

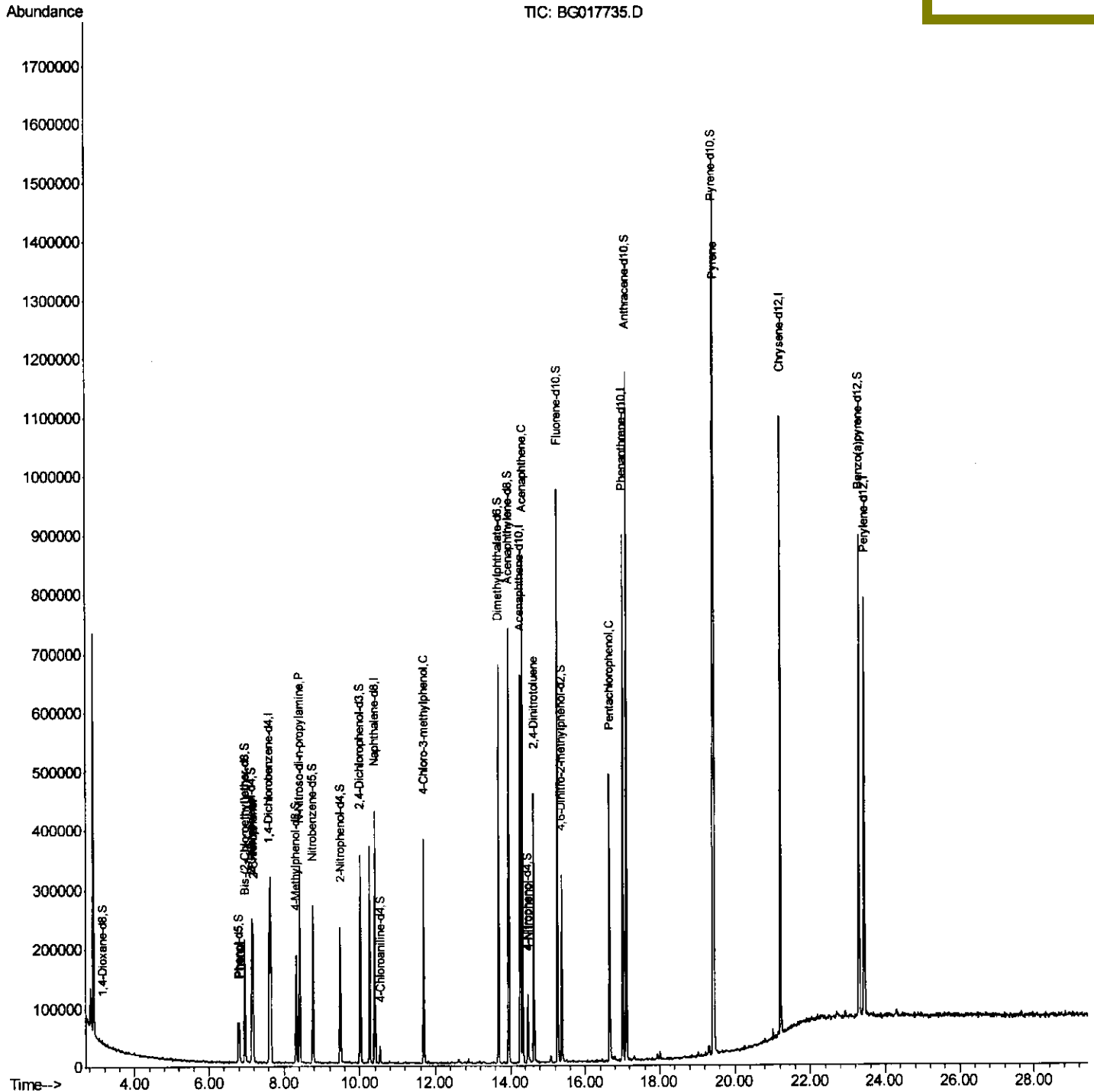
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061915\
Data File : BG017735.D
Acq On : 19 Jun 2015 18:32
Operator : TP/IZ
Sample : G2700-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
HOUN2MS

Quant Time: Jun 20 01:28:47 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG061915.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 19 17:30:37 2015
Response via : Initial Calibration

Manual Integrations
APPROVED

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Quantitation Report (Qedit)

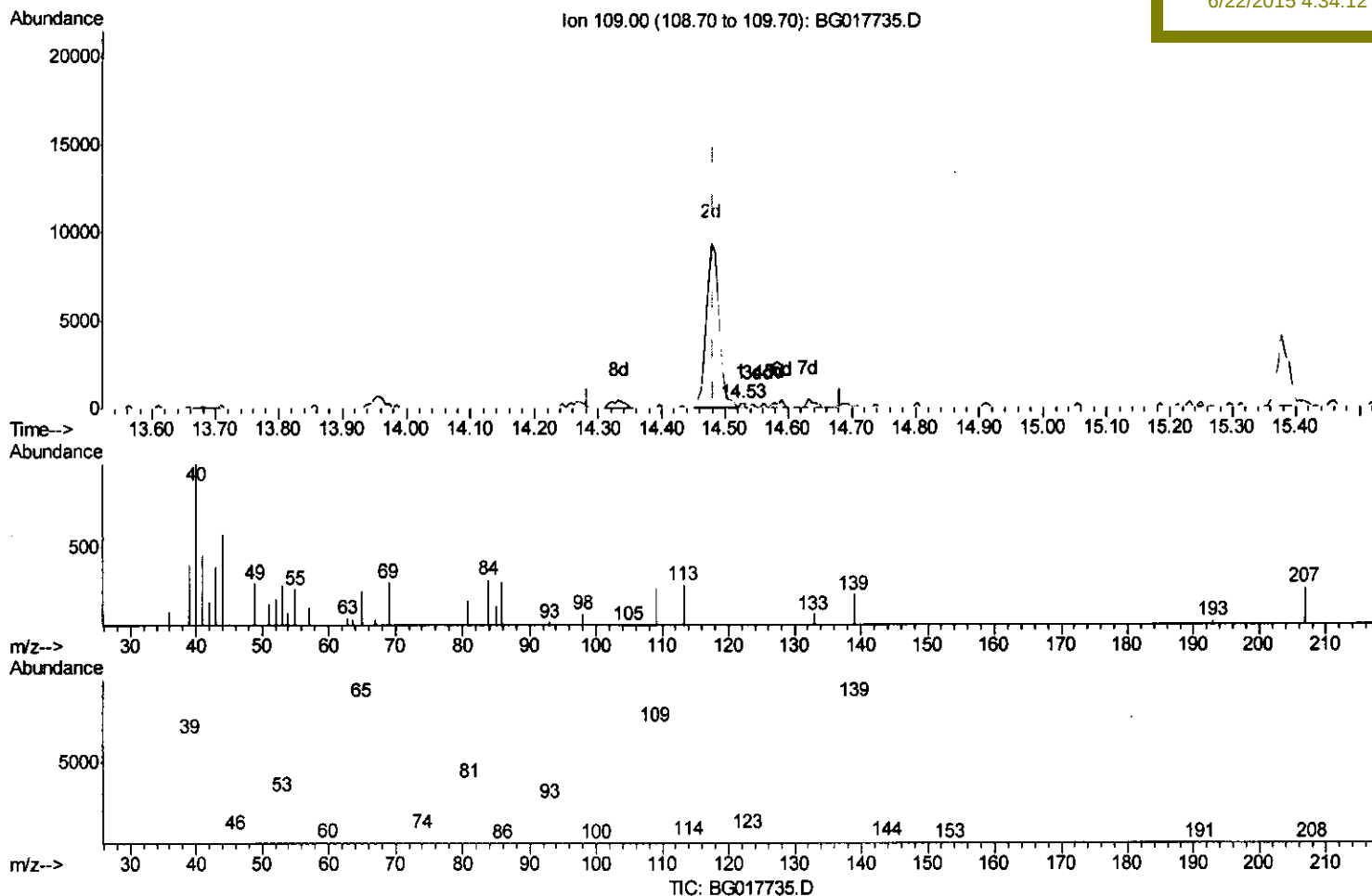
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ClientSampleId :
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Quant Time: Jun 20 01:14:16 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG061915.M
 Quant Title : SVOA CALIBRATION
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(52) 4-Nitrophenol

14.526min (+0.043) 0.07ng/ul

response 177

Ion	Exp%	Act%
109.00	100	100
139.00	102.10	92.60
65.00	117.90	97.11
0.00	0.00	0.00

Quantitation Report (Qedit)

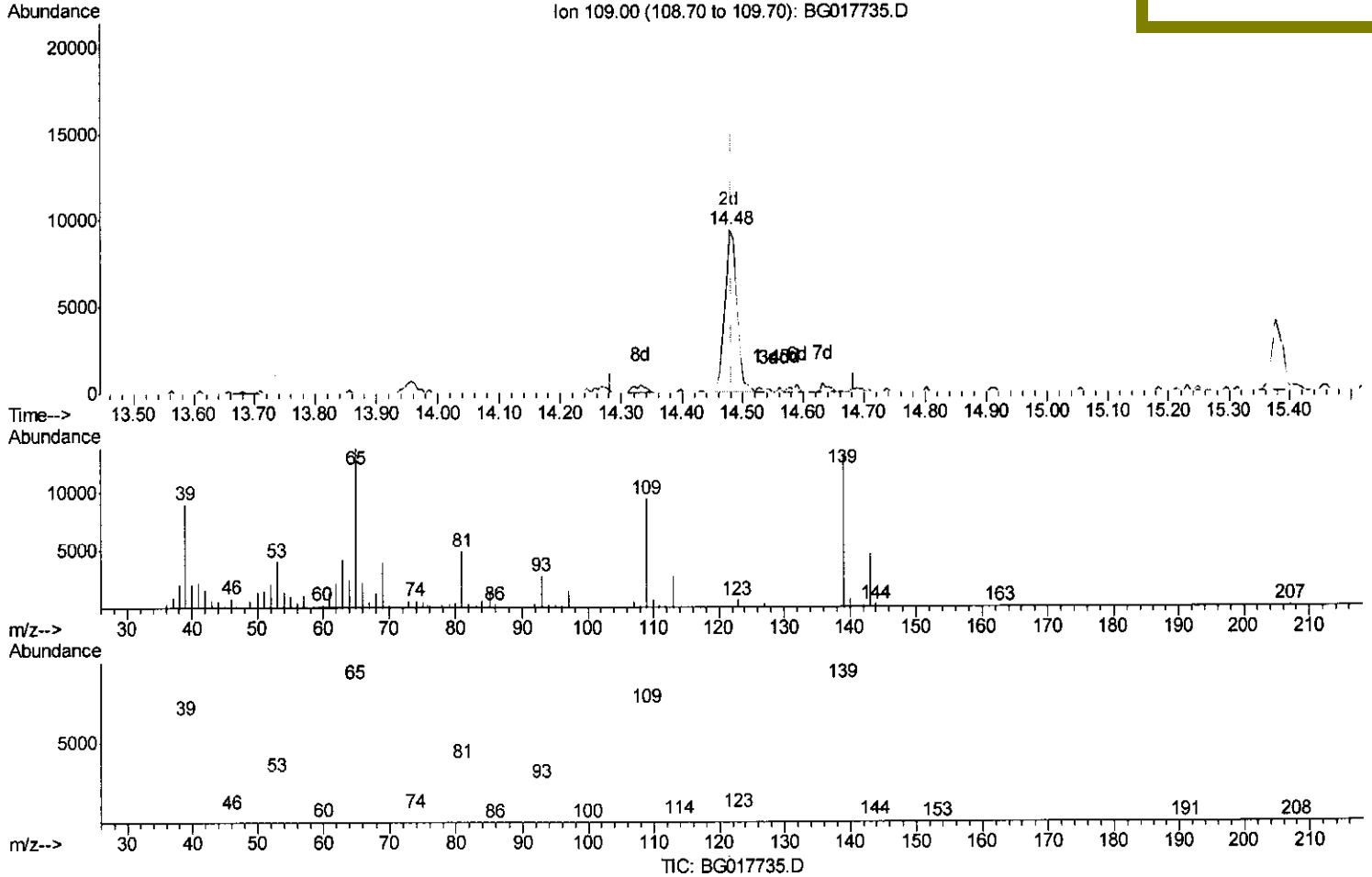
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(52) 4-Nitrophenol

14.479min (-0.004) 5.57ng/ul m

response 13374

Ion	Exp%	Act%
109.00	100	100
139.00	102.10	140.35#
65.00	117.90	147.17#
0.00	0.00	0.00

T.P.
 6/22/15

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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	7.62	152	71732	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	336854	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.27	164	210949	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.02	188	454829	20.00	ng/ul	0.00
77) Chrysene-d12	21.22	240	453314	20.00	ng/ul	0.00
85) Perylene-d12	23.44	264	442495	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.14	96	1838	1.14	ng/uL	0.00
5) Phenol-d5	6.79	99	31486	5.07	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	88419	24.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	100482	19.68	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	63290	12.34	ng/ul	0.00
19) Nitrobenzene-d5	8.76	128	63312	24.43	ng/ul	0.00
22) 2-Nitrophenol-d4	9.48	143	68334	25.72	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.01	165	109375	22.48	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	14245	2.15	ng/ul	0.00
43) Dimethylphthalate-d6	13.69	166	384142	26.77	ng/ul	0.00
46) Acenaphthylene-d8	13.96	160	484482	25.69	ng/ul	0.00
51) 4-Nitrophenol-d4	14.47	143	15887	5.15	ng/ul	0.00
57) Fluorene-d10	15.27	176	371970	26.44	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.38	200	59045	24.45	ng/ul	0.00
70) Anthracene-d10	17.12	188	571555	27.24	ng/ul	0.00
78) Pyrene-d10	19.42	212	652801	29.26	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.30	264	492552	25.04	ng/ul	0.00

Target Compounds

						Qvalue
6) Phenol	6.81	94	33244	5.02	ng/ul	96
10) 2-Chlorophenol	7.18	128	91939	17.98	ng/ul	94
15) N-Nitroso-di-n-propylamine	8.42	70	113061	24.88	ng/ul#	90
33) 4-Chloro-3-methylphenol	11.69	107	119672	21.00	ng/ul	88
49) Acenaphthene	14.33	153	323192	23.55	ng/ul	98
52) 4-Nitrophenol	14.48	109	13374m	5.57	ng/ul	
54) 2,4-Dinitrotoluene	14.64	165	108334	25.09	ng/ul#	89
68) Pentachlorophenol	16.66	266	71904	26.78	ng/ul	97
79) Pyrene	19.45	202	771658	26.91	ng/ul#	82

T.P.
 5/22/15

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