

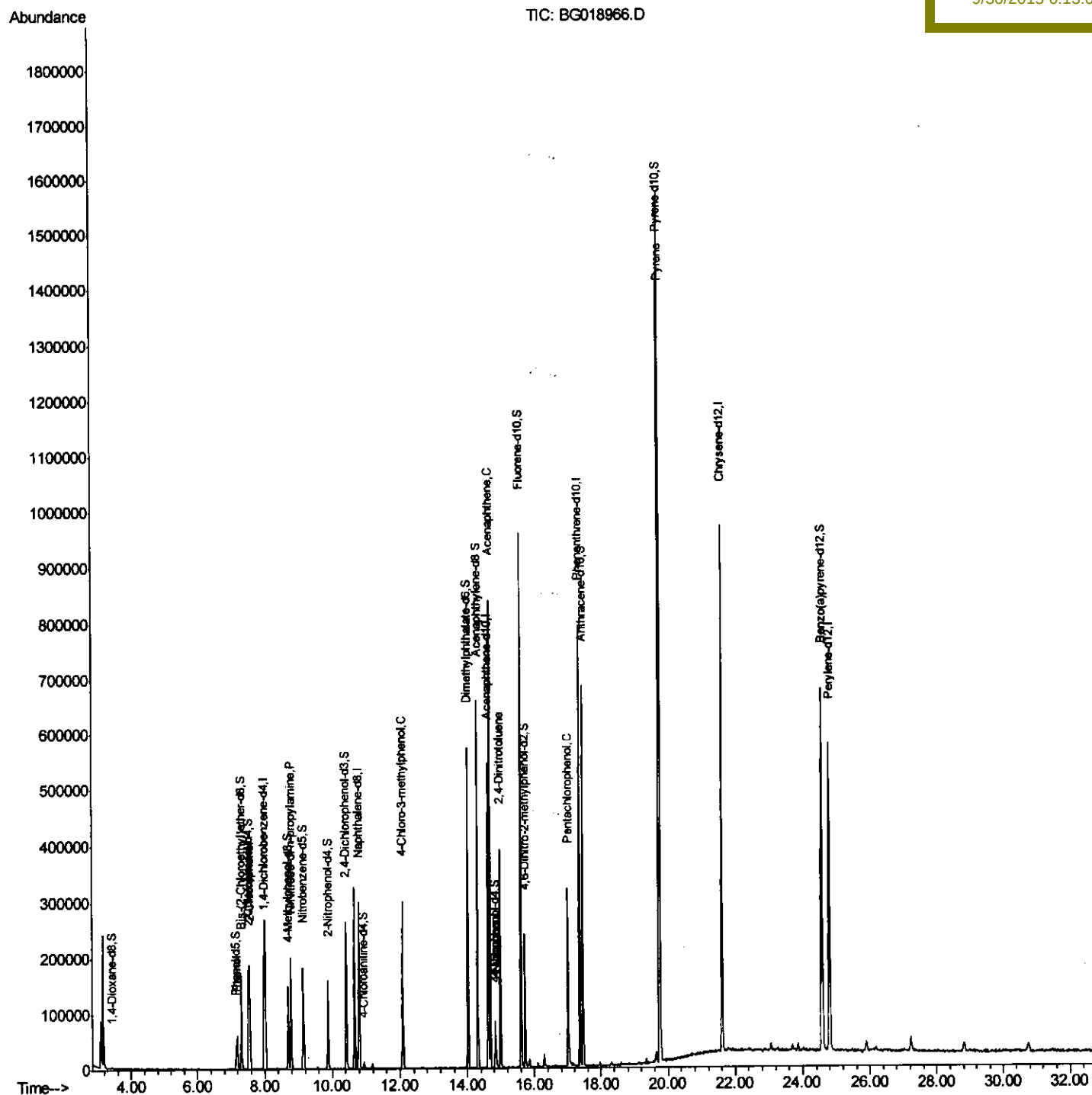
Data Path : Z:\HPCHEM1\BNA_G\Data\BG093015\
Data File : BG018966.D
Acq On : 30 Sep 2015 10:05
Operator : UM/NP
Sample : G3803-14MSD
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E43K2MSD

Quant Time: Sep 30 17:44:21 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Sep 30 04:08:14 2015
Response via : Initial Calibration

Manual Integrations
APPROVED

MMDadoda
9/30/2015 6:13:06 PM



Quantitation Report (Qedit)

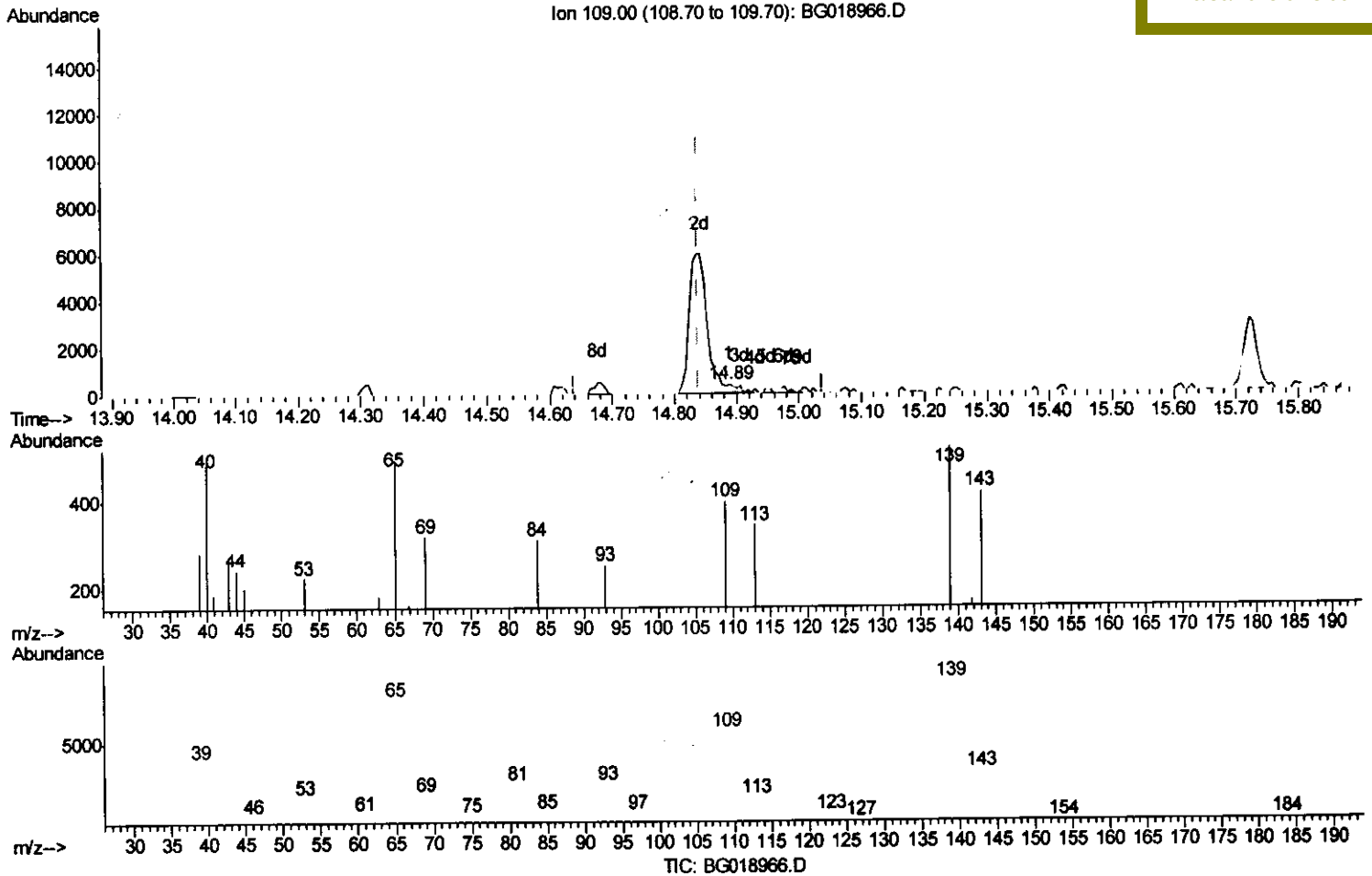
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(53) 4-Nitrophenol

14.888min (+0.050) 0.13ng/ul

response 259

Ion	Exp%	Act%
109.00	100	100
139.00	139.10	131.98
65.00	131.60	125.13
0.00	0.00	0.00

Quantitation Report (Qedit)

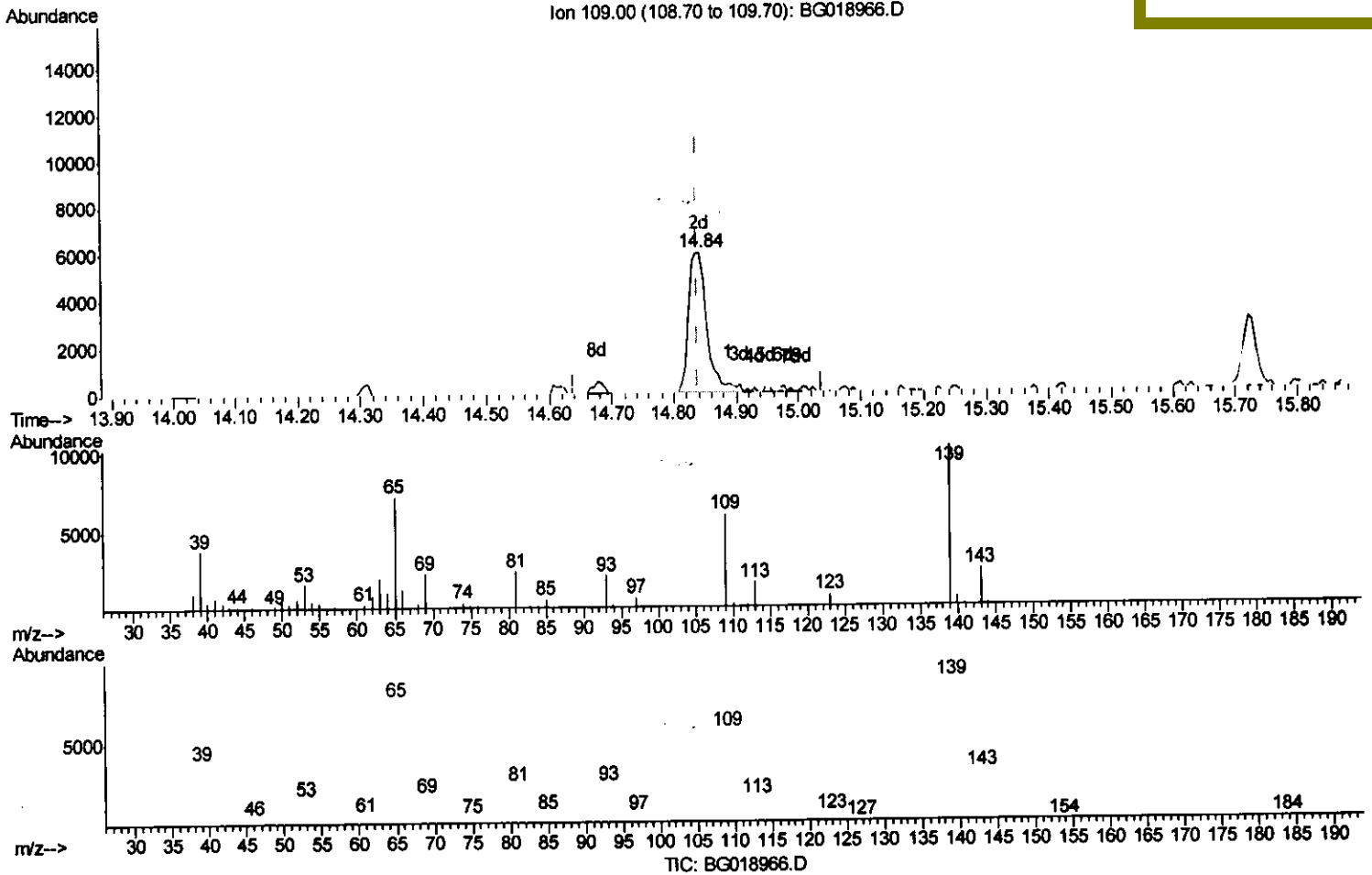
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(53) 4-Nitrophenol

14.841min (+0.003) 6.01ng/ul m

response 12220

Ion	Exp%	Act%
109.00	100	100
139.00	139.10	171.79#
65.00	131.60	120.07
0.00	0.00	0.00

U.M
 10/1/15

Data Path : Z:\HPCHEM1\BNA_G\Data\BG093015\
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 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.98	152	66037	20.00	ng/ul	0.00
18) Naphthalene-d8	10.79	136	281203	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.61	164	192048	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.36	188	518385	20.00	ng/ul	0.00
78) Chrysene-d12	21.62	240	570923	20.00	ng/ul	0.00
86) Perylene-d12	24.79	264	570916	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.42	96	1628	1.19	ng/uL	0.00
5) Phenol-d5	7.16	99	35880	6.62	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.31	67	77465	24.65	ng/ul	0.00
9) 2-Chlorophenol-d4	7.52	132	97352	23.70	ng/ul	0.00
13) 4-Methylphenol-d8	8.70	113	65475	14.88	ng/ul	0.00
19) Nitrobenzene-d5	9.15	128	61269	28.76	ng/ul	0.00
22) 2-Nitrophenol-d4	9.87	143	66636	31.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.41	165	122461	27.15	ng/ul	0.00
29) 4-Chloroaniline-d4	10.93	131	8989	1.71	ng/ul	0.00
44) Dimethylphthalate-d6	14.02	166	442057	28.60	ng/ul	0.00
47) Acenaphthylene-d8	14.31	160	527713	28.81	ng/ul	0.00
52) 4-Nitrophenol-d4	14.82	143	17886	6.03	ng/ul	0.00
58) Fluorene-d10	15.60	176	404646	29.94	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.72	200	68958	28.10	ng/ul	0.00
71) Anthracene-d10	17.46	188	406251	17.80	ng/ul	0.00
79) Pyrene-d10	19.74	212	790105	30.08	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.58	264	695336	25.03	ng/ul	0.00

Target Compounds

					Qvalue
6) Phenol	7.18	94	38910	6.94 ng/ul#	86
10) 2-Chlorophenol	7.55	128	95757	22.52 ng/ul	90
15) N-Nitroso-di-n-propylamine	8.78	70	81847	23.68 ng/ul#	82
33) 4-Chloro-3-methylphenol	12.07	107	122260	25.40 ng/ul	95
50) Acenaphthene	14.68	153	362675	27.63 ng/ul	99
53) 4-Nitrophenol	14.84	109	12220m	6.01 ng/ul	
55) 2,4-Dinitrotoluene	14.98	165	137978	30.76 ng/ul#	82
69) Pentachlorophenol	17.01	266	71614	21.81 ng/ul	92
80) Pyrene	19.77	202	943231	28.10 ng/ul	98

UM
 10/1/15

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