

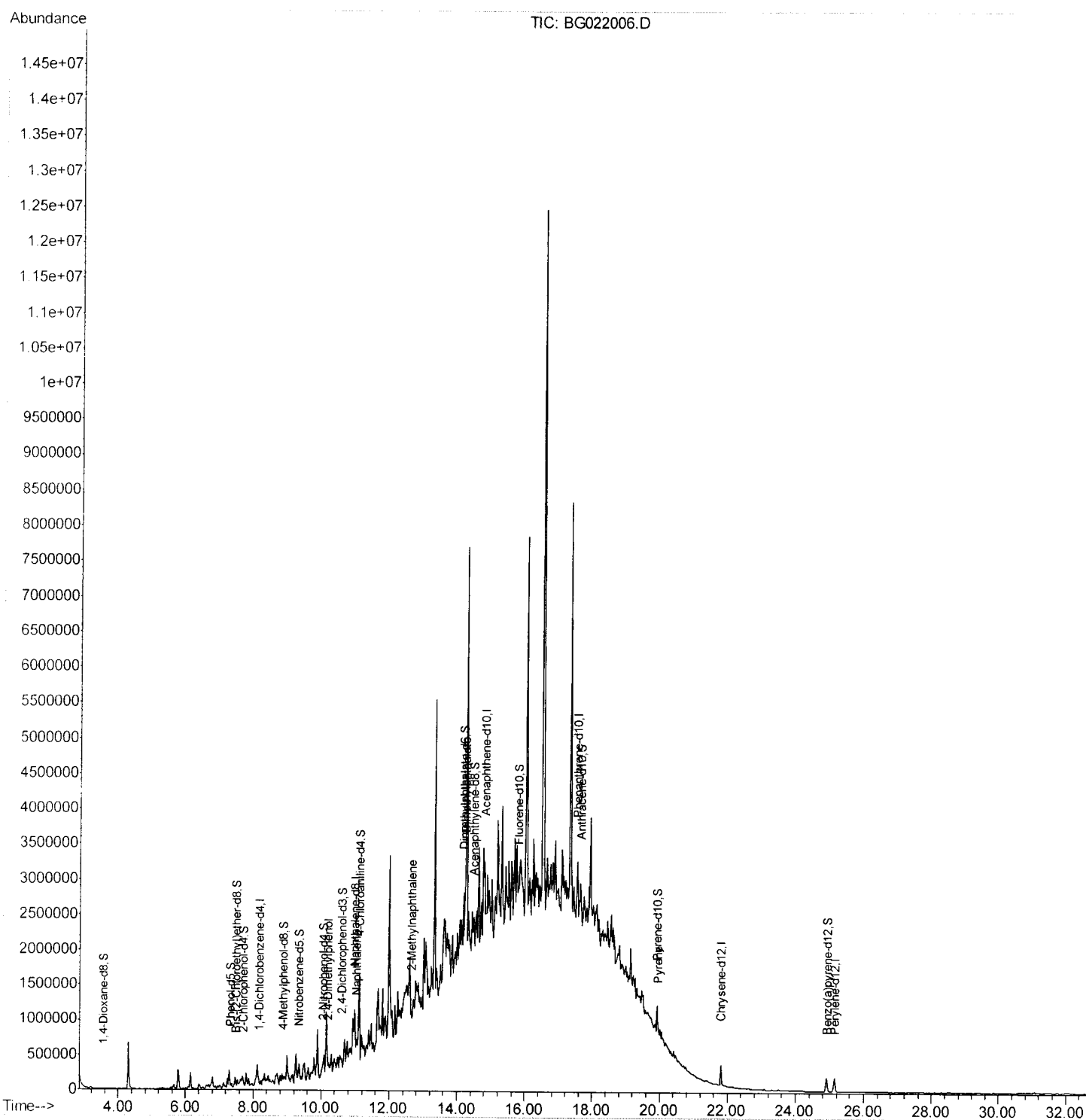
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052116\
Data File : BG022006.D
Acq On : 22 May 2016 10:41
Operator : UM/SJ
Sample : H3124-19
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JHGG5

Manual Integrations
APPROVED

umangi
5/23/2016 8:41:17 PM

Quant Time: May 23 18:42:12 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG052116.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon May 23 17:16:55 2016
Response via : Initial Calibration



Quantitation Report (Qedit)

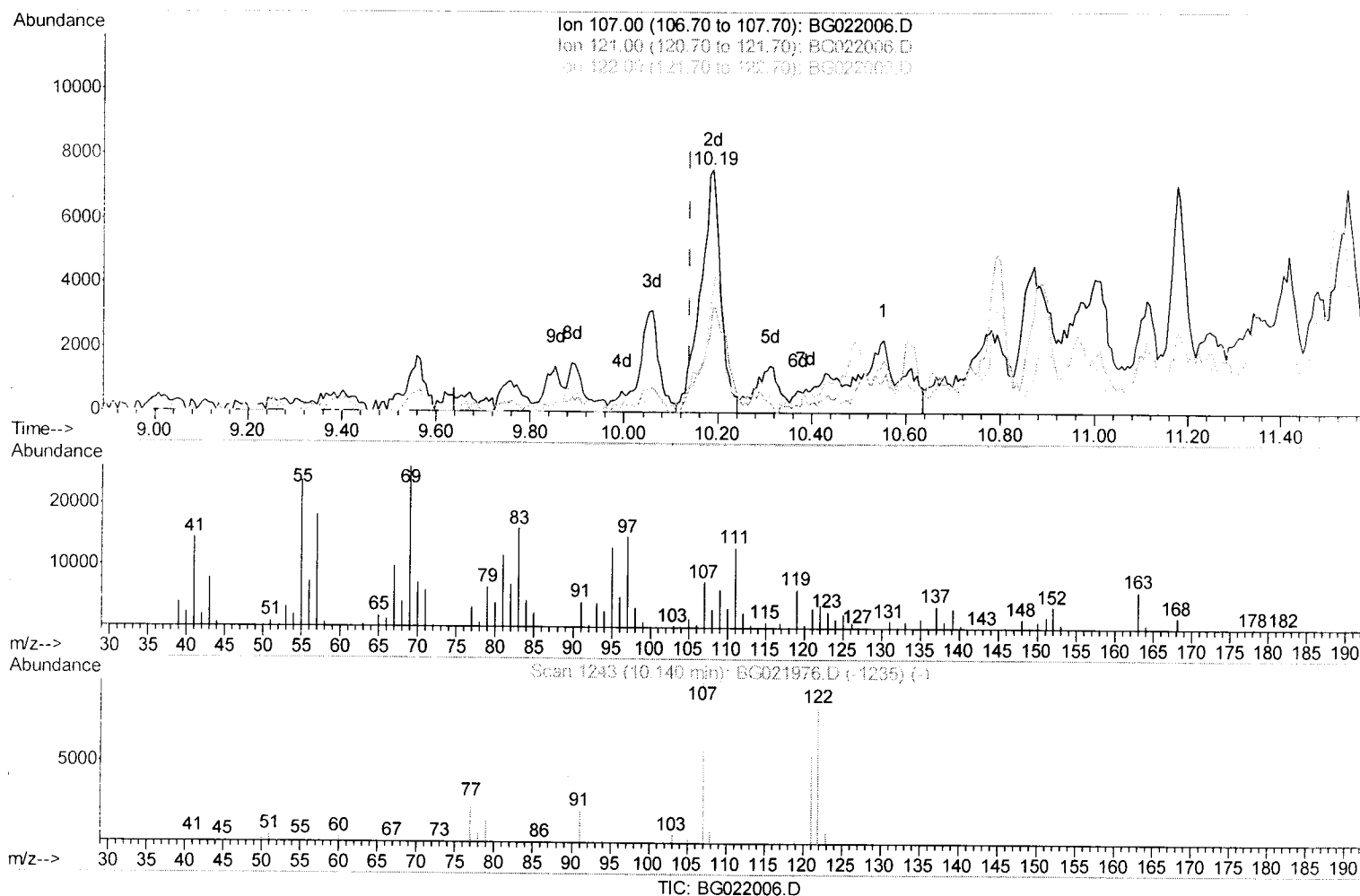
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052116\
 Data File : BG022006.D
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 Operator : UM/SJ
 Sample : H3124-19
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Instrument :
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Manual Integrations
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Quant Time: May 23 17:36:58 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG052116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon May 23 17:16:55 2016
 Response via : Initial Calibration



(24) 2,4-Dimethylphenol

10.189min (+0.048) 6.60ng/ul m

response 22762

Ion	Exp%	Act%
107.00	100	100
121.00	54.60	43.28#
122.00	87.90	51.26#
0.00	0.00	0.00

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 05/24/16

Quantitation Report (Qedit)

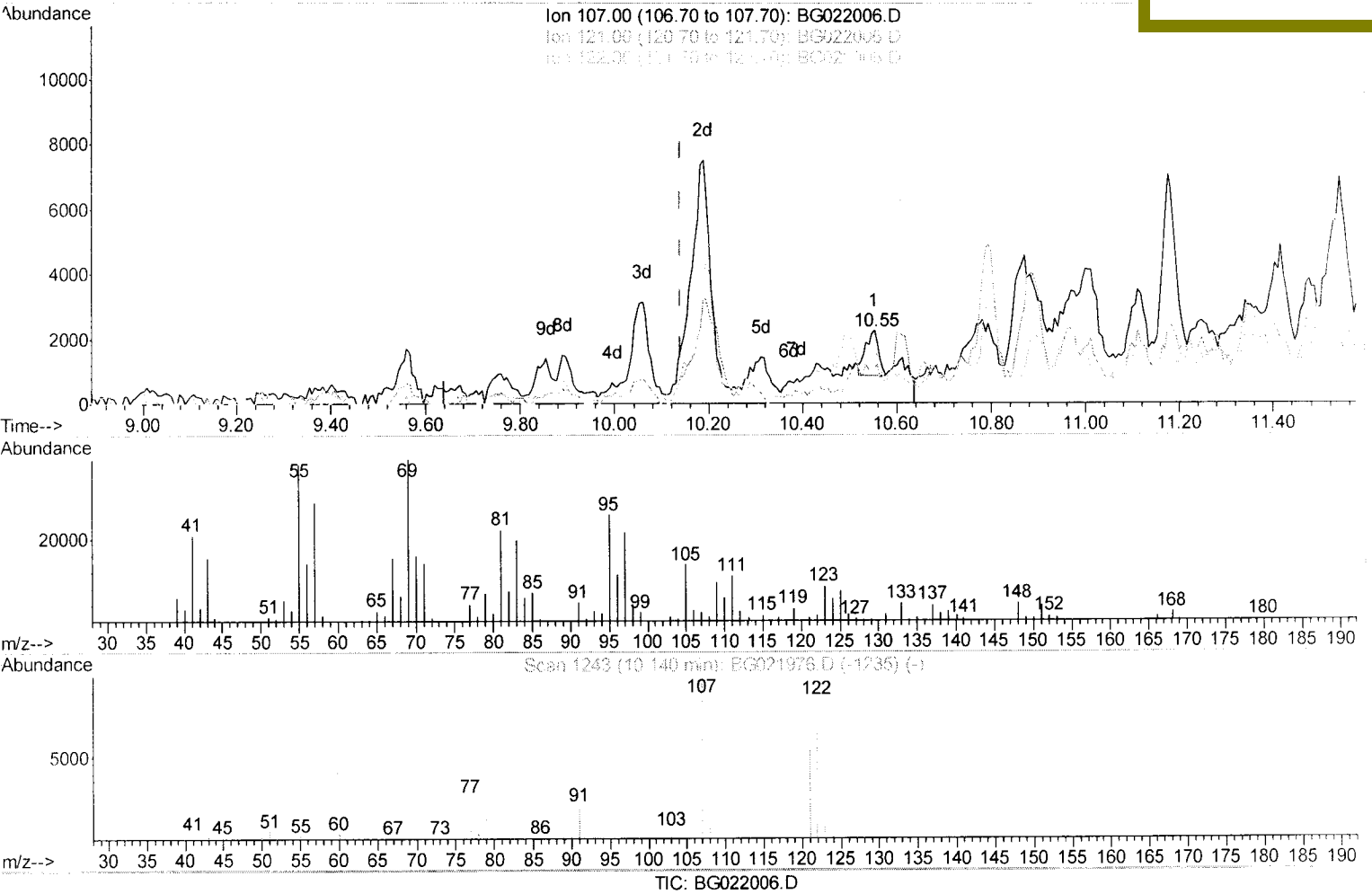
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052116\
 Data File : BG022006.D
 Acq On : 22 May 2016 10:41
 Operator : UM/SJ
 Sample : H3124-19
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
ClientSampled :
 JHGG5

Manual Integrations
APPROVED

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 5/23/2016 8:41:17 PM

Quant Time: May 23 17:36:58 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG052116.M
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 Response via : Initial Calibration



(24) 2,4-Dimethylphenol

10.553min (+0.413) 0.76ng/ul

response 2630

Ion	Exp%	Act%
107.00	100	100
121.00	54.60	46.05
122.00	87.90	73.00
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052116\
 Data File : BG022006.D
 Acq On : 22 May 2016 10:41
 Operator : UM/SJ
 Sample : H3124-19
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JHGG5

Quant Time: May 23 18:42:12 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG052116.M
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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	8.16	152	46334	20.00	ng/ul	0.00
18) Naphthalene-d8	10.98	136	213883	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.80	164	114116	20.00	ng/ul	0.02
61) Phenanthrene-d10	17.55	188	202381	20.00	ng/ul	0.02
75) Chrysene-d12	21.81	240	202928	20.00	ng/ul	0.00
83) Perylene-d12	25.14	264	204491	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.54	96	3301	3.71	ng/uL	0.00
5) Phenol-d5	7.31	99	75418	18.01	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.48	67	43750	19.93	ng/ul	0.00
9) 2-Chlorophenol-d4	7.69	132	70382	20.11	ng/ul	0.00
13) 4-Methylphenol-d8	8.86	113	60445	17.02	ng/ul	0.00
19) Nitrobenzene-d5	9.33	128	33523	21.25	ng/ul	0.00
22) 2-Nitrophenol-d4	10.05	143	35411	20.72	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.60	165	62355	20.69	ng/ul	0.00
29) 4-Chloroaniline-d4	11.11	131	90606	27.60	ng/ul	0.00
43) Dimethylphthalate-d6	14.19	166	177950	20.76	ng/ul	0.00
46) Acenaphthylene-d8	14.50	160	258847	21.63	ng/ul	0.01
51) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
57) Fluorene-d10	15.79	176	158609	19.76	ng/ul	0.02
62) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
70) Anthracene-d10	17.65	188	220613	22.71	ng/ul	0.02
76) Pyrene-d10	19.92	212	227675	19.90	ng/ul	0.01
87) Benzo(a)pyrene-d12	24.91	264	213489	22.50	ng/ul	0.00

Target Compounds

					Qvalue	
24) 2,4-Dimethylphenol	10.19	107	22762m	6.60	ng/ul	
28) Naphthalene	11.03	128	46836	4.32	ng/ul#	81
34) 2-Methylnaphthalene	12.63	142	51614	6.66	ng/ul	96
44) Dimethylphthalate	14.24	163	127949	14.86	ng/ul#	66
77) Pyrene	19.95	202	20869	1.46	ng/ul#	83

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 05/24/16

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