

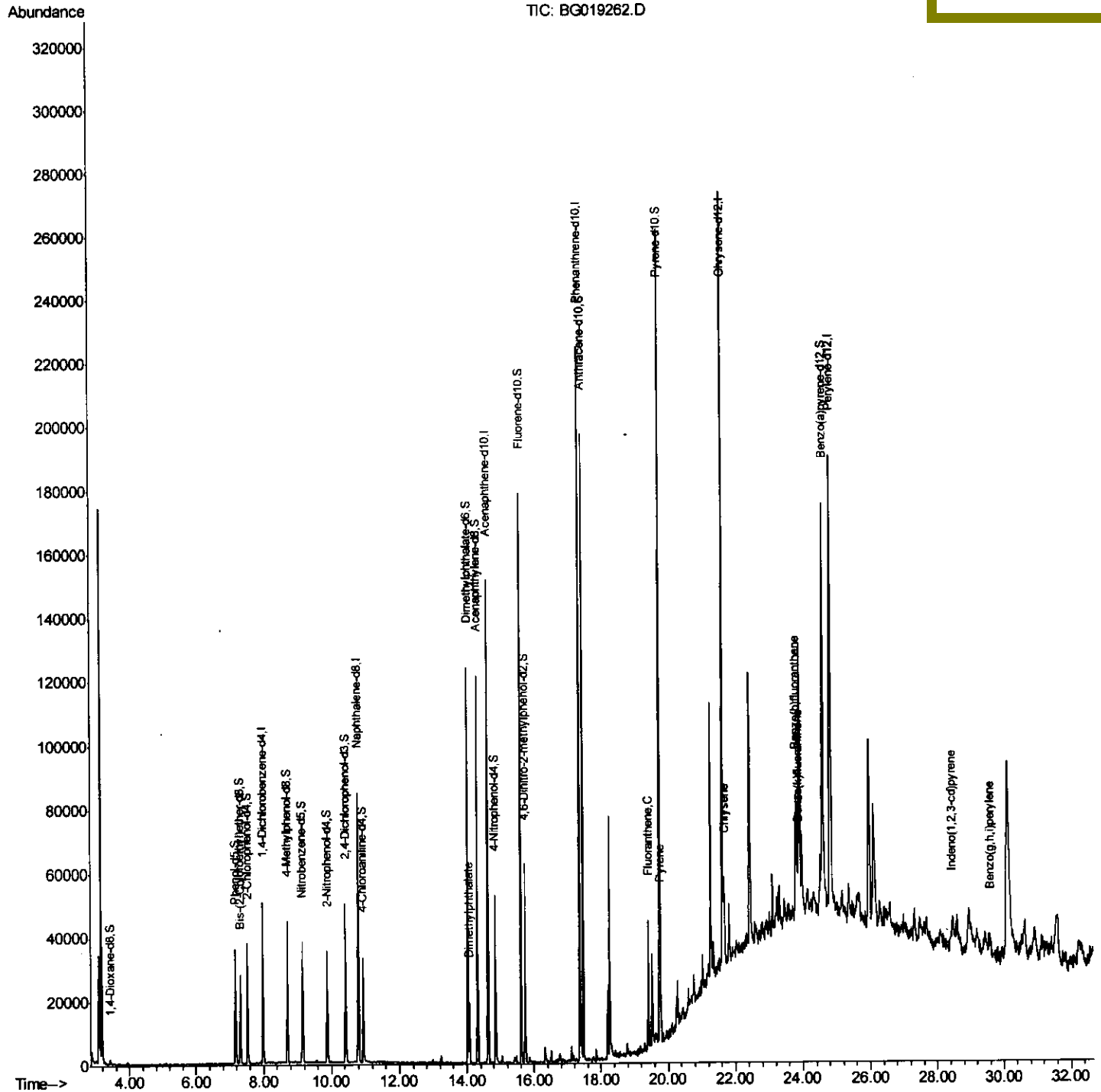
Data Path : Z:\HPCHEM1\BNA G\DATA\BG102015\
Data File : BG019262.D
Acq On : 20 Oct 2015 18:14
Operator : UM/NP
Sample : G4074-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Oct 21 02:06:18 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 21 01:23:32 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
JG9J8

Manual Integrations
APPROVED

MMdadoda
10/21/2015 2:33:01 PM



Quantitation Report (Qedit)

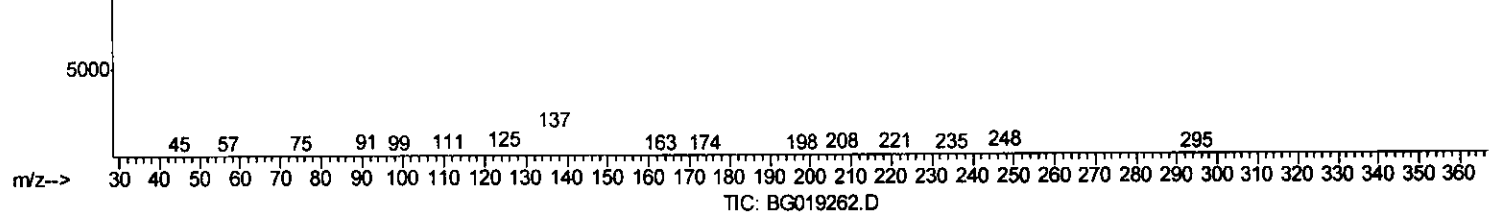
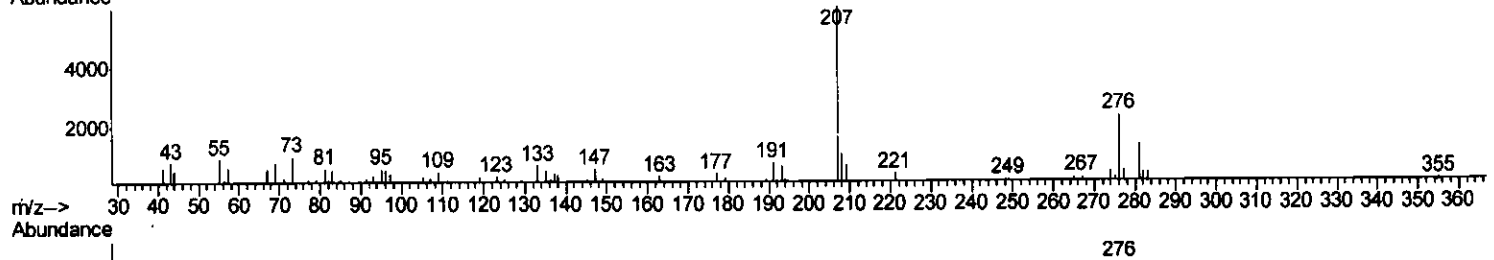
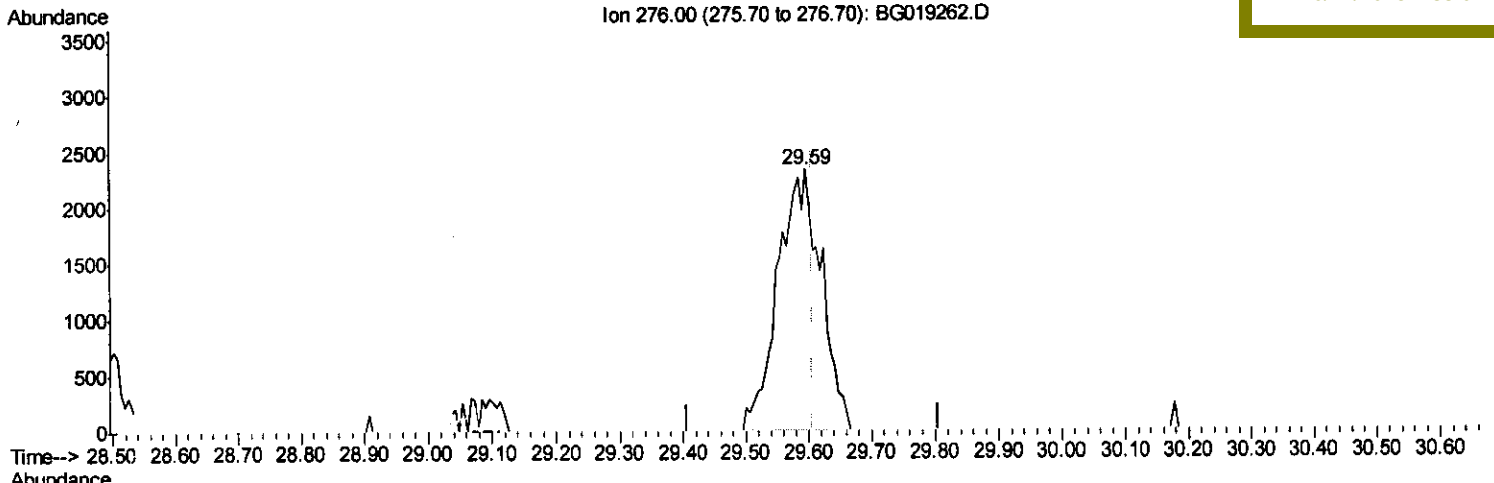
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Instrument :
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 ClientSampleId :
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Quant Time: Oct 21 02:03:41 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 21 01:23:32 2015
 Response via : Initial Calibration

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

29.595min (-0.011) 1.47ng/ul m

response 11074

Ion	Exp%	Act%
276.00	100	100
138.00	20.40	17.06
277.00	23.40	22.43
0.00	0.00	0.00

U.M
 10/31/15

Quantitation Report (Qedit)

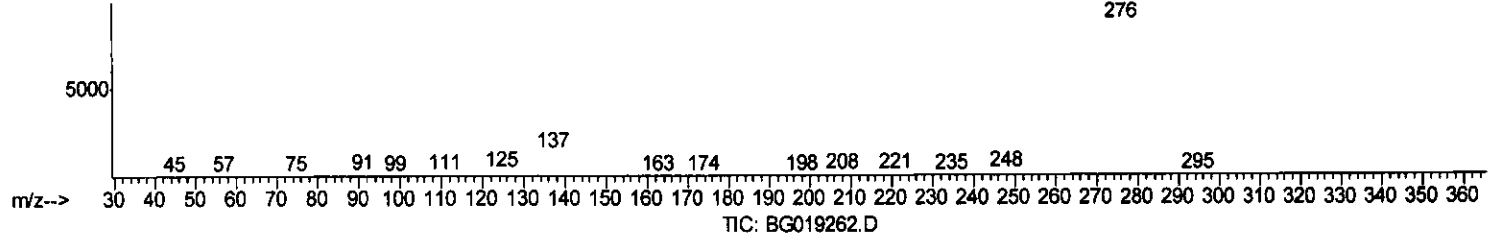
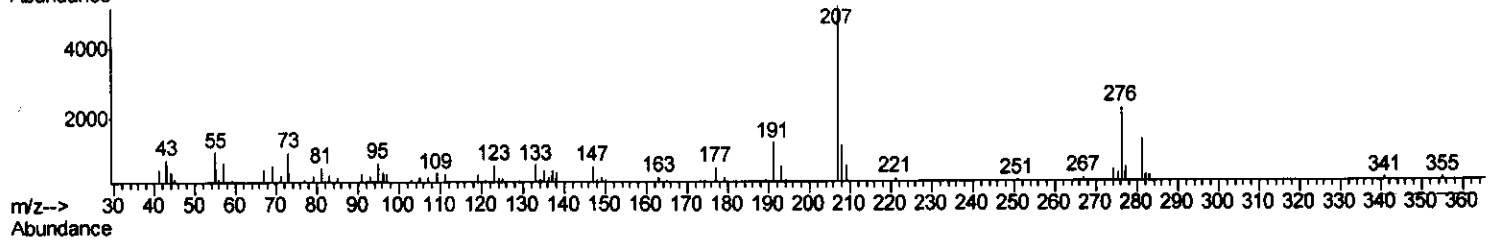
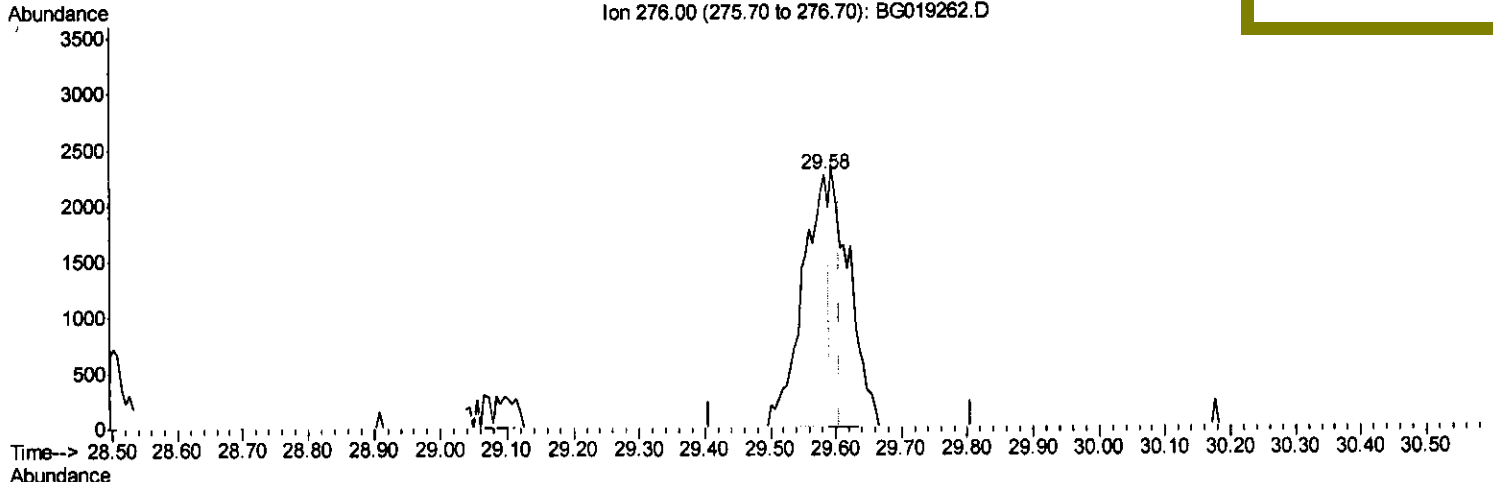
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(94) Benzo(g,h,i)perylene

29.583min (-0.023) 0.84ng/ul

response 6309

Ion	Exp%	Act%
276.00	100	100
138.00	20.40	19.27
277.00	23.40	24.81
0.00	0.00	0.00

Quantitation Report (Qedit)

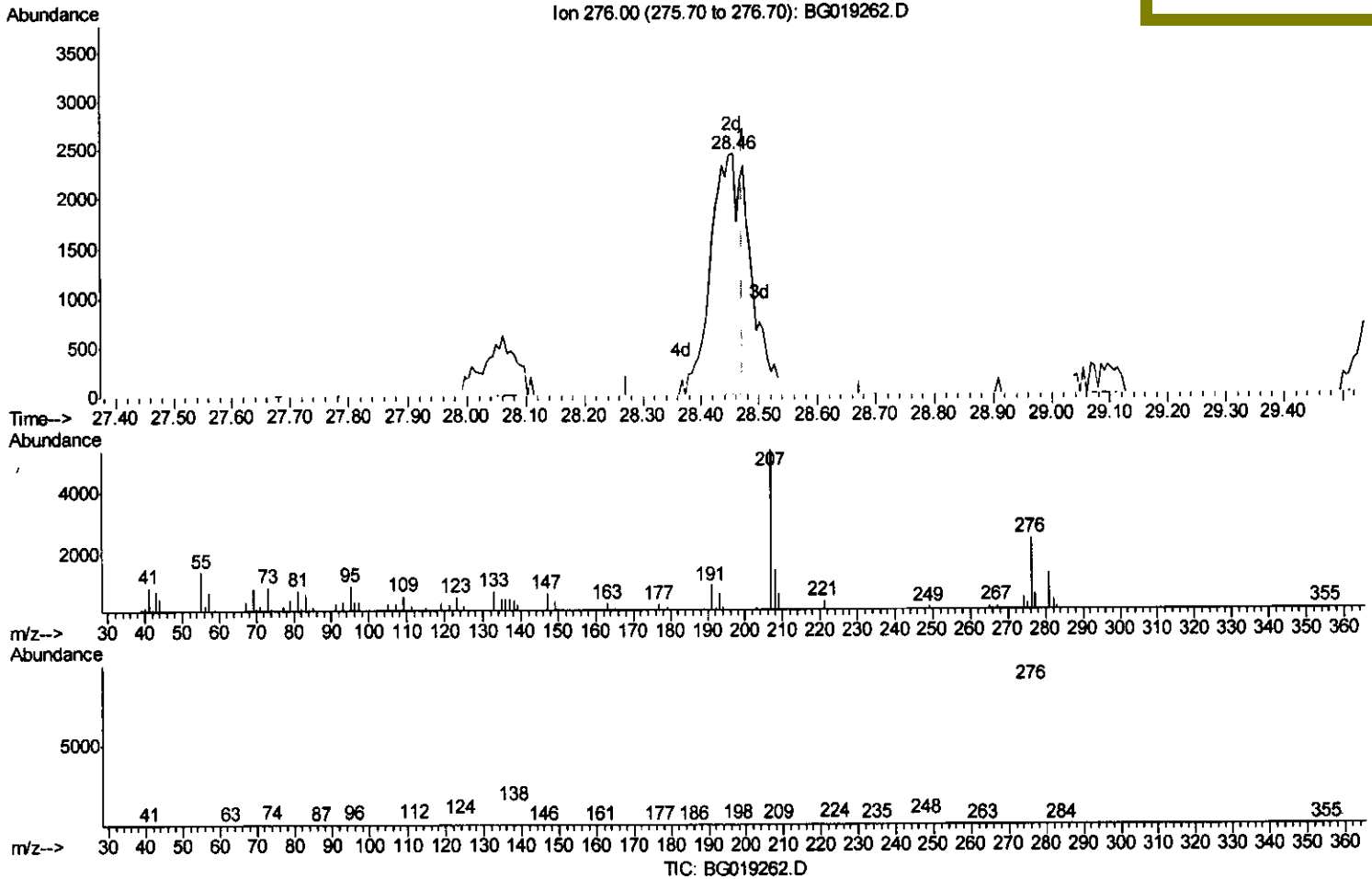
Data Path : Z:\HPCHEM1\BNA G\DATA\BG102015\
 Data File : BG019262.D
 Acq On : 20 Oct 2015 18:14
 Operator : UM/NP
 Sample : G4074-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JG9J8

Quant Time: Oct 21 02:03:41 2015
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Manual Integrations
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(92) Indeno(1,2,3-cd)pyrene

28.455min (-0.017) 1.24ng/ul m U.M

response 11382

10/31/15

Ion	Exp%	Act%
276.00	100	100
138.00	18.60	19.58
277.00	22.50	26.35
0.00	0.00	0.00

Quantitation Report (Qedit)

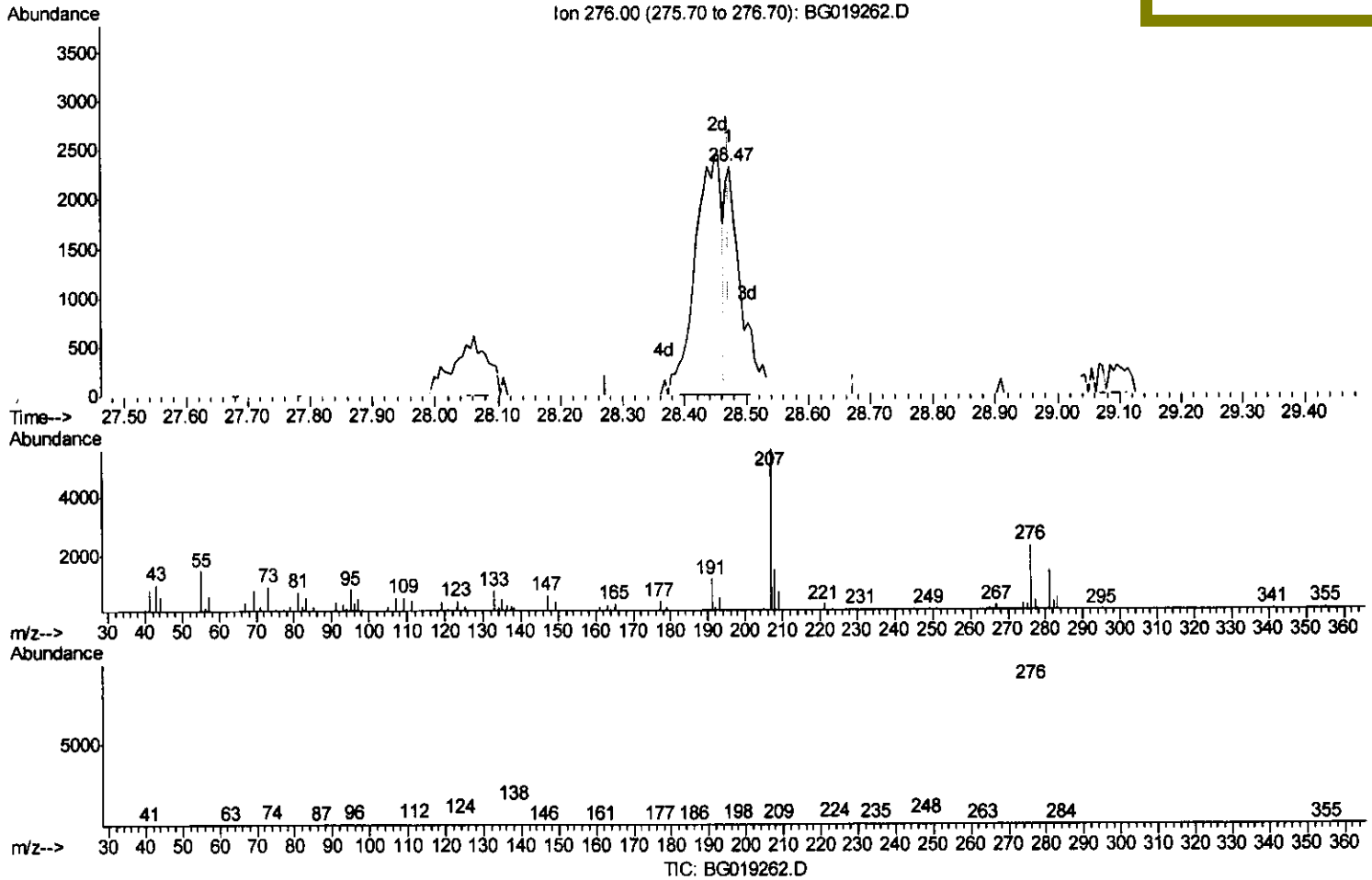
Data Path : Z:\HPCHEM1\BNA G\DATA\BG102015\
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 Operator : UM/NP
 Sample : G4074-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Quant Time: Oct 21 02:03:41 2015
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Manual Integrations
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(92) Indeno(1,2,3-cd)pyrene

28.473min (+0.001) 0.32ng/ul

response 2975

Ion	Exp%	Act%
276.00	100	100
138.00	18.60	21.25
277.00	22.50	20.39
0.00	0.00	0.00

Quantitation Report (Qedit)

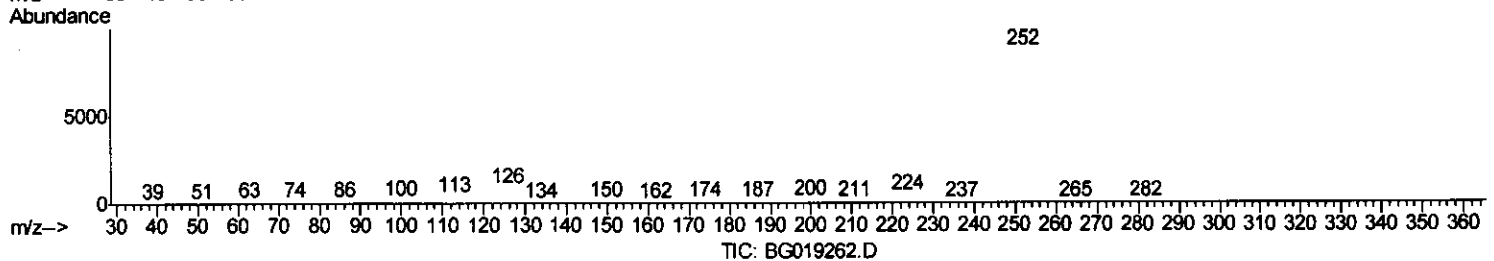
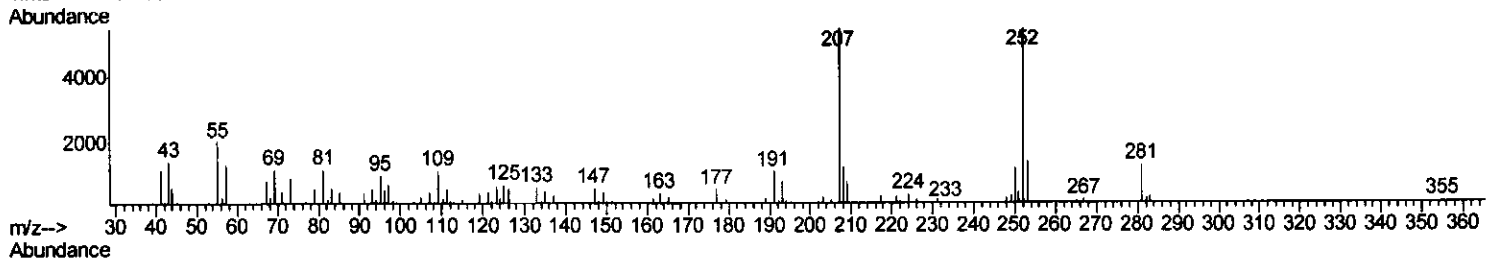
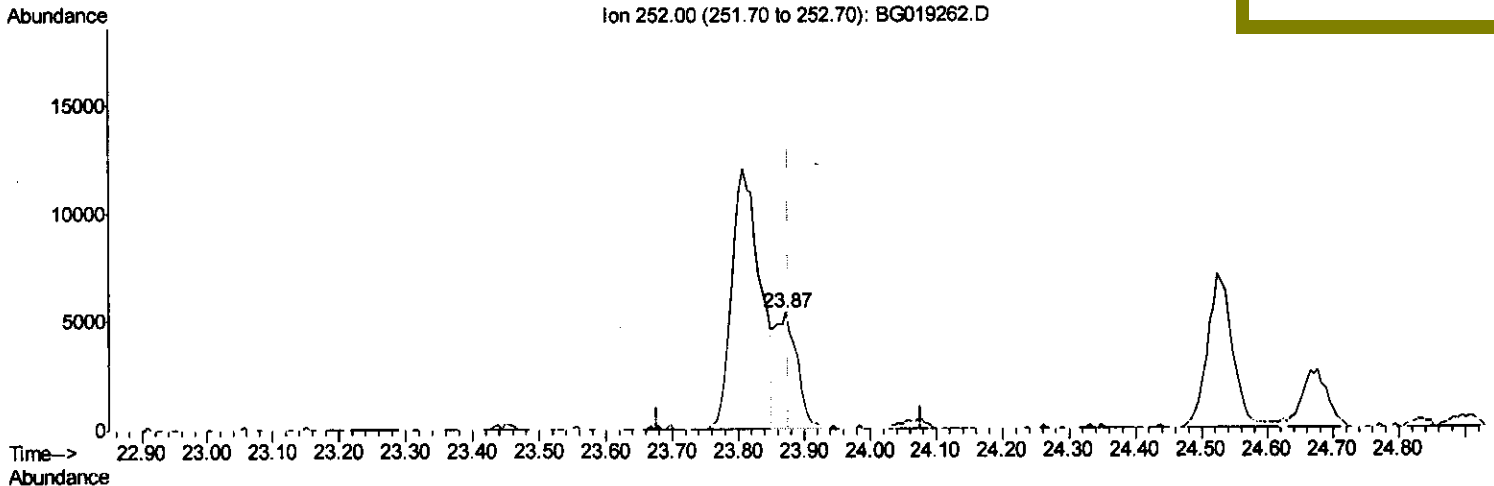
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 Data File : BG019262.D
 Acq On : 20 Oct 2015 18:14
 Operator : UM/NP
 Sample : G4074-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JG9J8

Quant Time: Oct 21 01:25:22 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 21 01:23:32 2015
 Response via : Initial Calibration

Manual Integrations
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(89) Benzo(k)fluoranthene

23.872min (-0.005) 1.59ng/ul m

U.M
 10/31/15

response 12517

Ion	Exp%	Act%
252.00	100	100
253.00	21.50	25.86#
125.00	10.50	13.18#
0.00	0.00	0.00

Quantitation Report (Qedit)

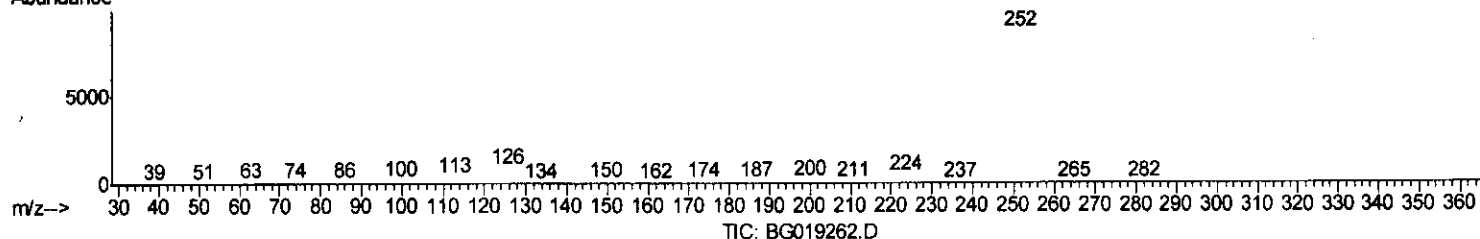
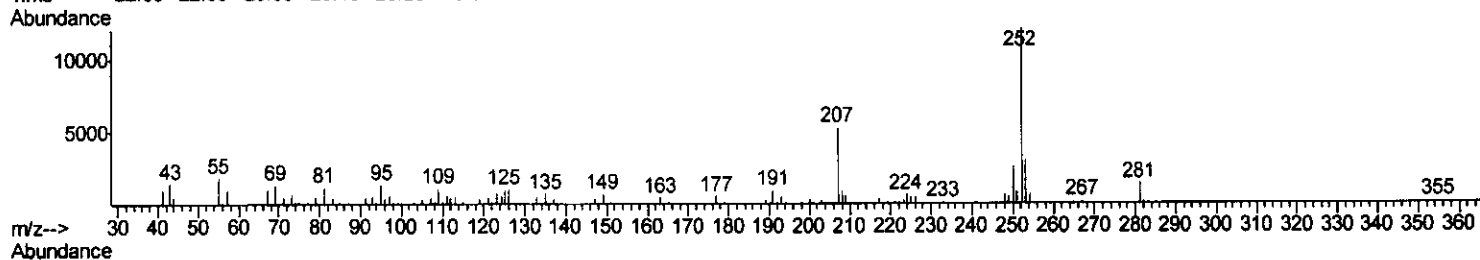
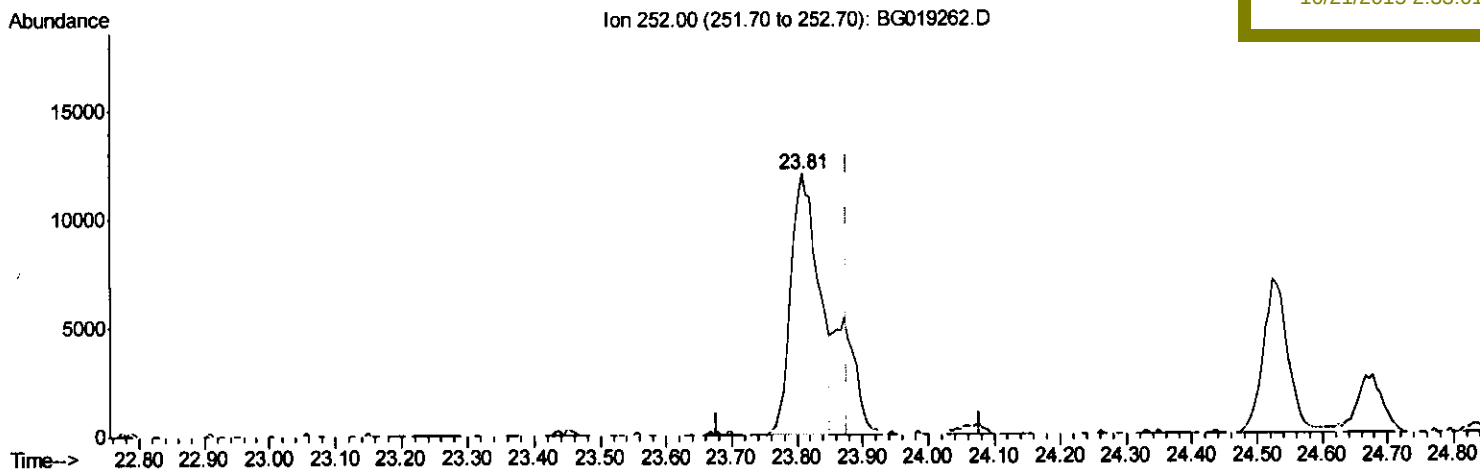
Data Path : Z:\HPCHEM1\BNA G\DATA\BG102015\
 Data File : BG019262.D
 Acq On : 20 Oct 2015 18:14
 Operator : UM/NP
 Sample : G4074-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JG9J8

Quant Time: Oct 21 01:25:22 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 21 01:23:32 2015
 Response via : Initial Calibration

Manual Integrations
 APPROVED

MMdadoda
 10/21/2015 2:33:01 PM



(89) Benzo(k)fluoranthene
 23.808min (-0.070) 4.50ng/ul
 response 35506

Ion	Exp%	Act%
252.00	100	100
253.00	21.50	25.53
125.00	10.50	9.39
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102015\
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Manual Integrations
 APPROVED

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	13588	20.00	ng/ul	0.00
18) Naphthalene-d8	10.79	136	64707	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.62	164	49619	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.37	188	128526	20.00	ng/ul	0.00
78) Chrysene-d12	21.63	240	139596	20.00	ng/ul	0.00
86) Perylene-d12	24.82	264	141617	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.42	96	609	2.50	ng/uL	0.00
5) Phenol-d5	7.17	99	18796	15.85	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.32	67	13300	17.43	ng/ul	0.00
9) 2-Chlorophenol-d4	7.53	132	15214	17.60	ng/ul	0.00
13) 4-Methylphenol-d8	8.71	113	15646	15.57	ng/ul	0.00
19) Nitrobenzene-d5	9.15	128	8093	16.23	ng/ul	0.00
22) 2-Nitrophenol-d4	9.88	143	9655	18.25	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.42	165	18197	17.36	ng/ul	0.00
29) 4-Chloroaniline-d4	10.93	131	17351	12.86	ng/ul	0.00
44) Dimethylphthalate-d6	14.03	166	76437	17.61	ng/ul	0.00
47) Acenaphthylene-d8	14.32	160	87673	18.69	ng/ul	0.00
52) 4-Nitrophenol-d4	14.84	143	11327	14.52	ng/ul	0.00
58) Fluorene-d10	15.61	176	70757	19.08	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	12469	15.73	ng/ul	0.00
71) Anthracene-d10	17.47	188	113691	18.72	ng/ul	0.00
79) Pyrene-d10	19.75	212	128573	20.31	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.61	264	132303	18.49	ng/ul	0.00

Target Compounds

					Ovalue
45) Dimethylphthalate	14.07	163	11684	2.67 ng/ul#	97
77) Fluoranthene	19.42	202	23116	2.57 ng/ul#	89
80) Pvrene	19.78	202	20271	2.45 ng/ul#	87
85) Chrysene	21.67	228	21374	2.87 ng/ul	98
88) Benzo(b)fluoranthene	23.81	252	35506	4.42 ng/ul	94
89) Benzo(k)fluoranthene	23.87	252	12517m	1.59 ng/ul	
92) Indeno(1,2,3-cd)pvrene	28.46	276	11382m	1.24 ng/ul	
94) Benzo(a,h,i)perylene	29.59	276	11074m	1.47 ng/ul	

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

UM
 10/31/15