

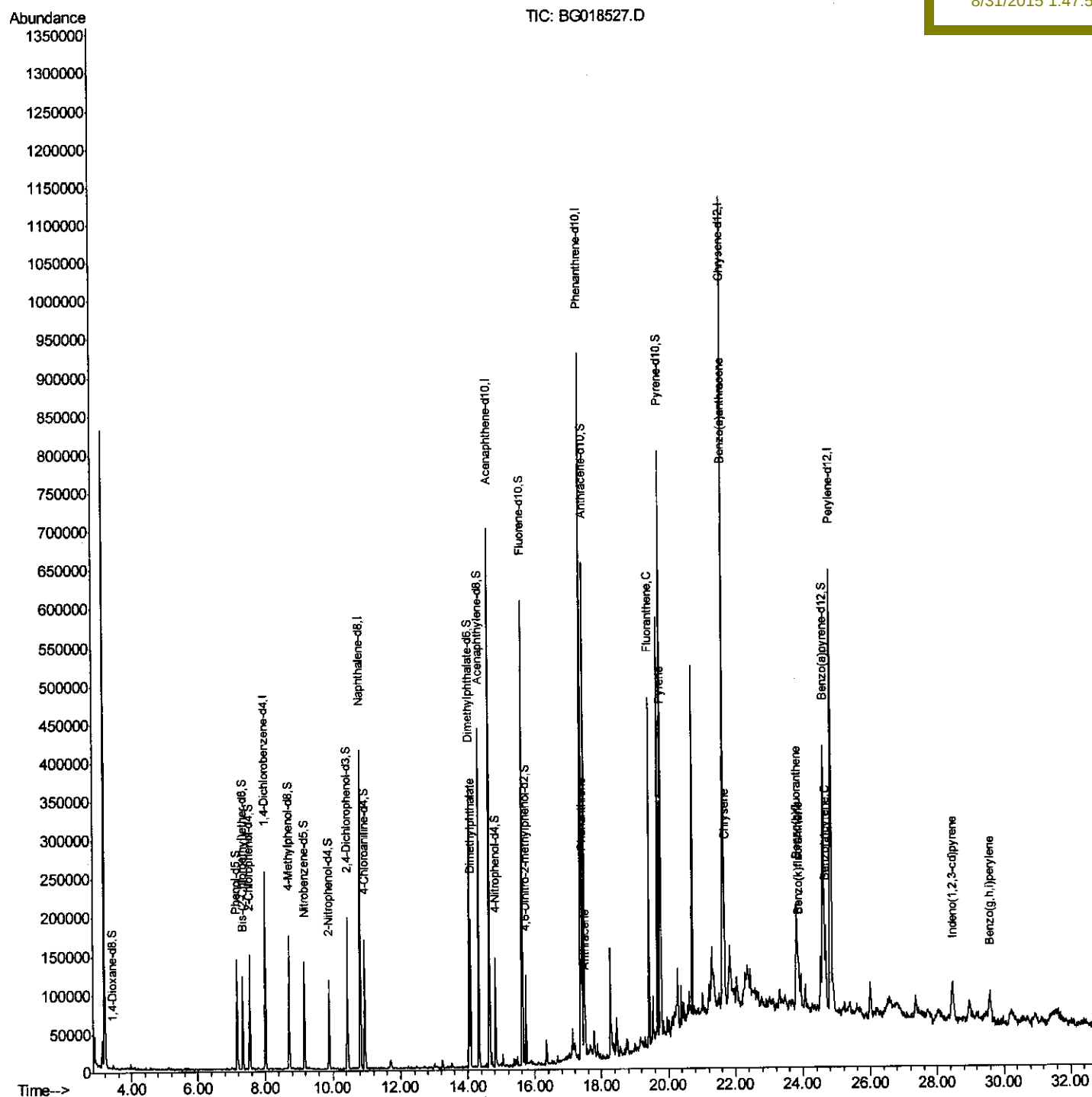
Data Path : Z:\HPCHEM1\BNA G\DATA\BG082815\
Data File : BG018527.D
Acq On : 29 Aug 2015 00:09
Operator : UM/MA
Sample : G3448-11
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AD0

Quant Time: Aug 31 04:29:25 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 31 03:29:45 2015
Response via : Initial Calibration

Manual Integrations
APPROVED

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8/31/2015 1:47:51 PM



Quantitation Report (Qedit)

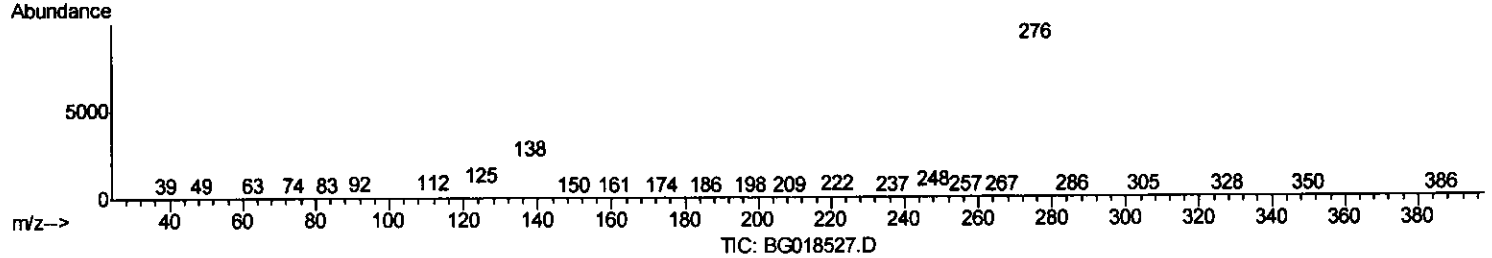
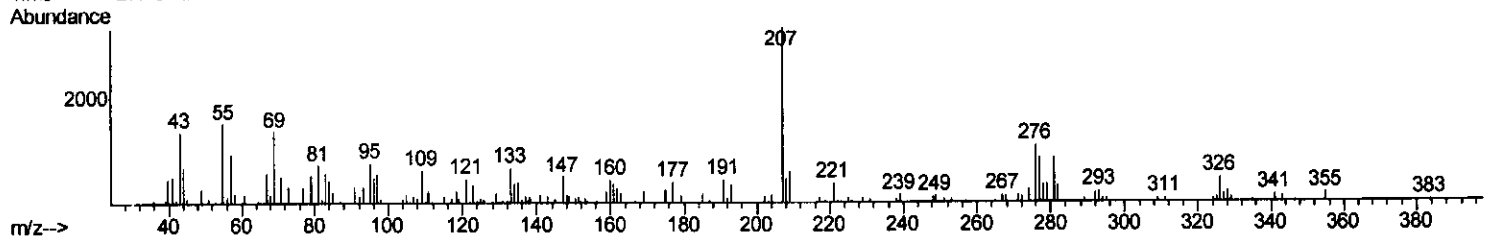
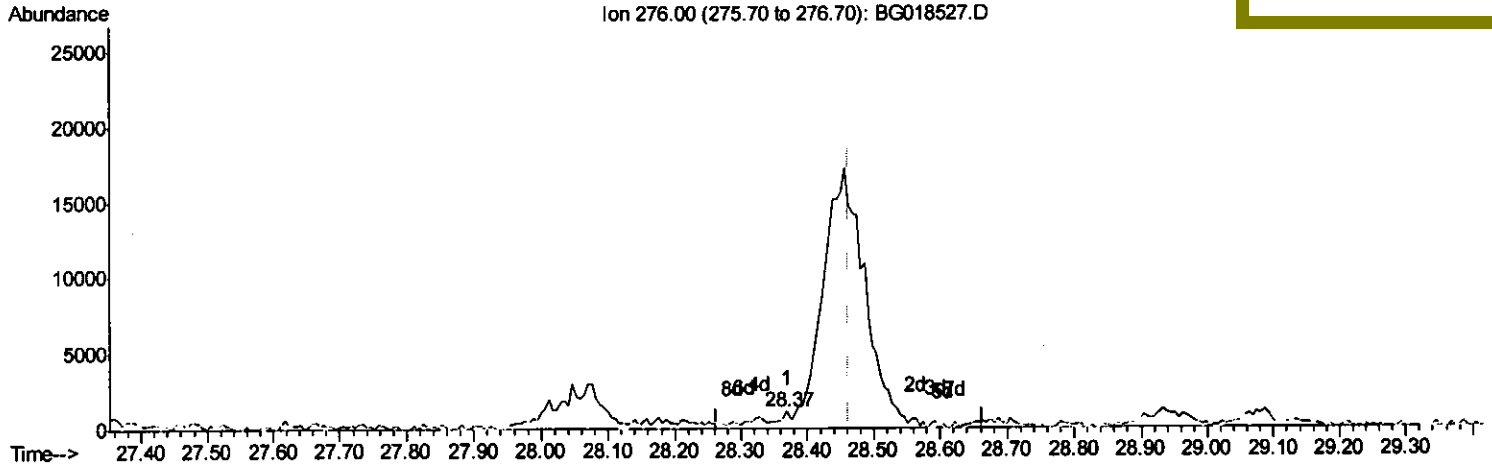
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(92) Indeno(1,2,3-cd)pyrene

28.369min (-0.094) 0.01ng/ul

response 424

Ion	Exp%	Act%
276.00	100	100
138.00	19.70	41.60#
277.00	23.60	81.03#
0.00	0.00	0.00

Quantitation Report (Qedit)

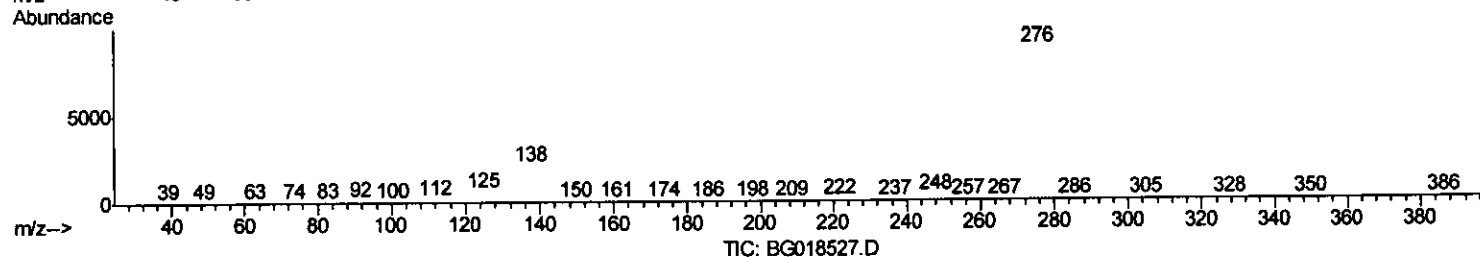
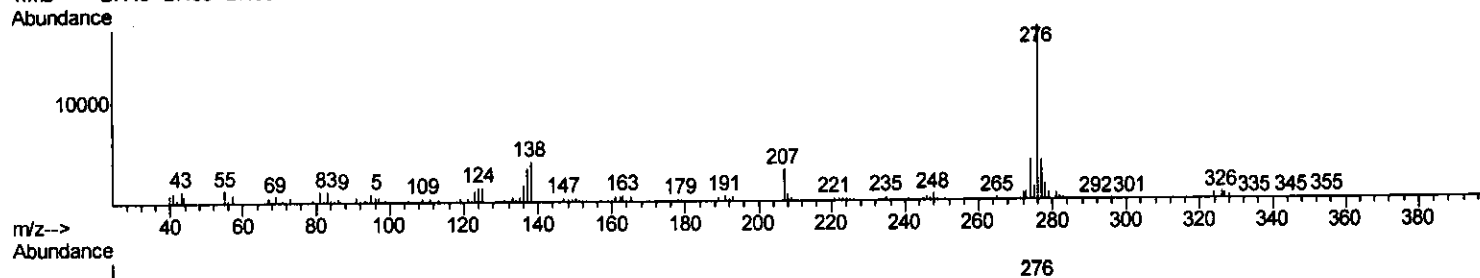
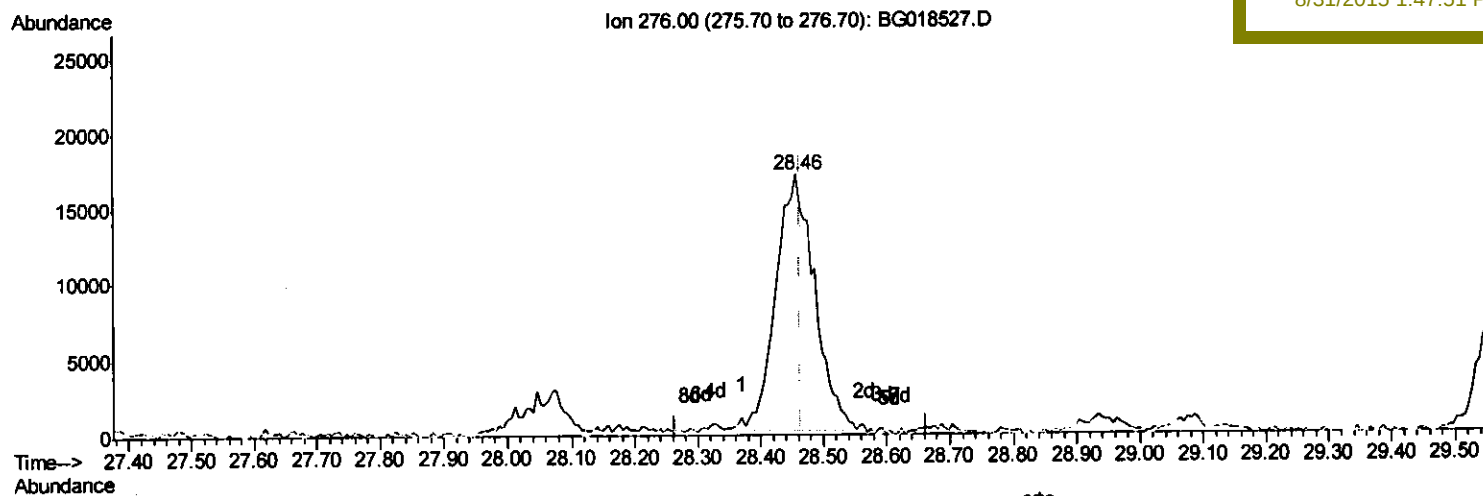
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(92) Indeno(1,2,3-cd)pyrene

28.457min (-0.006) 1.85ng/ul m

response 71635

Ion	Exp%	Act%
276.00	100	100
138.00	19.70	23.50
277.00	23.60	23.02
0.00	0.00	0.00

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09/01/15

Quantitation Report (Qedit)

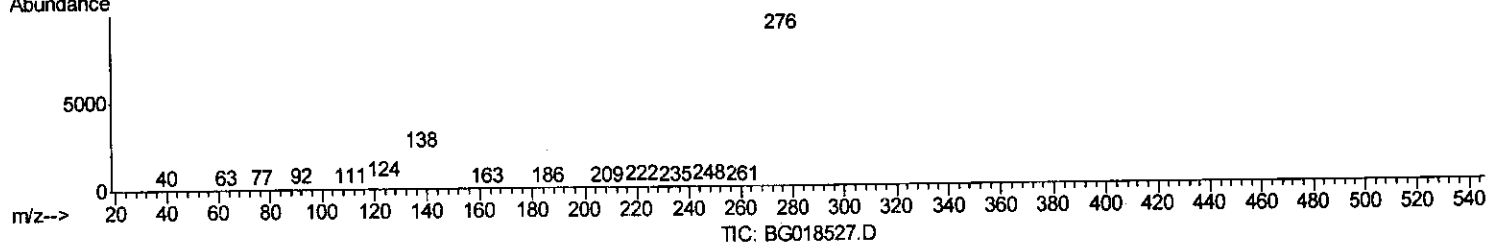
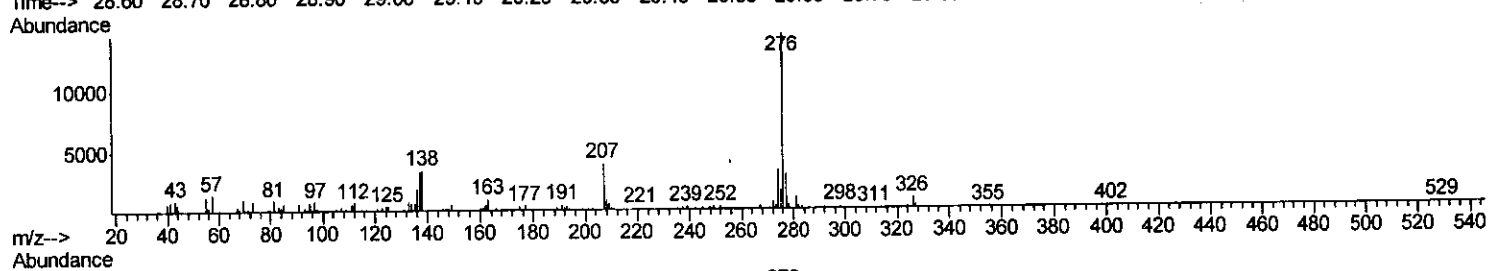
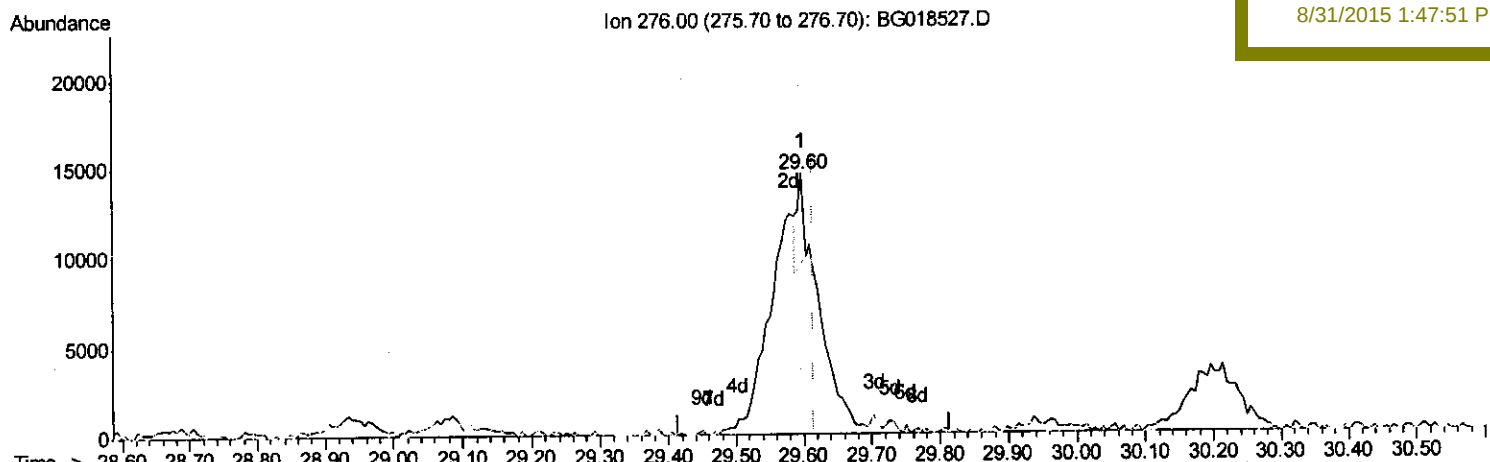
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Instrument :
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 ClientSampleId :
 C0AD0

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

29.597min (-0.018) 0.09ng/ui

response 3047

Ion	Exp%	Act%
276.00	100	100
138.00	20.60	23.55
277.00	23.80	21.04
0.00	0.00	0.00

Quantitation Report (Qedit)

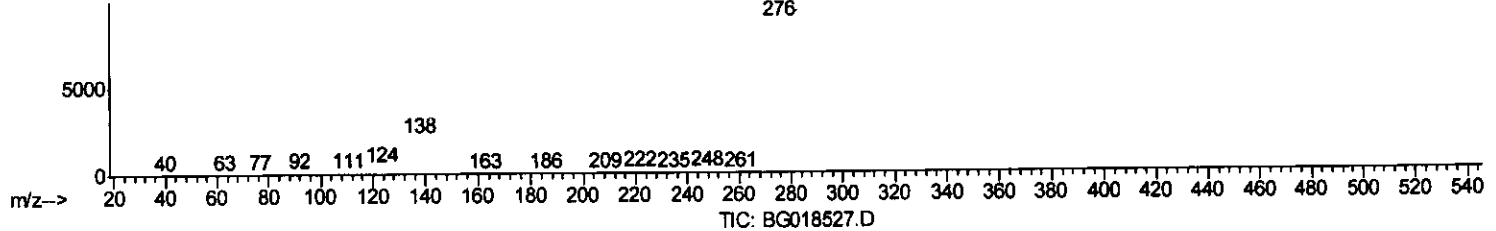
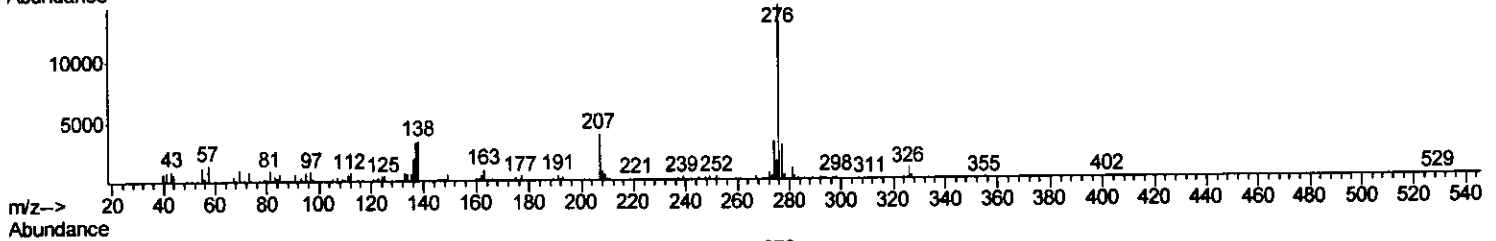
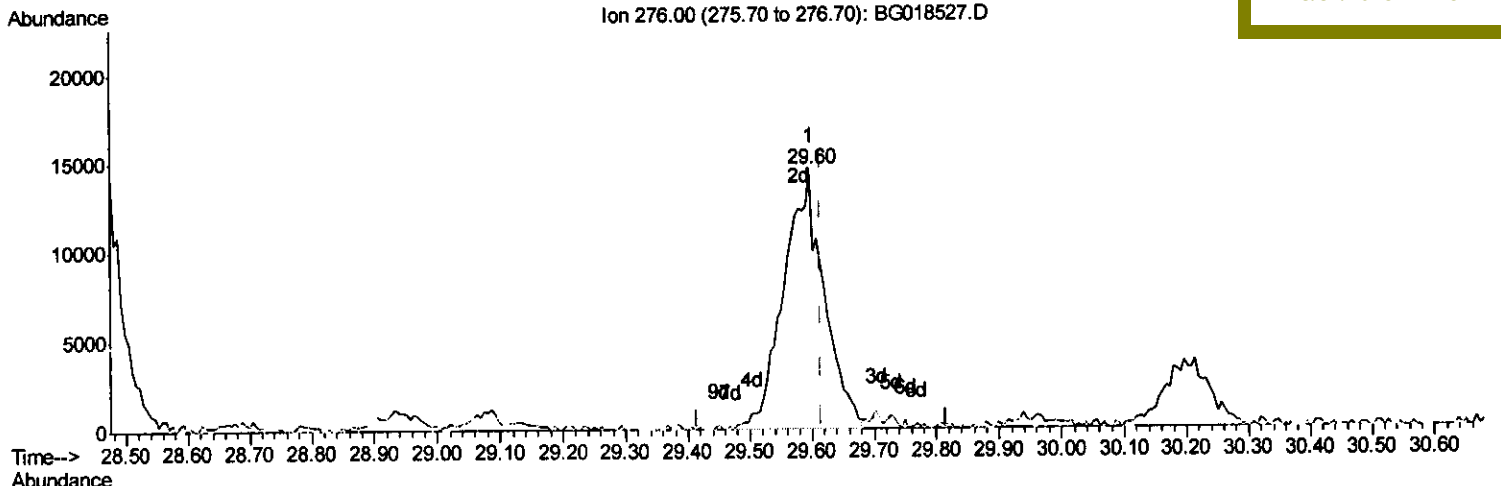
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Instrument :
 BNA_G
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Manual Integrations
 APPROVED

MMDadoda
 8/31/2015 1:47:51 PM



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

29.597min (-0.018) 2.03ng/ul m

response 65901

Ion	Exp%	Act%
276.00	100	100
138.00	20.60	23.55
277.00	23.80	21.04
0.00	0.00	0.00

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 09/01/15

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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	8.01	152	77257	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	351693	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.64	164	231281	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.38	188	564507	20.00	ng/ul	0.00
78) Chrysene-d12	21.64	240	554314	20.00	ng/ul	0.00
86) Perylene-d12	24.83	264	559083	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.45	96	2039	1.17	ng/uL	0.00
5) Phenol-d5	7.16	99	94133	13.42	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.34	67	57028	13.12	ng/ul	0.00
9) 2-Chlorophenol-d4	7.54	132	71700	13.35	ng/ul	0.00
13) 4-Methylphenol-d8	8.71	113	70641	11.99	ng/ul	0.00
19) Nitrobenzene-d5	9.16	128	37101	13.53	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	37740	13.52	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.43	165	80973	13.66	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	95502	14.26	ng/ul	0.00
44) Dimethylphthalate-d6	14.04	166	257308	13.22	ng/ul	0.00
47) Acenaphthylene-d8	14.33	160	327473	13.36	ng/ul	0.00
52) 4-Nitrophenol-d4	14.82	143	36978	10.56	ng/ul	0.00
58) Fluorene-d10	15.63	176	234646	13.84	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.74	200	24606	8.86	ng/ul	0.00
71) Anthracene-d10	17.48	188	367518	13.61	ng/ul	0.00
79) Pyrene-d10	19.76	212	360587	12.54	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.61	264	351613	11.85	ng/ul	0.00

Target Compounds

					Ovalue
45) Dimethylphthalate	14.09	163	128287	6.68 ng/ul	99
70) Phenanthrene	17.42	178	126634	4.13 ng/ul	95
72) Anthracene	17.51	178	32556	1.04 ng/ul	100
77) Fluoranthene	19.43	202	273101	8.35 ng/ul	98
80) Pyrene	19.79	202	228653	6.10 ng/ul	96
83) Benzo(a)anthracene	21.62	228	127618	3.71 ng/ul	96
85) Chrysene	21.68	228	113018	3.52 ng/ul	97
88) Benzo(b)fluoranthene	23.82	252	155832	4.44 ng/ul#	95
89) Benzo(k)fluoranthene	23.88	252	59020	1.75 ng/ul	96
91) Benzo(a)pyrene	24.68	252	109848	3.29 ng/ul#	95
92) Indeno(1,2,3-cd)pyrene	28.46	276	71635m	1.85 ng/ul	
94) Benzo(a,h,i)perylene	29.60	276	65901m	2.03 ng/ul	

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

U.M.
 09/01/15