

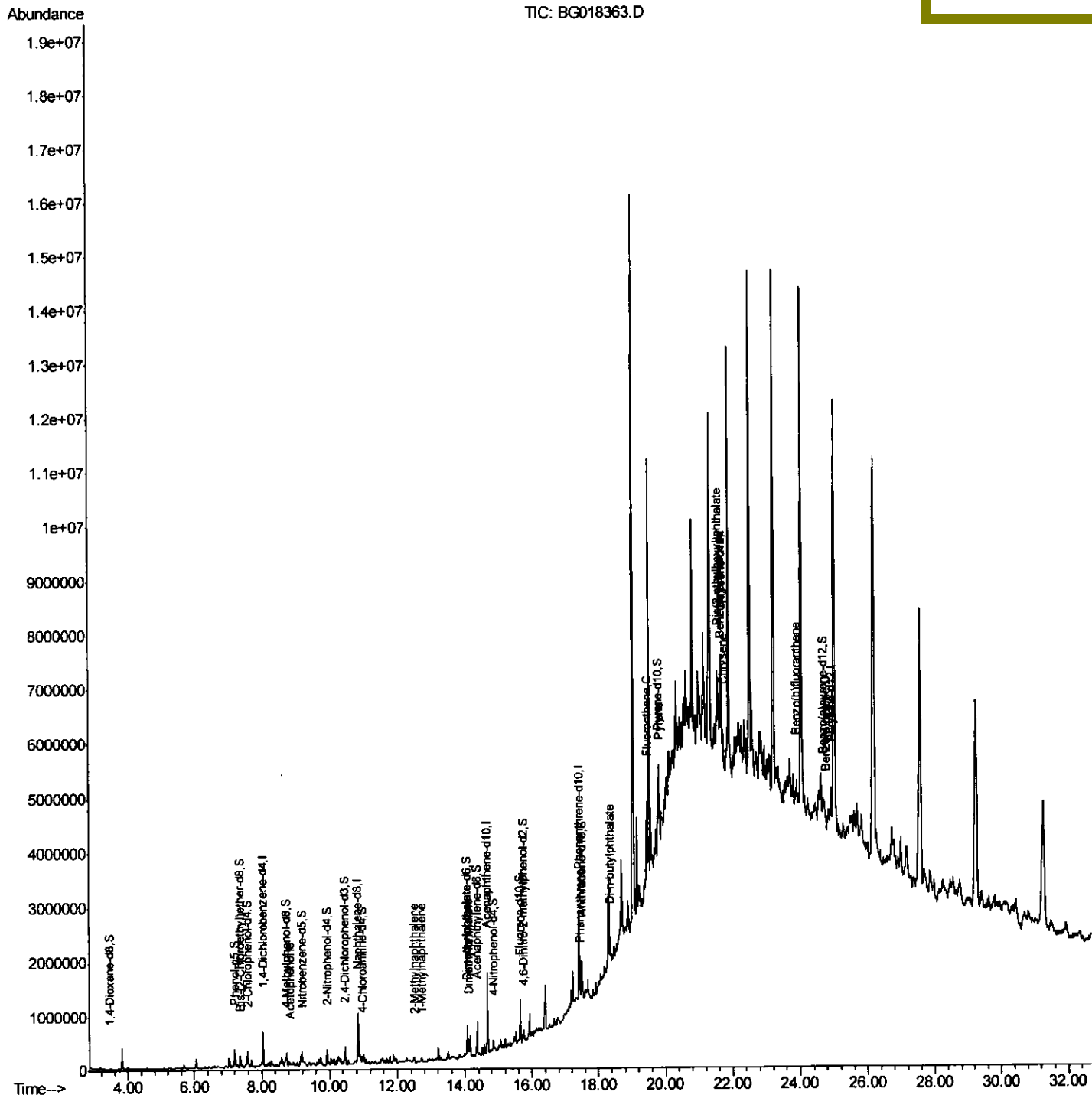
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081015\
Data File : BG018363.D
Acq On : 10 Aug 2015 20:56
Operator : UM/MA
Sample : G3136-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C01T8

Quant Time: Aug 11 02:14:34 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 11 01:39:57 2015
Response via : Initial Calibration

Manual Integrations APPROVED

apatel
8/11/2015 9:13:10 AM



Quantitation Report (Qedit)

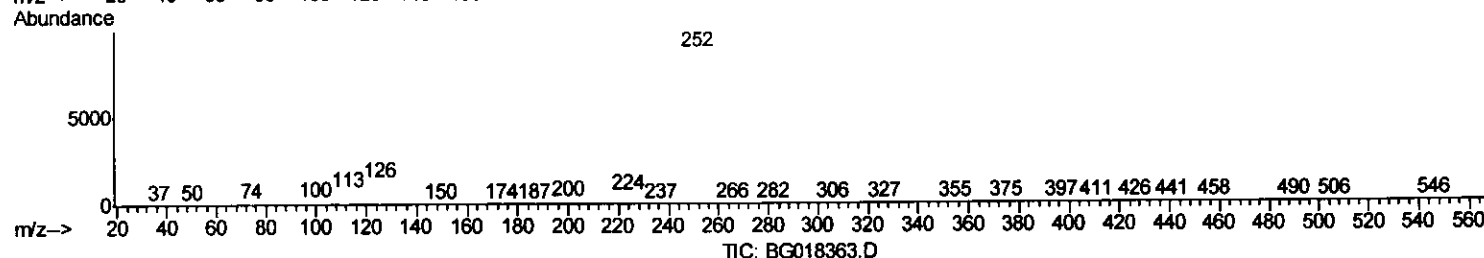
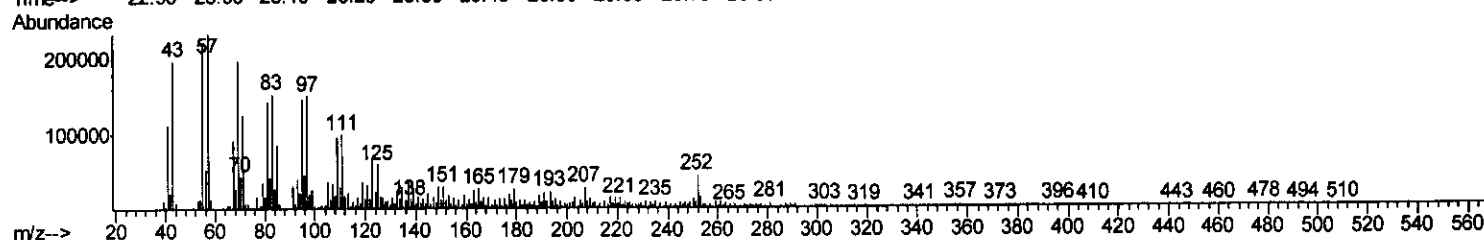
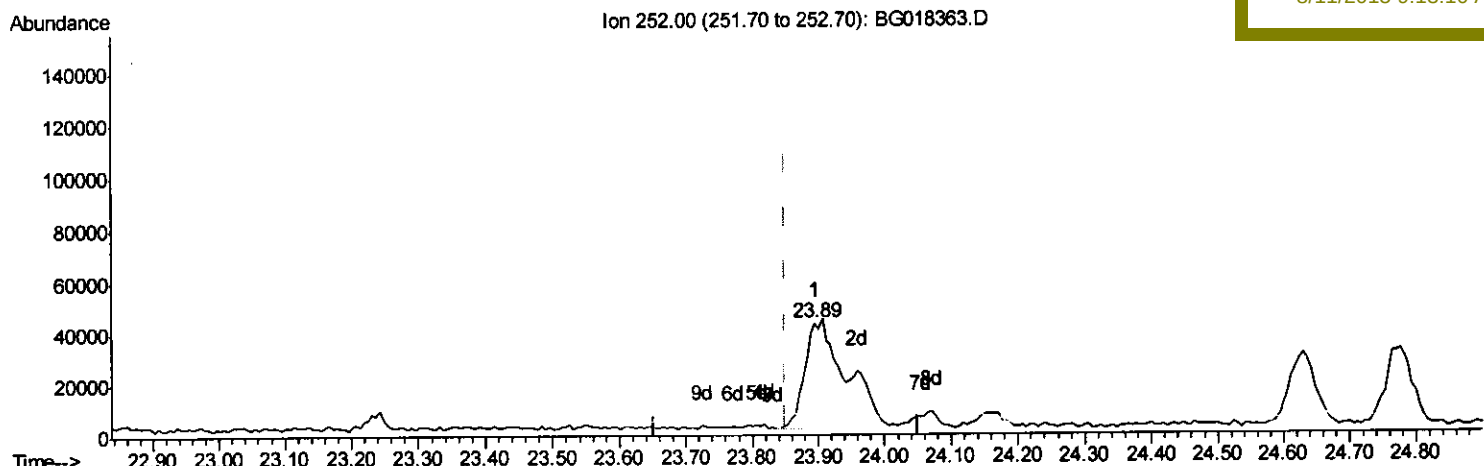
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(88) Benzo(b)fluoranthene

23.895min (+0.044) 1.41ng/ui

response 66748

Ion	Exp%	Act%
252.00	100	100
253.00	22.60	31.94#
125.00	10.40	134.85#
0.00	0.00	0.00

Quantitation Report (Qedit)

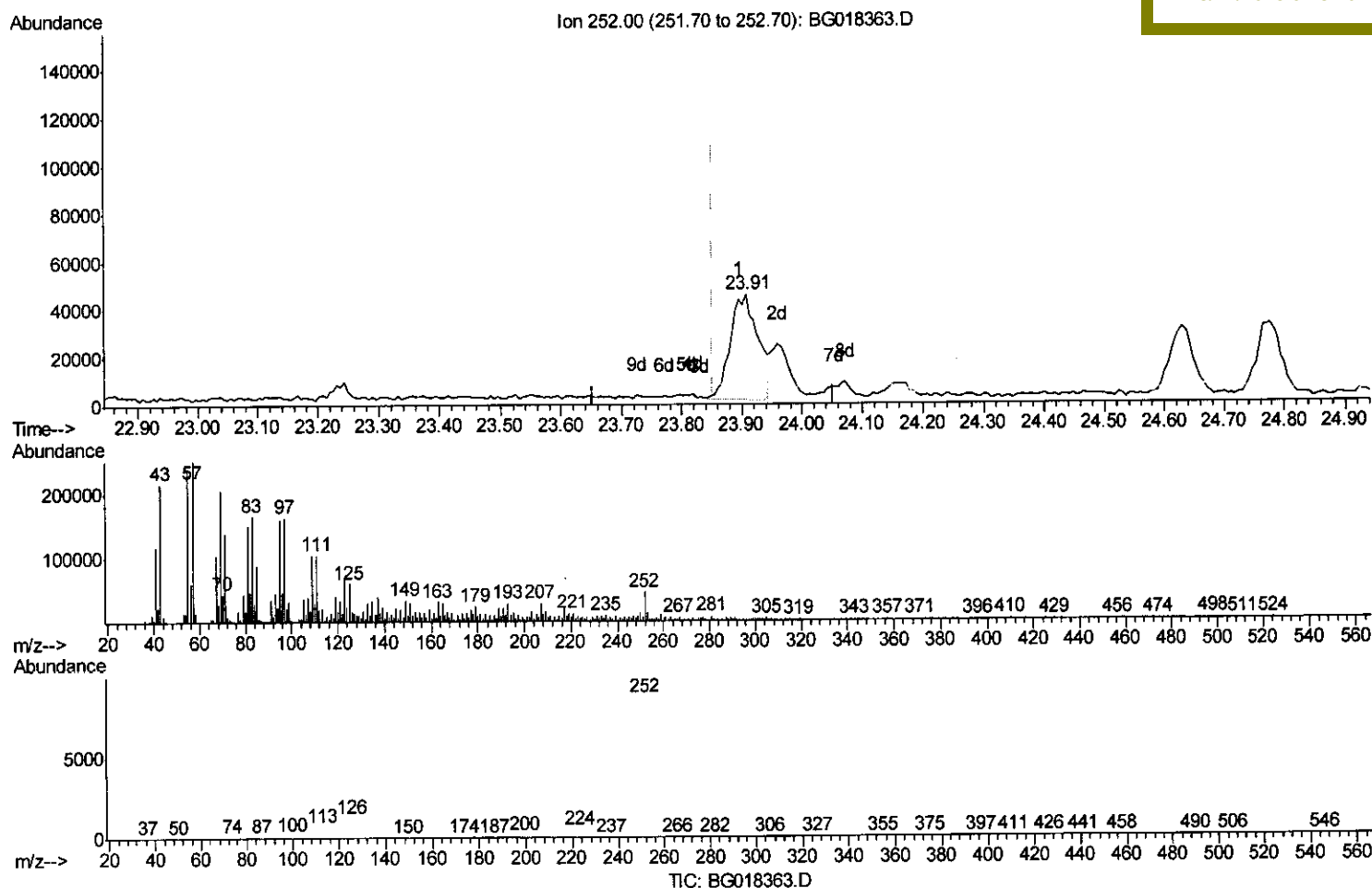
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Manual Integrations
APPROVED

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8/11/2015 9:13:10 AM



(88) Benzo(b)fluoranthene

23.907min (+0.055) 2.96ng/ul m U.M

response 140407

Ion	Exp%	Act%
252.00	100	100
253.00	22.60	32.08#
125.00	10.40	130.61#
0.00	0.00	0.00

8/12/15

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 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 Client Sampled :
 C01T8

Quant Time: Aug 11 02:14:34 2015
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Manual Integrations
 APPROVED

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 8/11/2015 9:13:10 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	162558	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	747784	20.00	ng/ul	0.01
36) Acenaphthene-d10	14.67	164	457701	20.00	ng/ul	0.01
62) Phenanthrene-d10	17.41	188	894969	20.00	ng/ul	0.01
78) Chrysene-d12	21.69	240	757610	20.00	ng/ul	0.03
86) Perylene-d12	24.93	264	753752	20.00	ng/ul	0.06

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	6890	1.87	ng/uL	0.00
5) Phenol-d5	7.19	99	150941	10.23	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.36	67	103216	11.29	ng/ul	0.00
9) 2-Chlorophenol-d4	7.57	132	121594	10.76	ng/ul	0.01
13) 4-Methylphenol-d8	8.74	113	87753	7.08	ng/ul	0.00
19) Nitrobenzene-d5	9.19	128	63949	10.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	68456	11.53	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.46	165	123458	9.80	ng/ul	0.00
29) 4-Chloroaniline-d4	10.97	131	66858	4.70	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	381091	9.90	ng/ul	0.00
47) Acenaphthylene-d8	14.36	160	439949	9.07	ng/ul	0.00
52) 4-Nitrophenol-d4	14.85	143	46136	6.66	ng/ul	0.00
58) Fluorene-d10	15.66	176	306000	9.12	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.76	200	31125	7.07	ng/ul	0.00
71) Anthracene-d10	17.51	188	401966	9.39	ng/ul	0.00
79) Pyrene-d10	19.79	212	379133	9.65	ng/ul	0.01
90) Benzo(a)pyrene-d12	24.71	264	394008	9.85	ng/ul	0.05

Target Compounds

						Qvalue
14) Acetophenone	8.85	105	40633	2.27	ng/ul#	40
34) 2-Methylnaphthalene	12.50	142	39277	1.37	ng/ul	91
35) 1-Methylnaphthalene	12.72	142	32278	1.15	ng/ul#	98
45) Dimethylphthalate	14.11	163	194394	5.12	ng/ul#	97
70) Phenanthrene	17.46	178	136697	2.82	ng/ul#	92
76) Di-n-butylphthalate	18.37	149	64067	1.13	ng/ul#	86
77) Fluoranthene	19.46	202	227887	4.39	ng/ul#	92
80) Pyrene	19.82	202	245983	4.80	ng/ul	99
83) Benzo(a)anthracene	21.67	228	95626	2.04	ng/ul#	73
84) Bis(2-ethylhexyl)phthalate	21.58	149	388401	12.69	ng/ul	96
85) Chrysene	21.73	228	111569	2.54	ng/ul#	81
88) Benzo(b)fluoranthene	23.91	252	140407m	2.96	ng/ul	
91) Benzo(a)pyrene	24.78	252	82993	1.84	ng/ul#	1

U.M
 8/12/15

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