

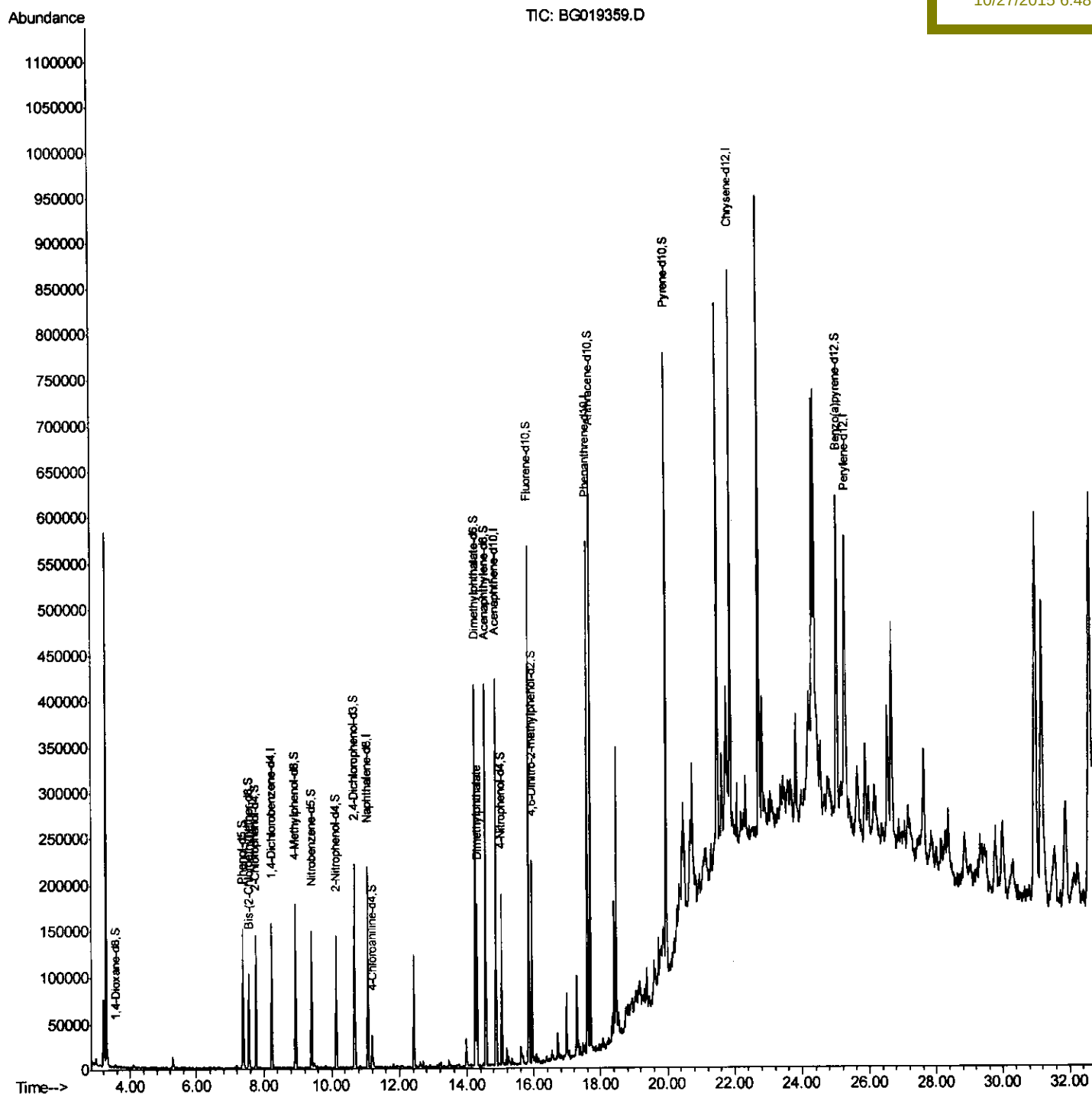
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102615\
Data File : BG019359.D
Acq On : 26 Oct 2015 14:41
Operator : UM/NP
Sample : G4074-14
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JG9H2

Quant Time: Oct 27 03:24:47 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102315.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 27 04:07:50 2015
Response via : Initial Calibration

Manual Integrations
APPROVED

MmDadoda
10/27/2015 6:48:42 PM



Quantitation Report (Qedit)

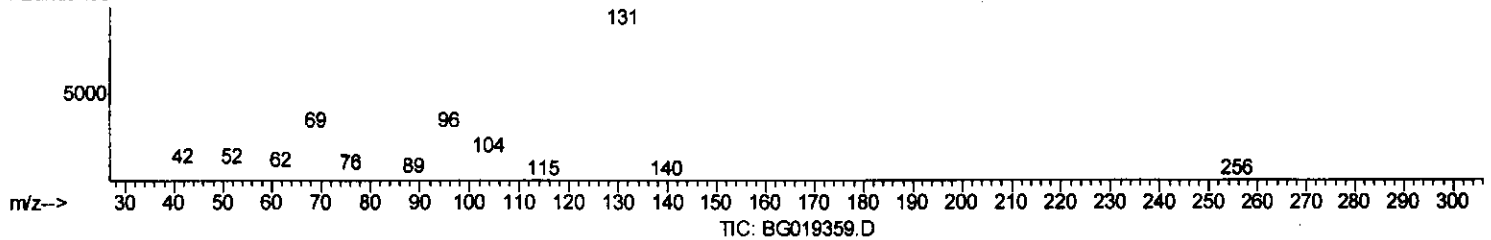
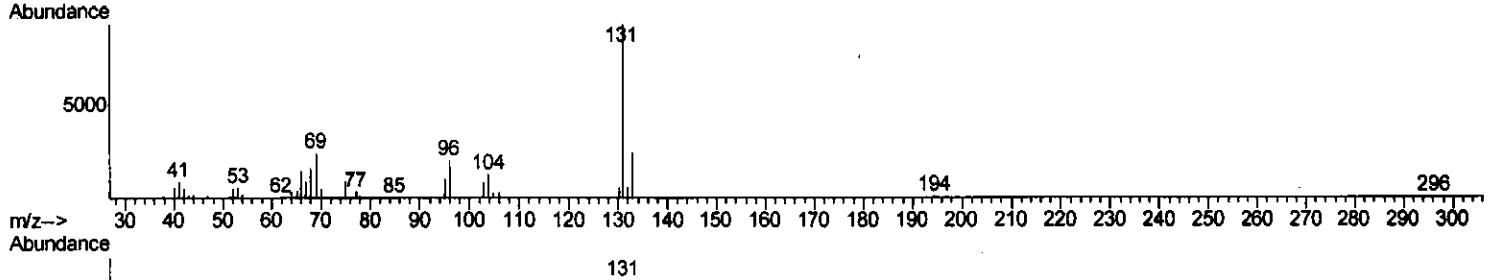
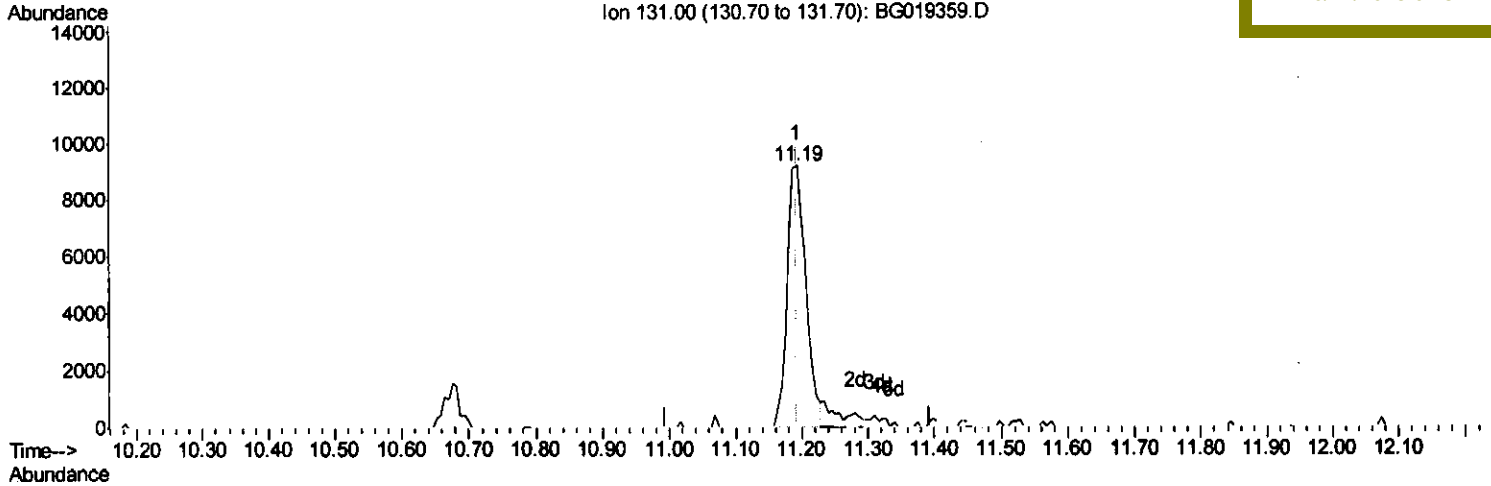
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Quant Time: Oct 27 03:09:21 2015
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(29) 4-Chloroaniline-d4 (S)

11.192min (-0.000) 5.80ng/ul

response 17972

Ion	Exp%	Act%
131.00	100	100
133.00	29.30	27.42
69.00	23.70	26.80
0.00	0.00	0.00

Quantitation Report (Qedit)

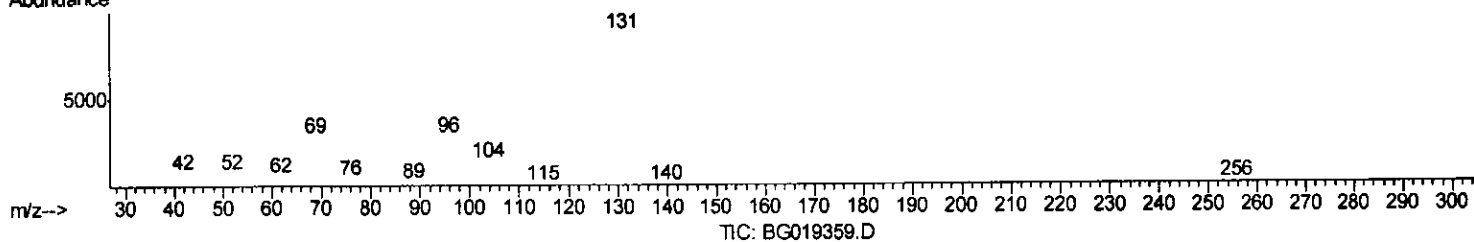
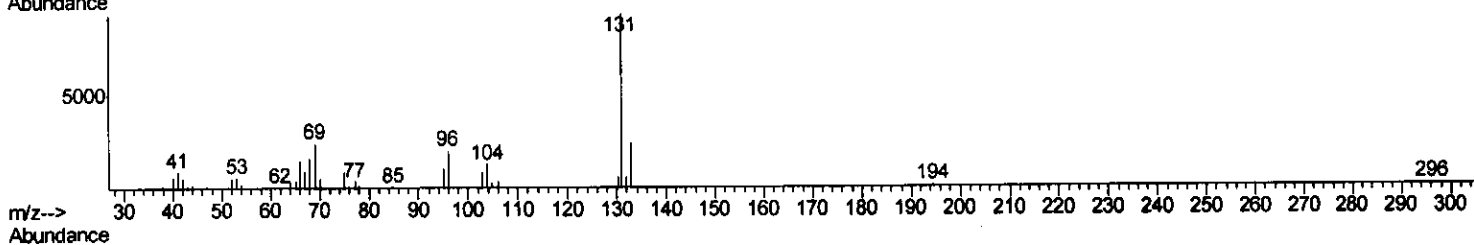
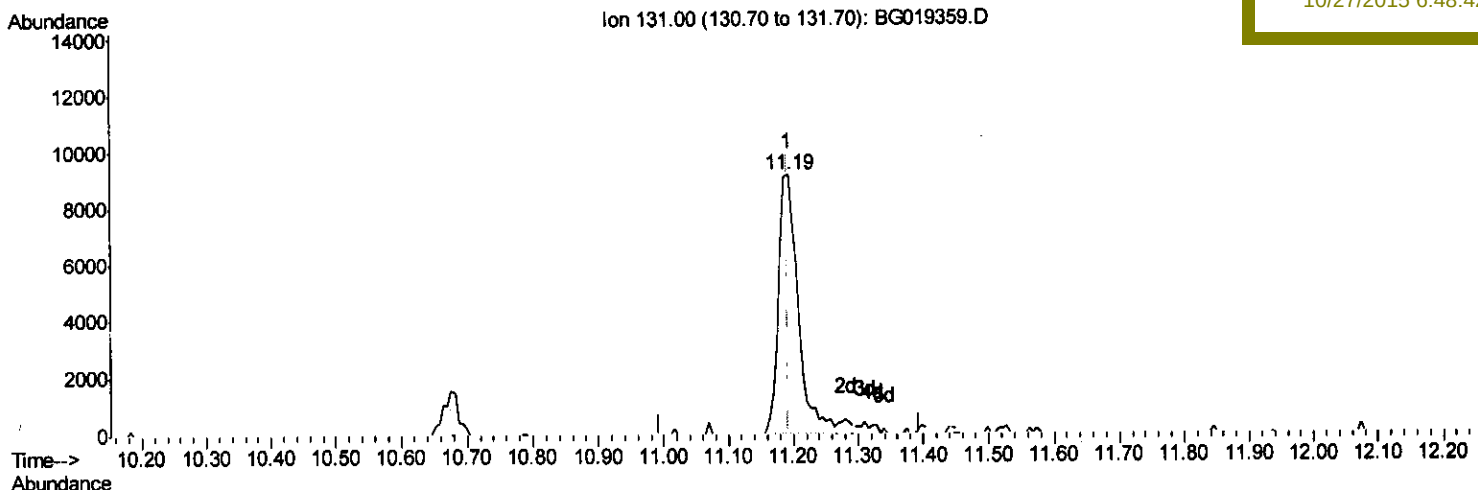
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102615\
 Data File : BG019359.D
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 Operator : UM/NP
 Sample : G4074-14
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JG9H2

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Manual Integrations
 APPROVED

MmDadoda
 10/27/2015 6:48:42 PM



(29) 4-Chloroaniline-d4 (S)

11.192min (-0.000) 6.17ng/ul m

response 19118

Ion	Exp%	Act%
131.00	100	100
133.00	29.30	27.42
69.00	23.70	26.80
0.00	0.00	0.00

U.M
 10/28/15

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102615\
 Data File : BG019359.D
 Acq On : 26 Oct 2015 14:41
 Operator : UM/NP
 Sample : G4074-14
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JG9H2

Quant Time: Oct 27 03:24:47 2015
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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	8.24	152	39128	20.00	ng/ul	0.00
18) Naphthalene-d8	11.07	136	164974	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.86	164	135484	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.60	188	341545	20.00	ng/ul	0.00
78) Chrysene-d12	21.89	240	390984	20.00	ng/ul	0.00
86) Perylene-d12	25.29	264	377363	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.55	96	2851	4.05	ng/uL	0.00
5) Phenol-d5	7.37	99	83860	24.36	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.55	67	51346	24.97	ng/ul	0.00
9) 2-Chlorophenol-d4	7.76	132	53413	24.03	ng/ul	0.00
13) 4-Methylphenol-d8	8.94	113	63783	23.76	ng/ul	0.00
19) Nitrobenzene-d5	9.41	128	27393	21.86	ng/ul	0.00
22) 2-Nitrophenol-d4	10.13	143	35482	24.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.68	165	77028	24.68	ng/ul	0.00
29) 4-Chloroaniline-d4	11.19	131	19118m	6.17	ng/ul	0.00
44) Dimethylphthalate-d6	14.25	166	253721	22.59	ng/ul	0.00
47) Acenaphthylene-d8	14.55	160	291444	23.44	ng/ul	0.00
52) 4-Nitrophenol-d4	15.03	143	33845	18.53	ng/ul	0.00
58) Fluorene-d10	15.85	176	234109	22.21	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.95	200	45952	20.44	ng/ul	0.00
71) Anthracene-d10	17.70	188	360651	23.06	ng/ul	0.00
79) Pyrene-d10	19.96	212	381936	22.30	ng/ul	0.00
90) Benzo(a)pyrene-d12	25.05	264	415771	21.60	ng/ul	0.00

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Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
45) Dimethylphthalate	14.30	163	101990	9.09	ng/ul#	95

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