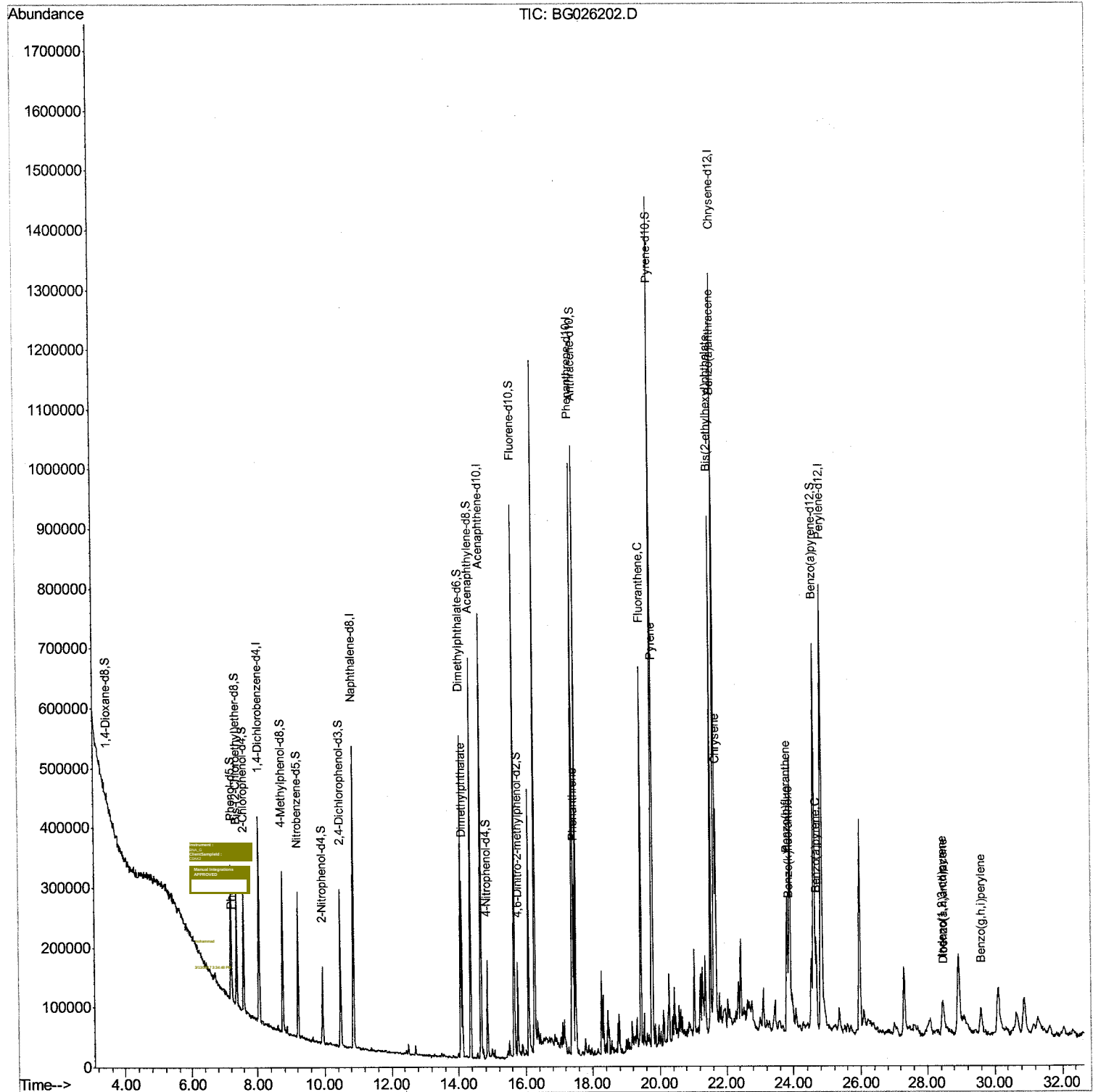


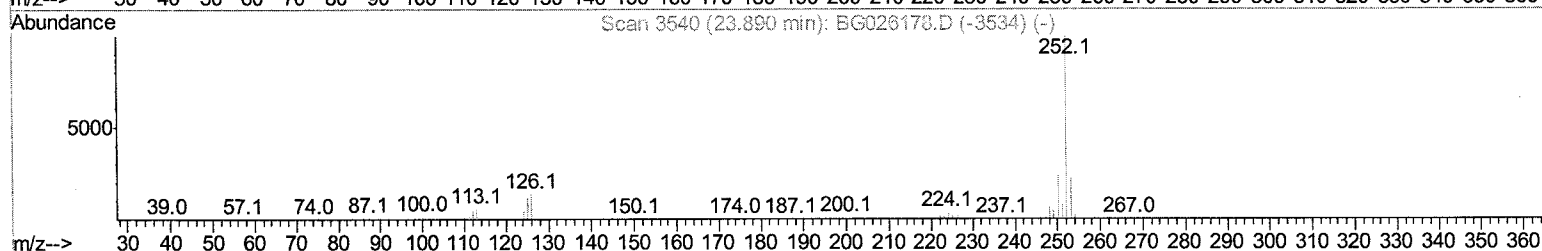
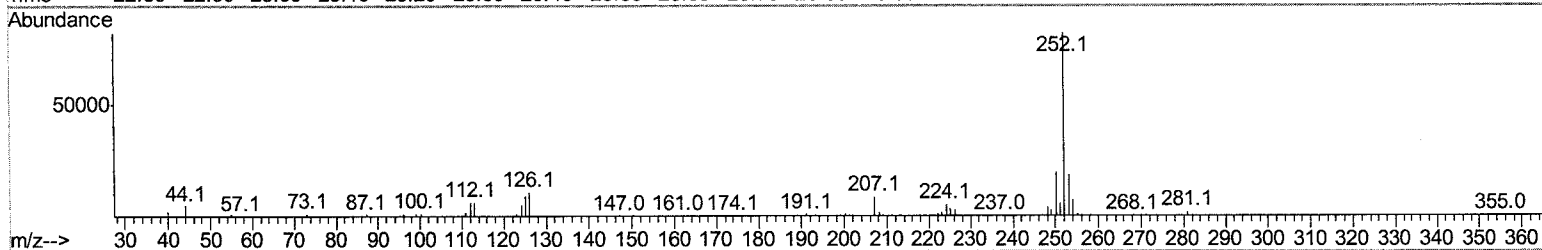
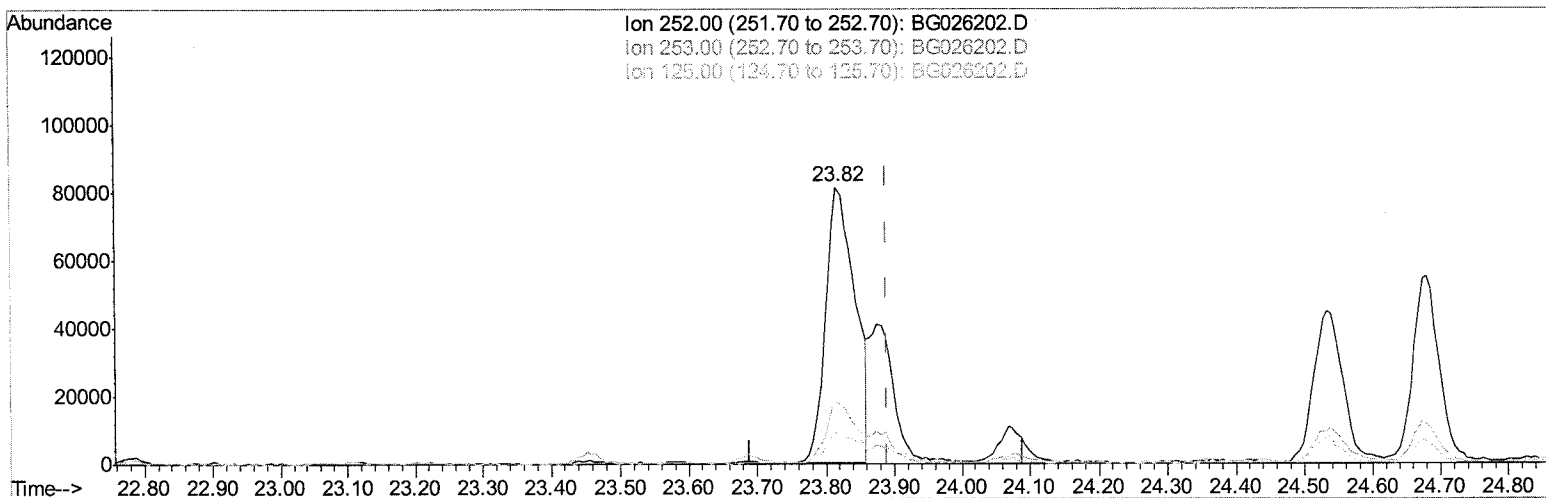
Data Path : Z:\HPCHEM1\BNA G\Data\BG031017\
Data File : BG026202.D
Acq On : 11 Mar 2017 10:38
Operator : SJ/MA
Sample : I2072-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Quant Time: Mar 13 12:11:13 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG030717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Mar 11 00:32:15 2017
Response via : Initial Calibration



Data Path : Z:\HPCHEM1\BNA G\Data\BG031017\
Data File : BG026202.D
Acq On : 11 Mar 2017 10:38
Operator : SJ/MA
Sample : I2072-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Quant Time: Mar 13 12:05:58 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG030717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Mar 11 00:32:15 2017
Response via : Initial Calibration



TIC: BG026202.D

(88) Benzo(k)fluoranthene

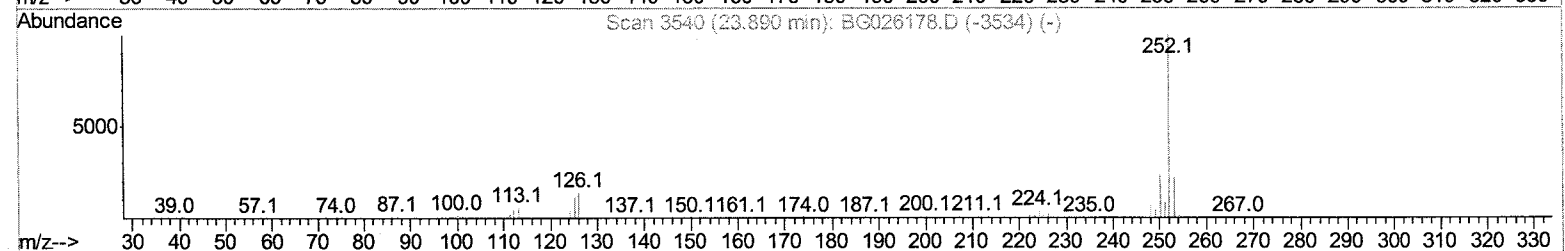
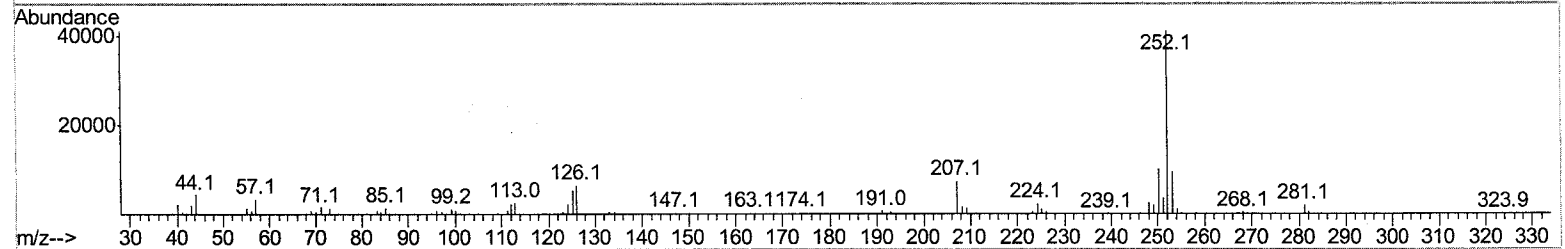
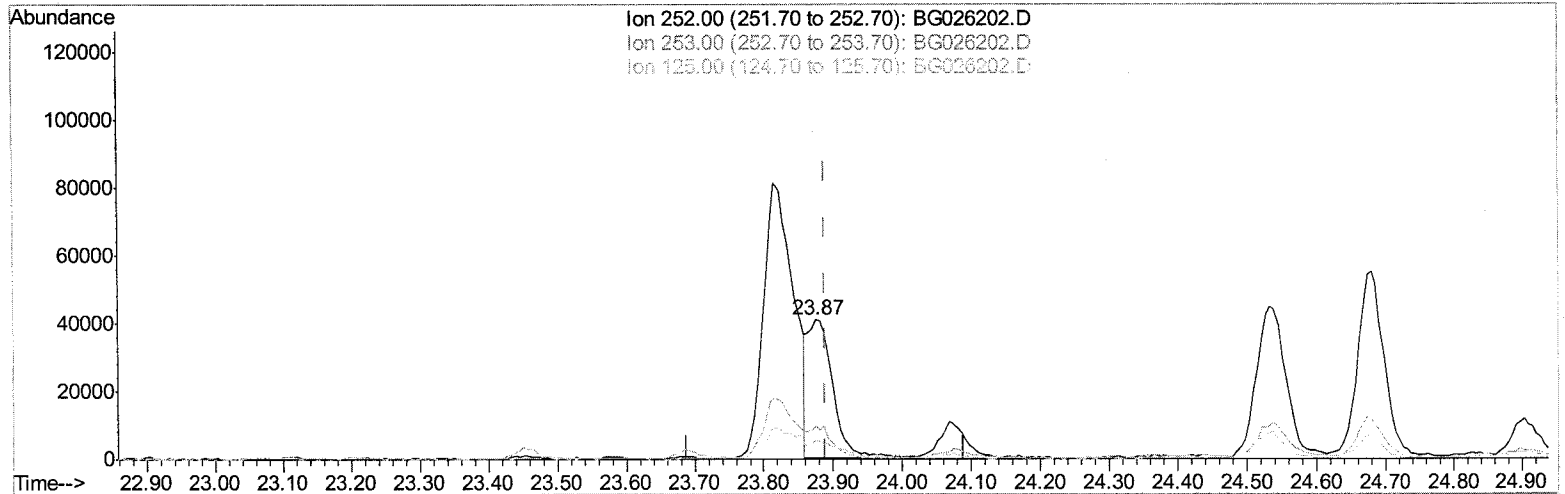
23.816min (-0.073) 6.61ng/ul

response 240432

Ion	Exp%	Found%
252.00	100	100
253.00	21.60	22.46
125.00	10.20	11.44
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA G\Data\BG031017\
Data File : BG026202.D
Acq On : 11 Mar 2017 10:38
Operator : SJ/MA
Sample : I2072-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Quant Time: Mar 13 12:05:58 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG030717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Mar 11 00:32:15 2017
Response via : Initial Calibration



TIC: BG026202.D

(88) Benzo(k)fluoranthene

23.875min (-0.015) 2.75ng/ul m

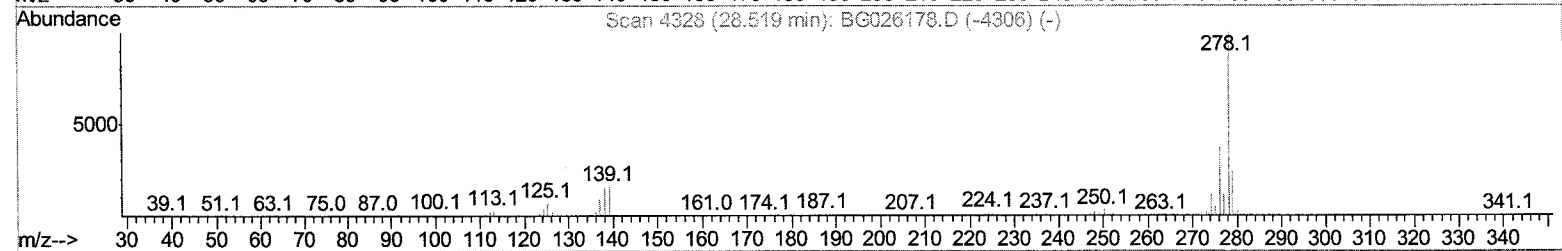
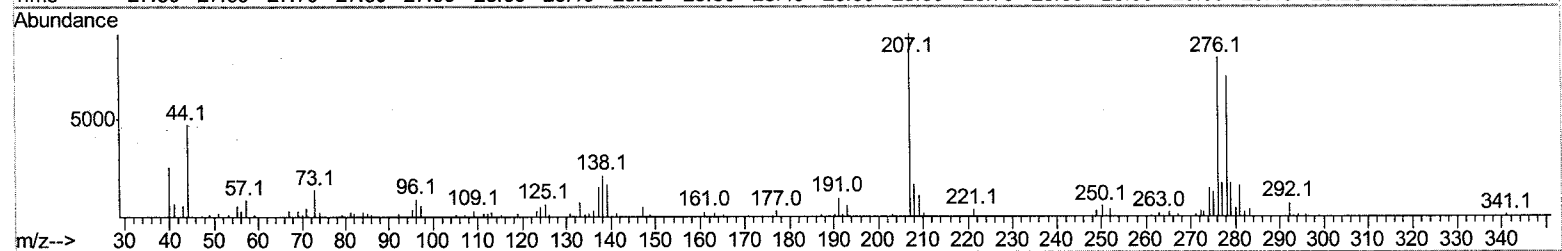
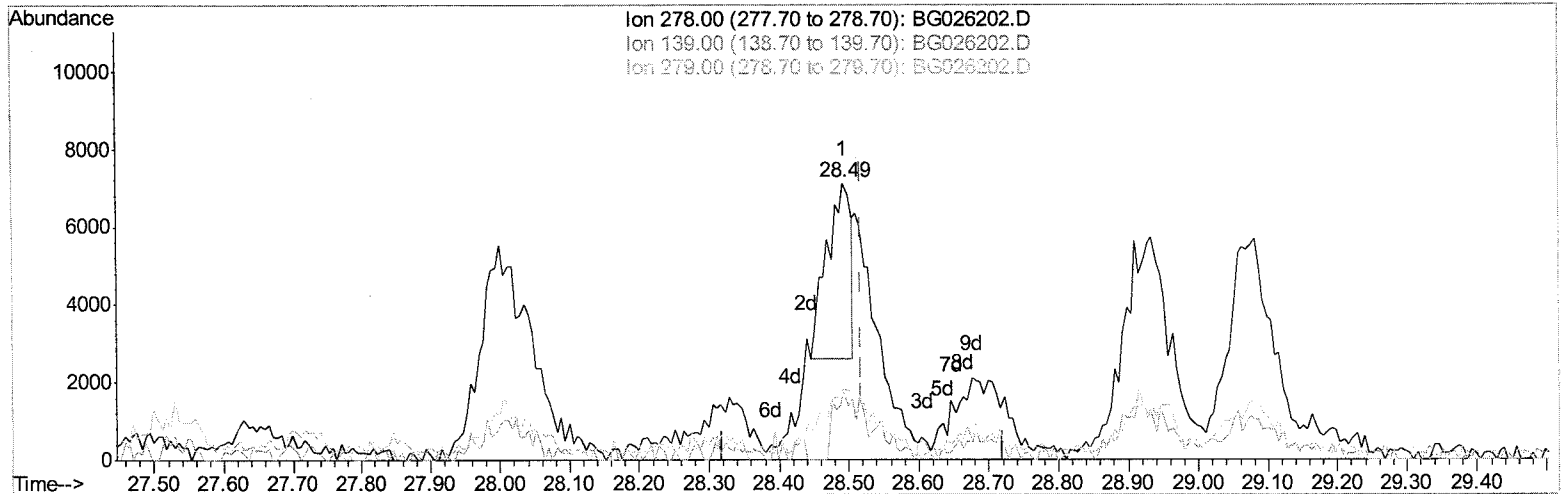
SJ 3/13/17

response 99920

Ion	Exp%	Manual Integration
252.00	100	100
253.00	21.60	23.84
125.00	10.20	13.53#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA G\Data\BG031017\
Data File : BG026202.D
Acq On : 11 Mar 2017 10:38
Operator : SJ/MA
Sample : I2072-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Quant Time: Mar 13 12:05:58 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG030717MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Mar 11 00:32:15 2017
Response via : Initial Calibration



TIC: BG026202.D

(92) Dibenzo(a,h)anthracene

28.493min (-0.026) 0.30ng/ul

response 10804

Ion	Exp%	Manual Integration
278.00	100	100
139.00	17.40	24.61#
279.00	23.70	25.76
0.00	0.00	0.00

```
Data Path : Z:\HPCHEM1\BNA G\Data\BG031017\  
Data File : BG026202.D  
Acq On    : 11 Mar 2017   10:38  
Operator   : SJ/MA  
Sample     : I2072-07  
Misc      :  
ALS Vial   : 37      Sample Multiplier: 1
```

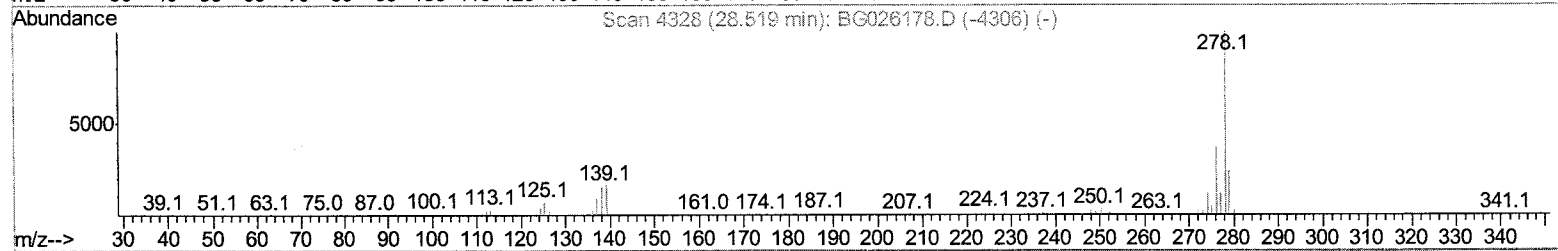
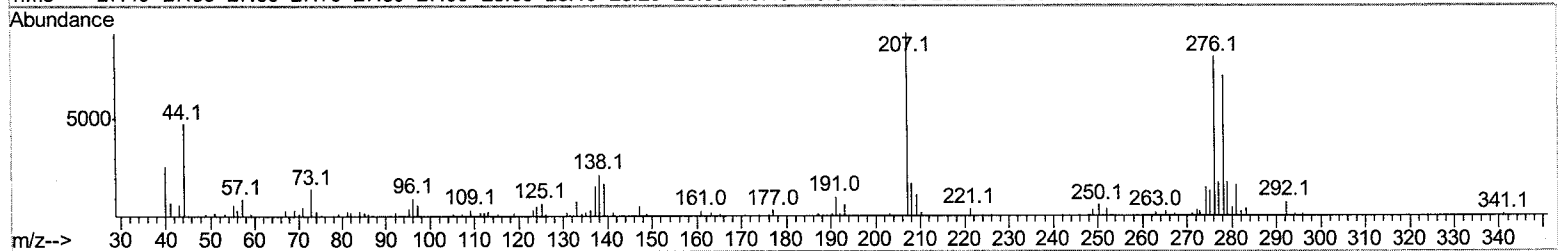
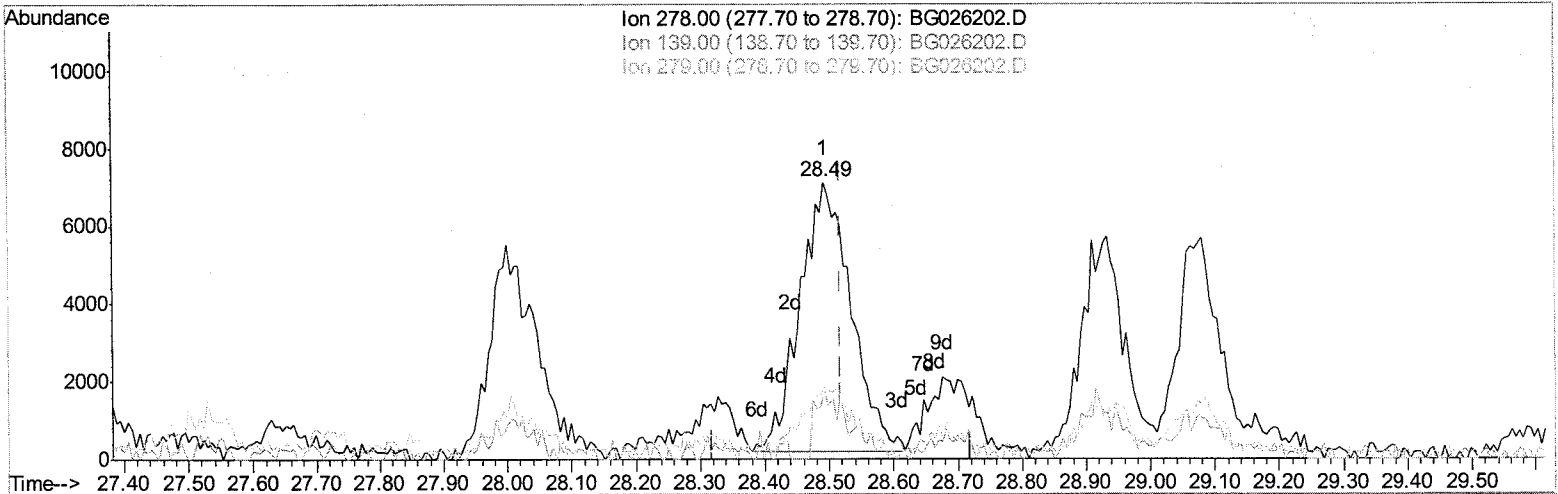
Quant Time: Mar 13 12:05:58 2017

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG030717MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat Mar 11 00:32:15 2017

Response via : Initial Calibration



TIC: BG026202.D

(92) Dibenzo(a,h)anthracene

28.493min (-0.026) 1.03ng/ul m

SD 3/13/17

response 37578

Ion	Exp%	Manual Integrations APPROVED
1	100	
2	100	
3	100	
4	100	
5	100	
6	100	
7	100	
8	100	
9	100	
10	100	
11	100	
12	100	
13	100	
14	100	
15	100	
16	100	
17	100	
18	100	
19	100	
20	100	
21	100	
22	100	
23	100	
24	100	
25	100	
26	100	
27	100	
28	100	
29	100	
30	100	
31	100	
32	100	
33	100	
34	100	
35	100	
36	100	
37	100	
38	100	
39	100	
40	100	
41	100	
42	100	
43	100	
44	100	
45	100	
46	100	
47	100	
48	100	
49	100	
50	100	
51	100	
52	100	
53	100	
54	100	
55	100	
56	100	
57	100	
58	100	
59	100	
60	100	
61	100	
62	100	
63	100	
64	100	
65	100	
66	100	
67	100	
68	100	
69	100	
70	100	
71	100	
72	100	
73	100	
74	100	
75	100	
76	100	
77	100	
78	100	
79	100	
80	100	
81	100	
82	100	
83	100	
84	100	
85	100	
86	100	
87	100	
88	100	
89	100	
90	100	
91	100	
92	100	
93	100	
94	100	
95	100	
96	100	
97	100	
98	100	
99	100	
100	100	

278.00	100	100
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139.00 17.40 ~~mohammad~~ 24.61#

279.00 23.70 25.76

0.00	0.00	0.00
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Data Path : Z:\HPCHEM1\BNA G\Data\BG031017\
 Data File : BG026202.D
 Acq On : 11 Mar 2017 10:38
 Operator : SJ/MA
 Sample : I2072-07
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Quant Time: Mar 13 12:11:13 2017

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG030717MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat Mar 11 00:32:15 2017

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	102306	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	457444	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.67	164	259428	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.39	188	619524	20.00	ng/ul	0.00
77) Chrysene-d12	21.64	240	702483	20.00	ng/ul	0.00
85) Perylene-d12	24.83	264	673239	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.49	96	5089	2.14	ng/ul	0.00
5) Phenol-d5	7.22	99	146985	19.32	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	108017	22.50	ng/ul	0.00
9) 2-Chlorophenol-d4	7.59	132	117361	19.10	ng/ul	0.00
13) 4-Methylphenol-d8	8.76	113	107603	18.10	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	62932	19.04	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	49646	18.25	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.48	165	113207	18.66	ng/ul	0.00
29) 4-Chloroaniline-d4	10.88	131	106	0.01	ng/ul	-0.11
43) Dimethylphthalate-d6	14.06	166	391107	23.18	ng/ul	0.00
46) Acenaphthylene-d8	14.36	160	527273	23.28	ng/ul	0.00
51) 4-Nitrophenol-d4	14.85	143	47196	14.20	ng/ul	0.00
57) Fluorene-d10	15.65	176	392807	23.61	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.76	200	34784	11.97	ng/ul	0.00
70) Anthracene-d10	17.49	188	588282	20.93	ng/ul	0.00
78) Pyrene-d10	19.77	212	690933	23.91	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.61	264	640389	21.58	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.24	94	12231	1.55 ng/ul	94
44) Dimethylphthalate	14.11	163	200135	12.10 ng/ul#	98
69) Phenanthrene	17.44	178	185469	5.65 ng/ul	99
76) Fluoranthene	19.44	202	369860	9.99 ng/ul	99
79) Pyrene	19.80	202	343518	9.59 ng/ul	98
82) Benzo(a)anthracene	21.62	228	225720	5.87 ng/ul	97
83) Bis(2-ethylhexyl)phthalate	21.52	149	314178	17.17 ng/ul	100
84) Chrysene	21.69	228	261385	7.11 ng/ul	97
87) Benzo(b)fluoranthene	23.82	252	240432	6.44 ng/ul	98
88) Benzo(k)fluoranthene	23.87	252	99920m>	2.75 ng/ul	
90) Benzo(a)pyrene	24.68	252	141878	3.87 ng/ul	98
91) Indeno(1,2,3-cd)pyrene	28.45	276	83600	1.89 ng/ul	93
92) Dibenzo(a,h)anthracene	28.49	278	37578m>	1.03 ng/ul	
93) Benzo(g,h,i)perylene	29.58	276	79883	2.17 ng/ul#	91

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

SJ 3/13/17