

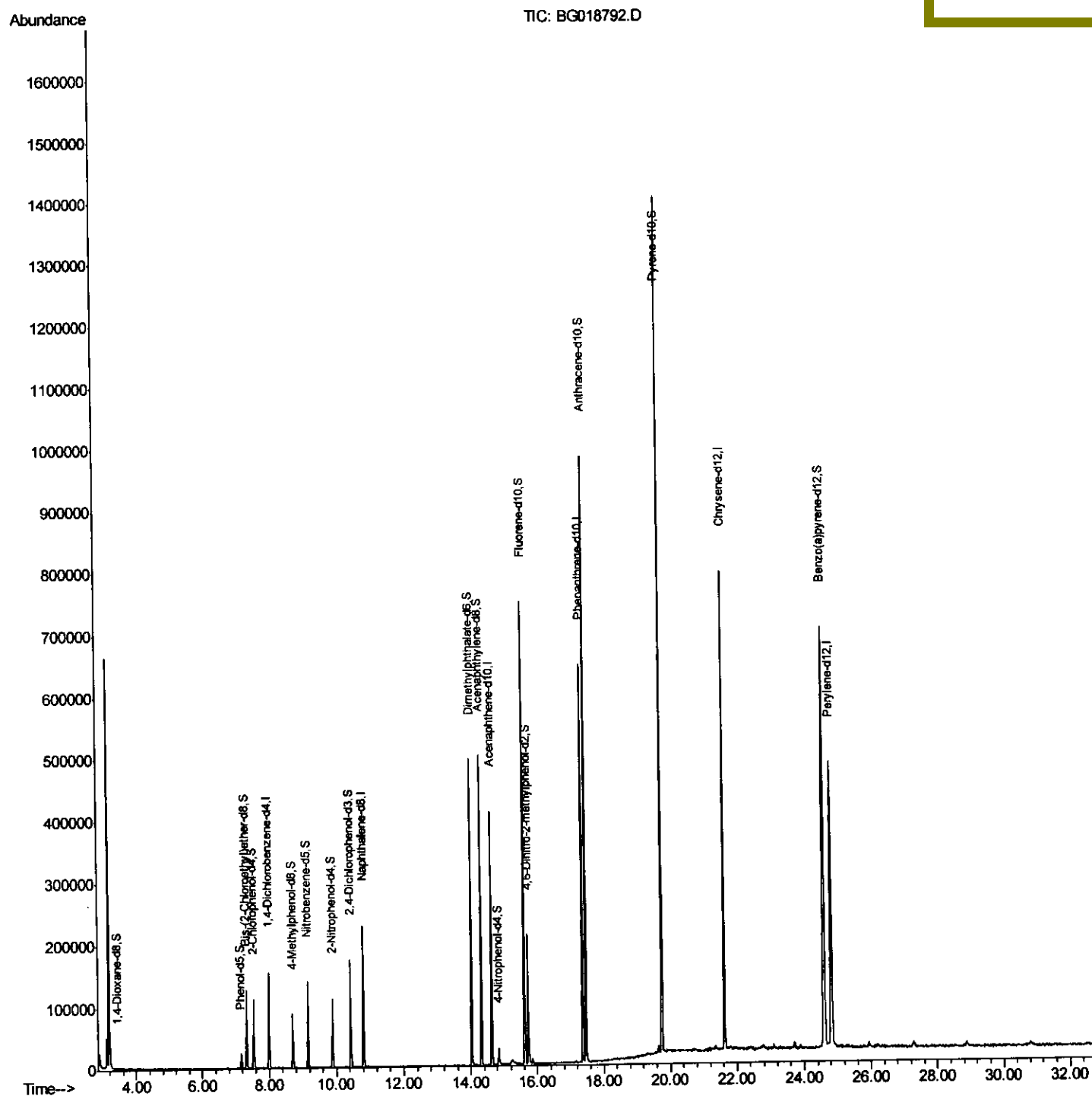
Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
Data File : BG018792.D
Acq On : 21 Sep 2015 21:11
Operator : TP
Sample : G3696-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AN4

Quant Time: Sep 22 02:37:28 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 22 02:06:40 2015
Response via : Initial Calibration

Manual Integrations
APPROVED

MMDadoda
9/22/2015 2:14:30 PM



Quantitation Report (Qedit)

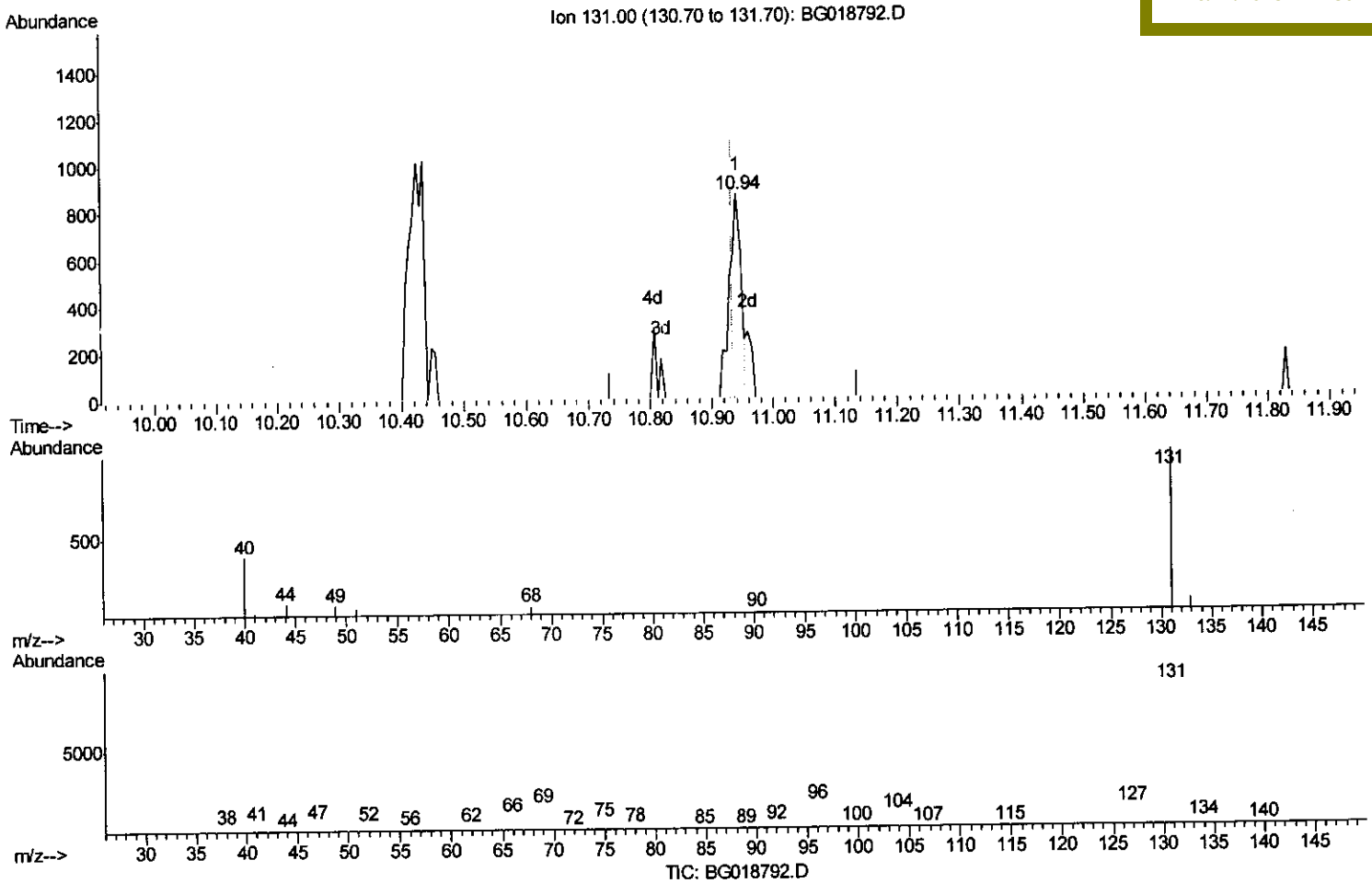
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 Data File : BG018792.D
 Acq On : 21 Sep 2015 21:11
 Operator : TP
 Sample : G3696-13
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Quant Time: Sep 22 02:09:08 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
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Manual Integrations
 APPROVED

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(29) 4-Chloroaniline-d4 (S)

10.941min (+0.006) 0.30ng/ul

response 1151

Ion	Exp%	Act%
131.00	100	100
133.00	32.00	23.12#
69.00	18.90	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

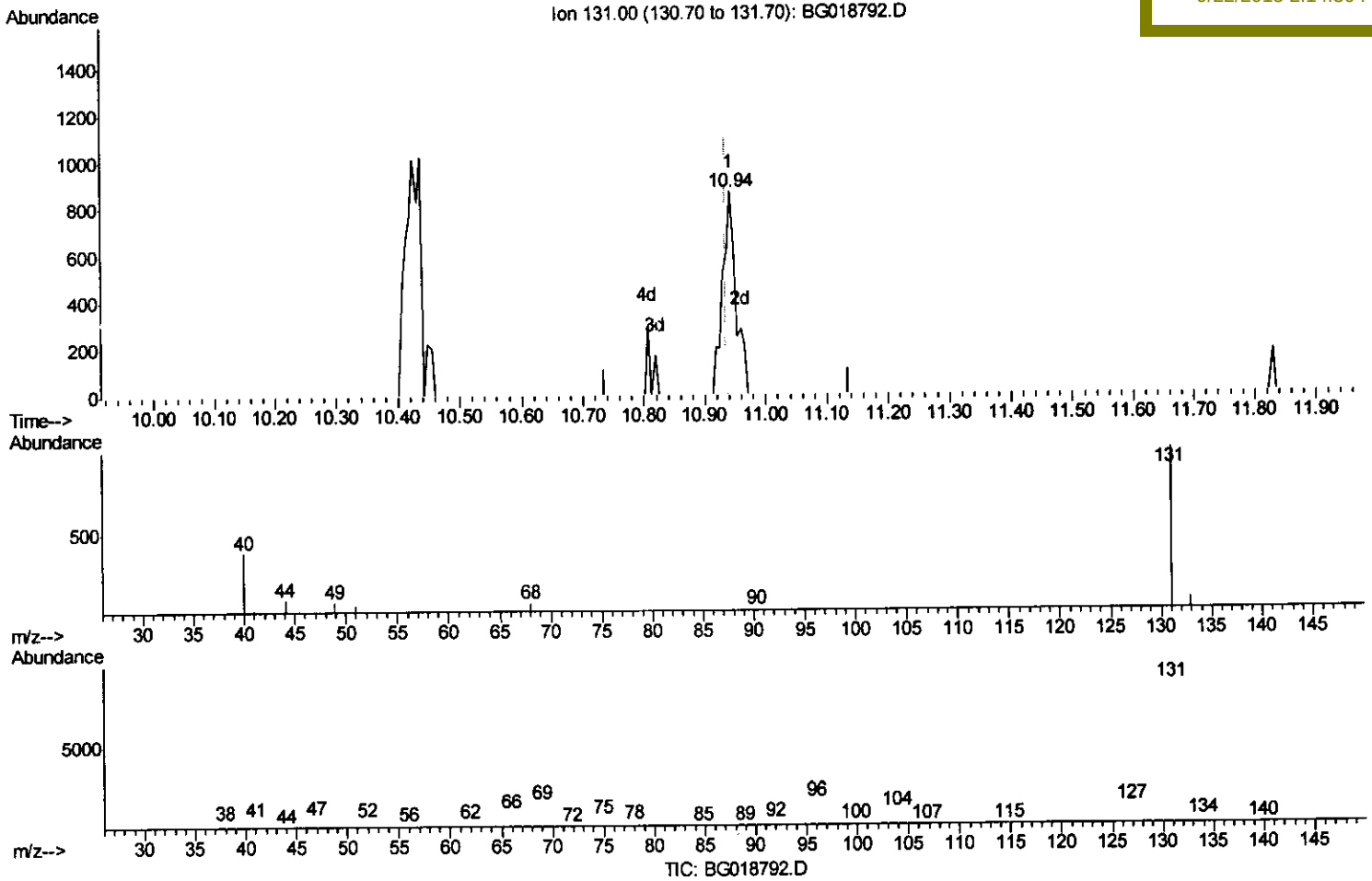
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 Acq On : 21 Sep 2015 21:11
 Operator : TP
 Sample : G3696-13
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 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
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Manual Integrations
 APPROVED

MMDadoda
 9/22/2015 2:14:30 PM



(29) 4-Chloroaniline-d4 (S)

10.941min (+0.006) 0.35ng/ul m

response 1323

Ion	Exp%	Act%
131.00	100	100
133.00	32.00	23.12#
69.00	18.90	0.00#
0.00	0.00	0.00

T.P
 9/23/15

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092115\
 Data File : BG018792.D
 Acq On : 21 Sep 2015 21:11
 Operator : TP
 Sample : G3696-13
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COAN4

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Manual Integrations
 APPROVED

MMDadoda
 9/22/2015 2:14:30 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	45938	20.00	ng/ul	0.00
18) Naphthalene-d8	10.80	136	205104	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.63	164	139485	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.37	188	396968	20.00	ng/ul	0.00
78) Chrysene-d12	21.62	240	456120	20.00	ng/ul	0.00
86) Perylene-d12	24.81	264	459960	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	1256	1.31	ng/uL	-0.01
5) Phenol-d5	7.17	99	18773	4.98	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.32	67	58104	26.58	ng/ul	0.00
9) 2-Chlorophenol-d4	7.53	132	61876	21.66	ng/ul	0.00
13) 4-Methylphenol-d8	8.70	113	38086	12.44	ng/ul	0.00
19) Nitrobenzene-d5	9.16	128	41478	26.69	ng/ul	0.00
22) 2-Nitrophenol-d4	9.88	143	44959	28.77	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.42	165	77074	23.43	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	1323m	0.35	ng/ul	0.00
44) Dimethylphthalate-d6	14.03	166	358411	31.93	ng/ul	0.00
47) Acenaphthylene-d8	14.32	160	390236	29.34	ng/ul	0.00
52) 4-Nitrophenol-d4	14.83	143	10553	4.90	ng/ul	0.00
58) Fluorene-d10	15.62	176	320442	32.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	55435	29.50	ng/ul	0.00
71) Anthracene-d10	17.46	188	604452	34.59	ng/ul	0.00
79) Pyrene-d10	19.75	212	734525	35.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.60	264	729252	32.58	ng/ul	0.00

T.P
 9/23/15

Target Compounds Ovalue

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