaniline	14.88	138	4496	3.127	ng/ul#	74
50) Acenaphthene	15.04	153	22777	3.449	ng/ul	95
51) 2,4-Dinitrophenol	15.10	184	1798	1.628	ng/ul#	79
53) 4-Nitrophenol	15.17	109	3304	2.875	ng/ul#	81
54) Dibenzofuran	15.37	168	35687	3.554	ng/ul	94
55) 2,4-Dinitrotoluene	15.33	165	8870	3.433	ng/ul	89
56) 2,3,4,6-Tetrachlorophenol	15.60	232	7412	2.741	ng/ul	95
57) Diethylphthalate	15.76	149	30577	3.380	ng/ul	96
59) Fluorene	16.02	166	30028	3.447	ng/ul	96
60) 4-Chlorophenyl-phenylether	16.00	204	19306	3.710	ng/ul	95
61) 4-Nitroaniline	16.04	138	5298	3.103	ng/ul#	79
64) 4,6-Dinitro-2-methylphenol	16.08	198	6735	3.263	ng/ul#	87
65) N-Nitrosodiphenylamine	16.21	169	25667	3.361	ng/ul	98
66) 4-Bromophenyl-phenylether	16.90	248	12859	3.305	ng/ul	93
67) Hexachlorobenzene	17.02	284	15686	3.546	ng/ul	94
68) Atrazine	17.15	200	10982	3.014	ng/ul	93
69) Pentachlorophenol	17.38	266	4428	1.917	ng/ul	94
70) Phenanthrene	17.76	178	55249	3.673	ng/ul	97
72) Anthracene	17.86	178	56825	3.628	ng/ul	95
73) 1,2,3,4-Tetrachlorobenzene	13.78	216	14324	3.395	ng/uL	94
74) Pentachlorobenzene	15.29	250	21559	3.438	ng/uL	94
75) Carbazole	18.13	167	44640	3.481	ng/ul	98
76) Di-n-butylphthalate	18.65	149	46670	3.277	ng/ul	98
77) Fluoranthene	19.76	202	70073	3.645	ng/ul#	95
80) Pyrene	20.13	202	77893	3.752	ng/ul#	93
81) Butylbenzylphthalate	20.99	149	19094	3.095	ng/ul#	84
82) 3,3'-Dichlorobenzidine	21.94	252	22451	3.068	ng/ul	98
83) Benzo(a)anthracene	22.04	228	76023	3.578	ng/ul	98
84) Bis(2-ethylhexyl)phthalate	21.90	149	26133	2.895	ng/ul#	88
85) Chrysene	22.11	228	69027	3.552	ng/ul	99
87) Di-n-octyl phthalate	23.21	149	44658	2.808	ng/ul	100
88) Benzo(b)fluoranthene	24.45	252	76710m	3.415	ng/ul	
89) Benzo(k)fluoranthene	24.52	252	74211	3.389	ng/ul	98
91) Benzo(a)pyrene	25.40	252	74105	3.429	ng/ul#	96
92) Indeno(1,2,3-cd)pyrene	29.59	276	89037m	3.395	ng/ul	
93) Dibenzo(a,h)anthracene	29.66	278	73927m	3.286	ng/ul	
94) Benzo(g,h,i)perylene	30.86	276	74045m	3.405	ng/ul	

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

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

