

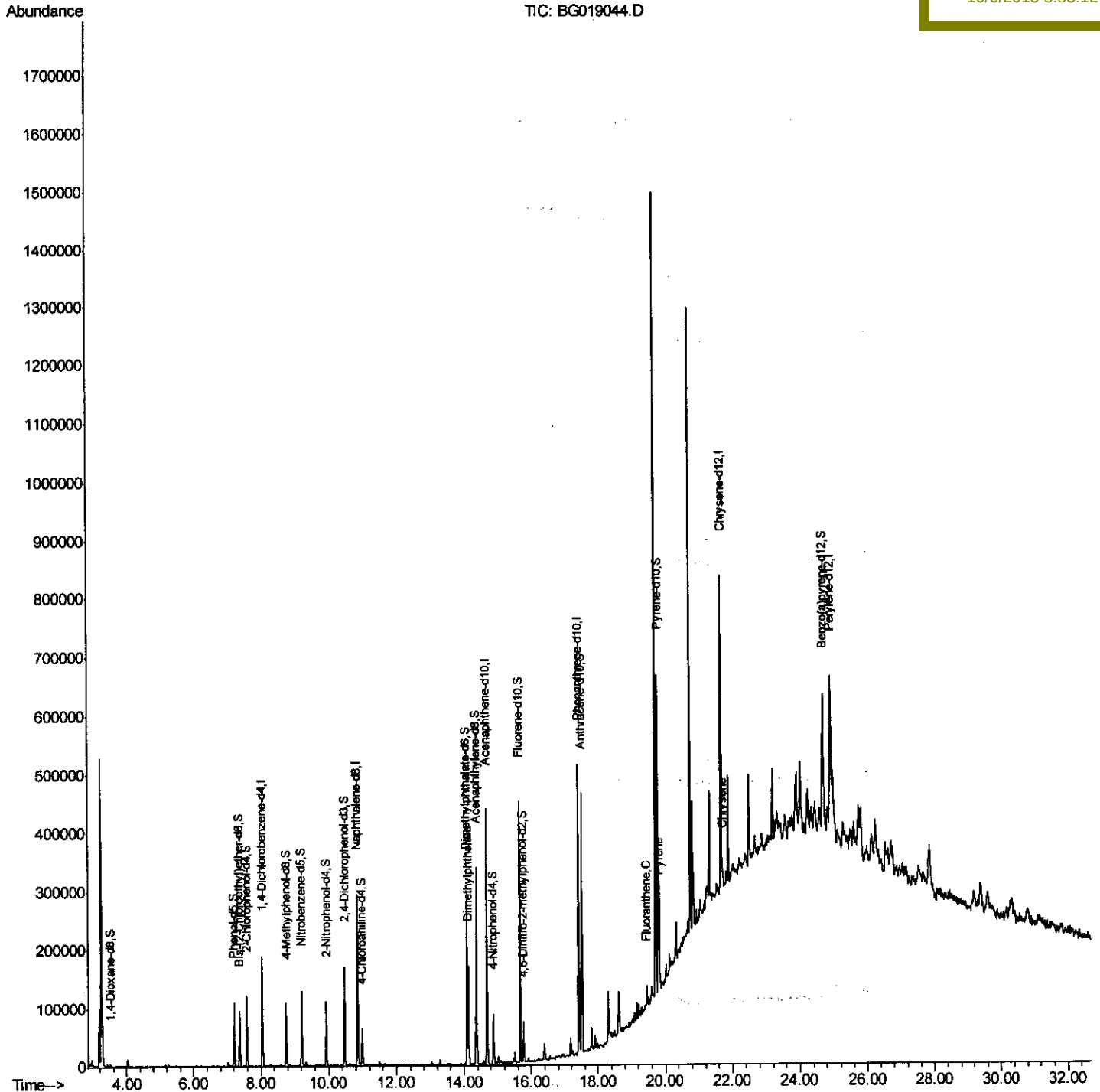
Data Path : Z:\HPCHEM1\BNA_G\Data\BG100515\
Data File : BG019044.D
Acq On : 6 Oct 2015 1:54
Operator : UM/NP
Sample : G3865-10
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E5MY7

Quant Time: Oct 06 09:17:26 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG100415.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 06 04:27:31 2015
Response via : Initial Calibration

Manual Integrations
APPROVED

MMDadoda
10/6/2015 5:55:12 PM



Quantitation Report (Qedit)

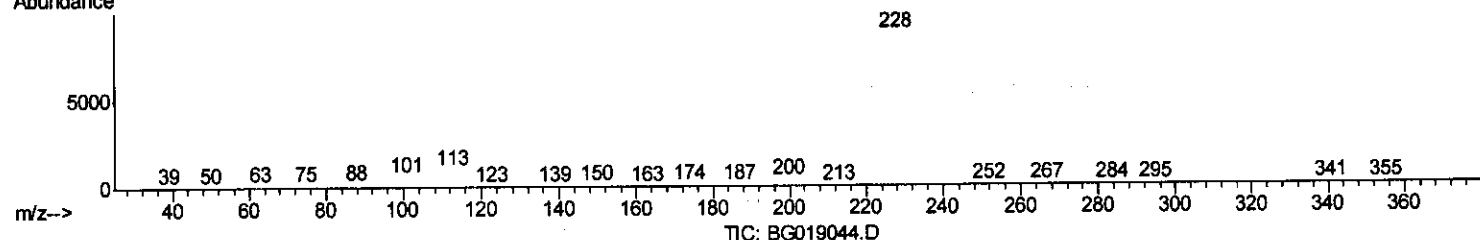
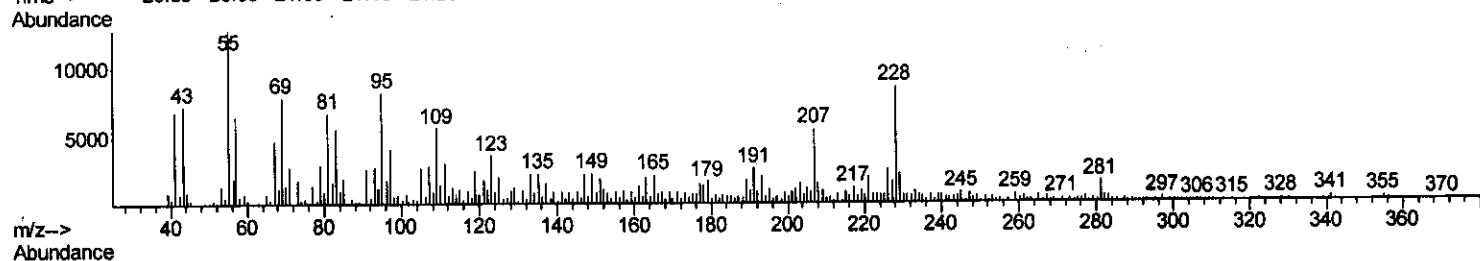
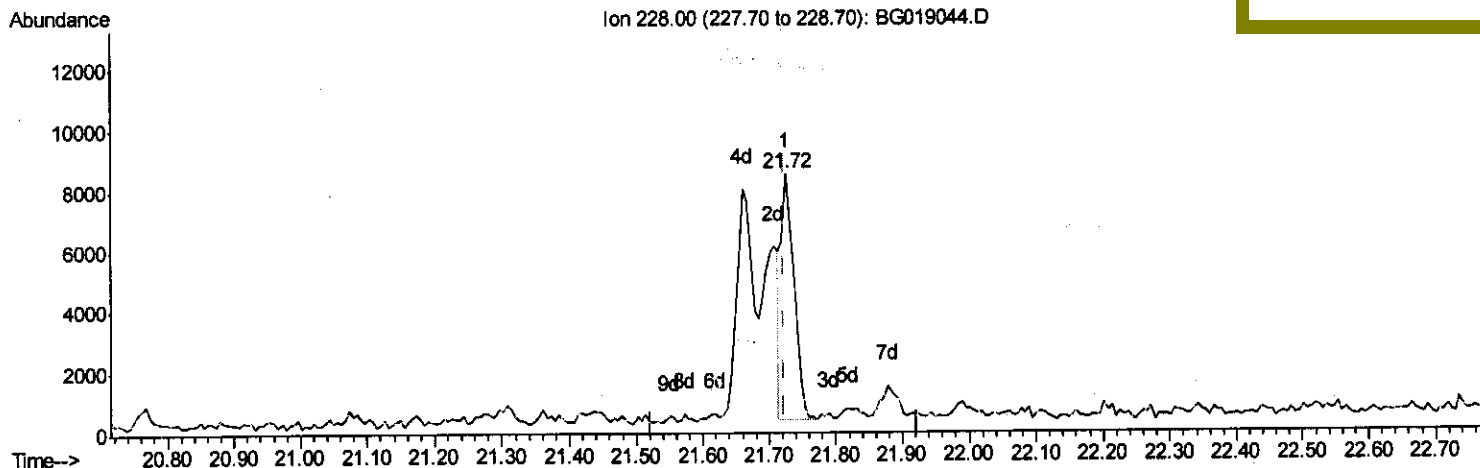
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Instrument :
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 ClientSampleId :
 E5MY7

Quant Time: Oct 07 01:08:33 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG100415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 06 04:27:31 2015
 Response via : Initial Calibration

Manual Integrations
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MMDadoda
 10/6/2015 5:55:12 PM



(85) Chrysene

21.725min (+0.003) 0.61ng/ul

response 10635

Ion	Exp%	Act%
228.00	100	100
226.00	30.20	30.35
229.00	19.50	27.20#
0.00	0.00	0.00

Quantitation Report (Qedit)

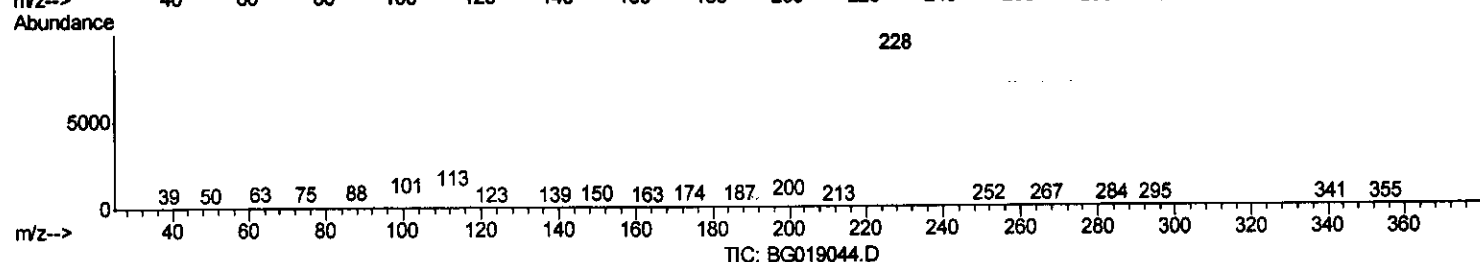
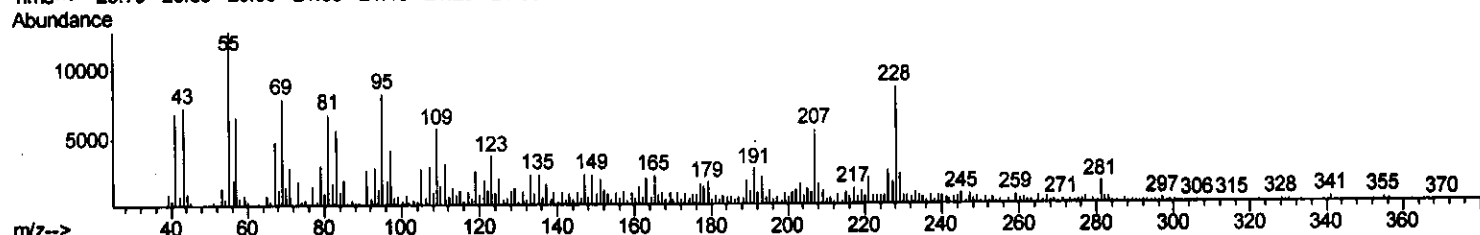
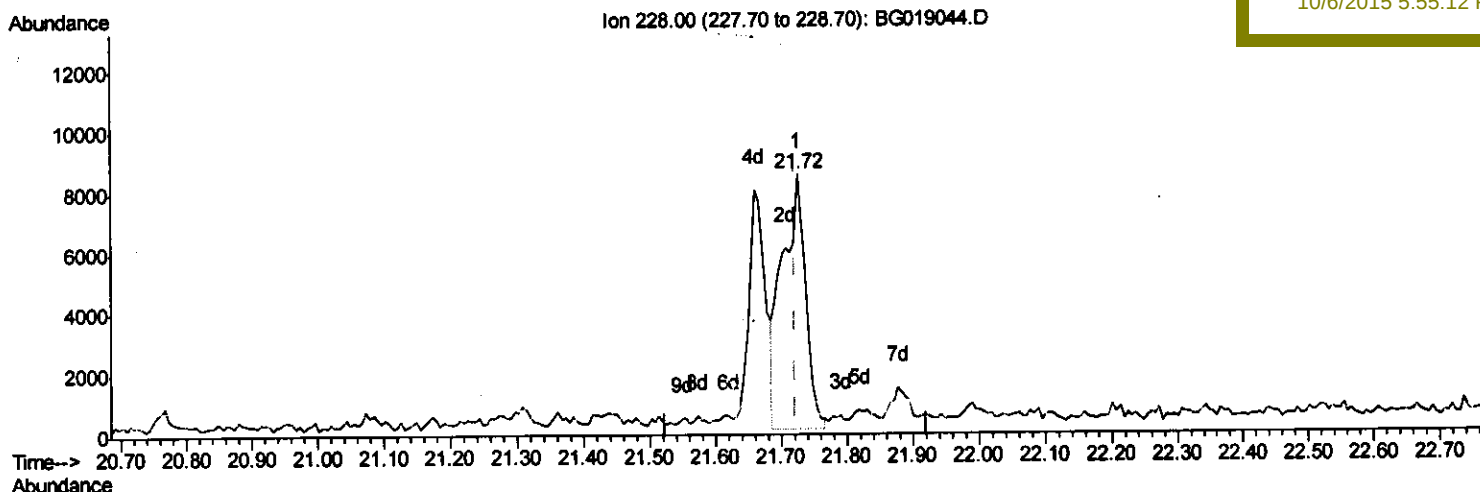
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Manual Integrations
 APPROVED

MMDadoda
 10/6/2015 5:55:12 PM



(85) Chrysene

21.725min (+0.003) 1.19ng/ul m U.M

response 20817

10/7/15

Ion	Exp%	Act%
228.00	100	100
226.00	30.20	30.35
229.00	19.50	27.20#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG100515\
 Data File : BG019044.D
 Acq On : 6 Oct 2015 1:54
 Operator : UM/NP
 Sample : G3865-10
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 Client Sampled :
 E5MY7

Quant Time: Oct 06 09:17:26 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG100415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 06 04:27:31 2015
 Response via : Initial Calibration

Manual Integrations
 APPROVED

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 10/6/2015 5:55:12 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	54153	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	239392	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	144196	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	300229	20.00	ng/ul	0.00
78) Chrysene-d12	21.68	240	321945	20.00	ng/ul	0.00
86) Perylene-d12	24.92	264	296113	20.00	ng/ul	0.03

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	2432	2.03	ng/uL	0.00
5) Phenol-d5	7.21	99	59557	12.76	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.38	67	47321	15.67	ng/ul	0.00
9) 2-Chlorophenol-d4	7.58	132	53887	15.38	ng/ul	0.00
13) 4-Methylphenol-d8	8.75	113	39649	10.81	ng/ul	0.00
19) Nitrobenzene-d5	9.21	128	30532	16.87	ng/ul	0.00
22) 2-Nitrophenol-d4	9.93	143	34622	18.60	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.47	165	61051	16.38	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	34131	7.51	ng/ul	0.00
44) Dimethylphthalate-d6	14.08	166	197617	17.38	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	239545	17.03	ng/ul	0.00
52) 4-Nitrophenol-d4	14.86	143	19249	8.38	ng/ul	0.01
58) Fluorene-d10	15.66	176	172846	16.79	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.77	200	13918	8.55	ng/ul	0.00
71) Anthracene-d10	17.51	188	248401	17.57	ng/ul	0.00
79) Pyrene-d10	19.80	212	266597	17.88	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.70	264	278873	18.41	ng/ul	0.03

Target Compounds

						Qvalue
45) Dimethylphthalate	14.12	163	109371	9.42	ng/ul	98
77) Fluoranthene	19.46	202	20669	1.05	ng/ul	99
80) Pyrene	19.82	202	21514	1.11	ng/ul	98
85) Chrysene	21.72	228	20817m	1.19	ng/ul	

U.M
 10/7/15

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