

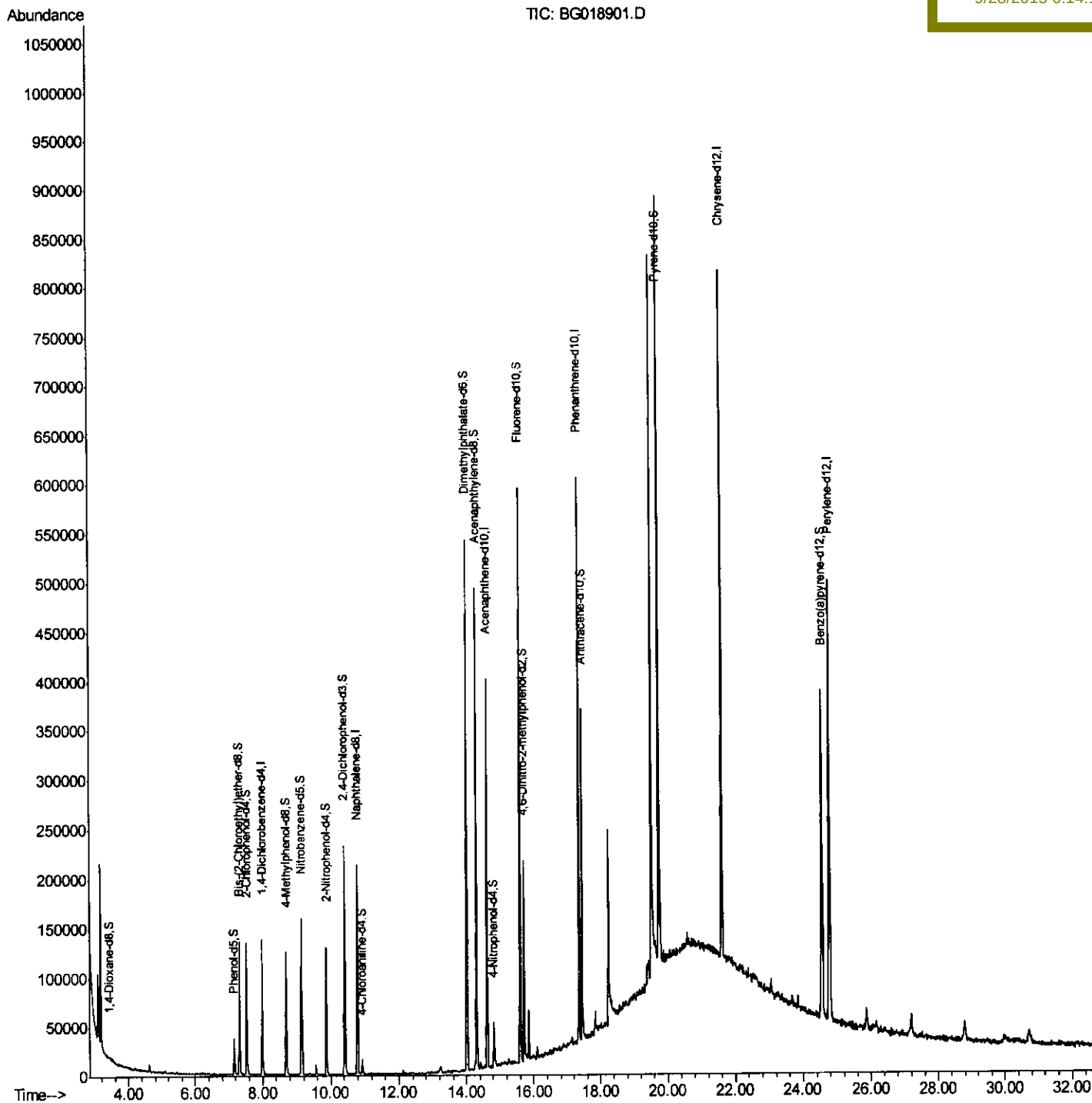
Data Path : Z:\HPCHEM1\BNA_G\Data\BG092515\
Data File : BG018901.D
Acq On : 26 Sep 2015 5:40
Operator : UM/NP
Sample : G3800-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C03F9

Quant Time: Sep 26 06:58:09 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 26 04:18:39 2015
Response via : Initial Calibration

Manual Integrations
APPROVED

MMDadoda
9/28/2015 6:14:15 PM



Quantitation Report (Qedit)

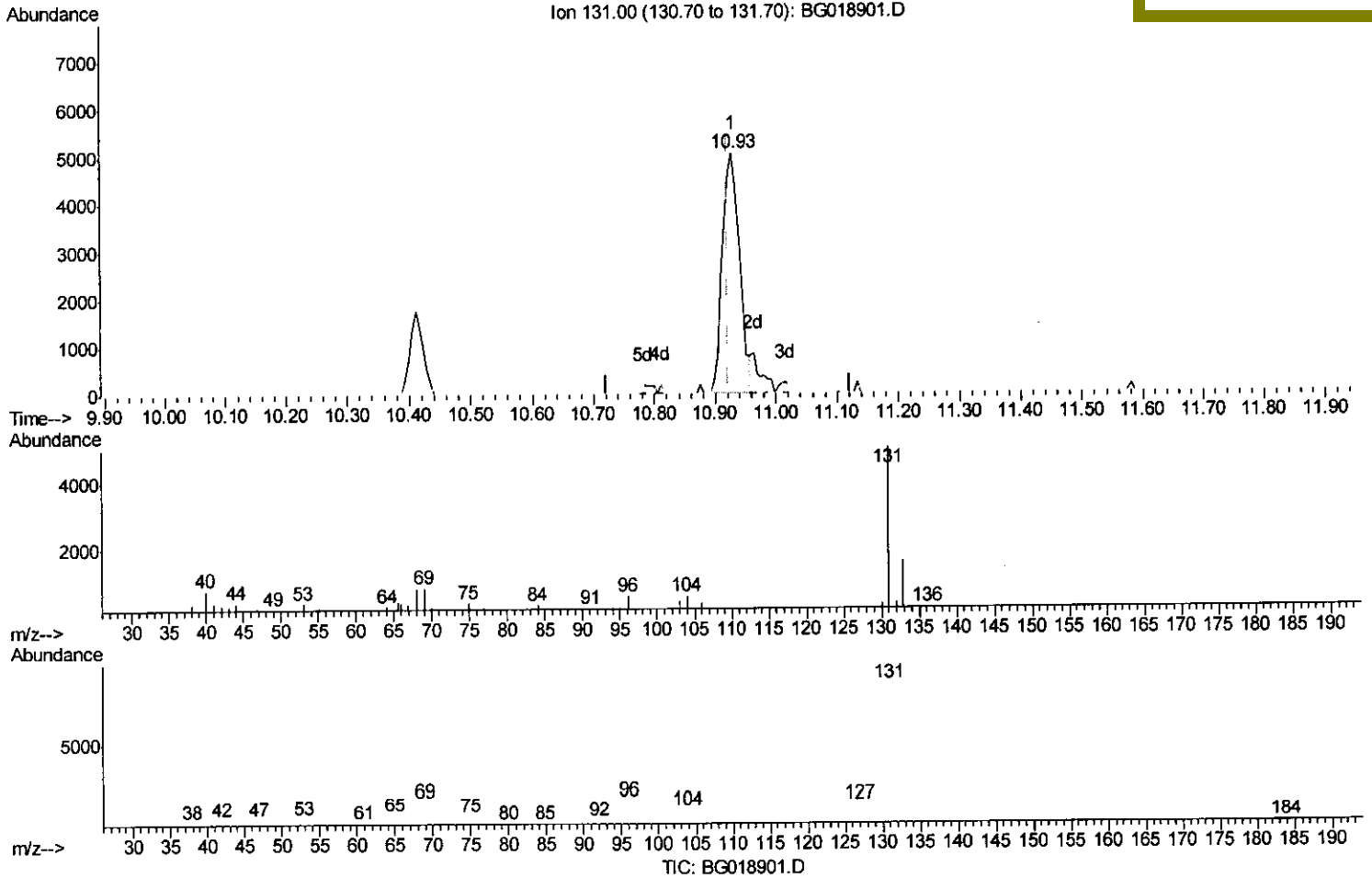
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(29) 4-Chloroaniline-d4 (S)

10.928min (+0.006) 2.69ng/ul

response 9915

Ion	Exp%	Act%
131.00	100	100
133.00	32.00	31.77
69.00	18.90	16.02
0.00	0.00	0.00

Quantitation Report (Qedit)

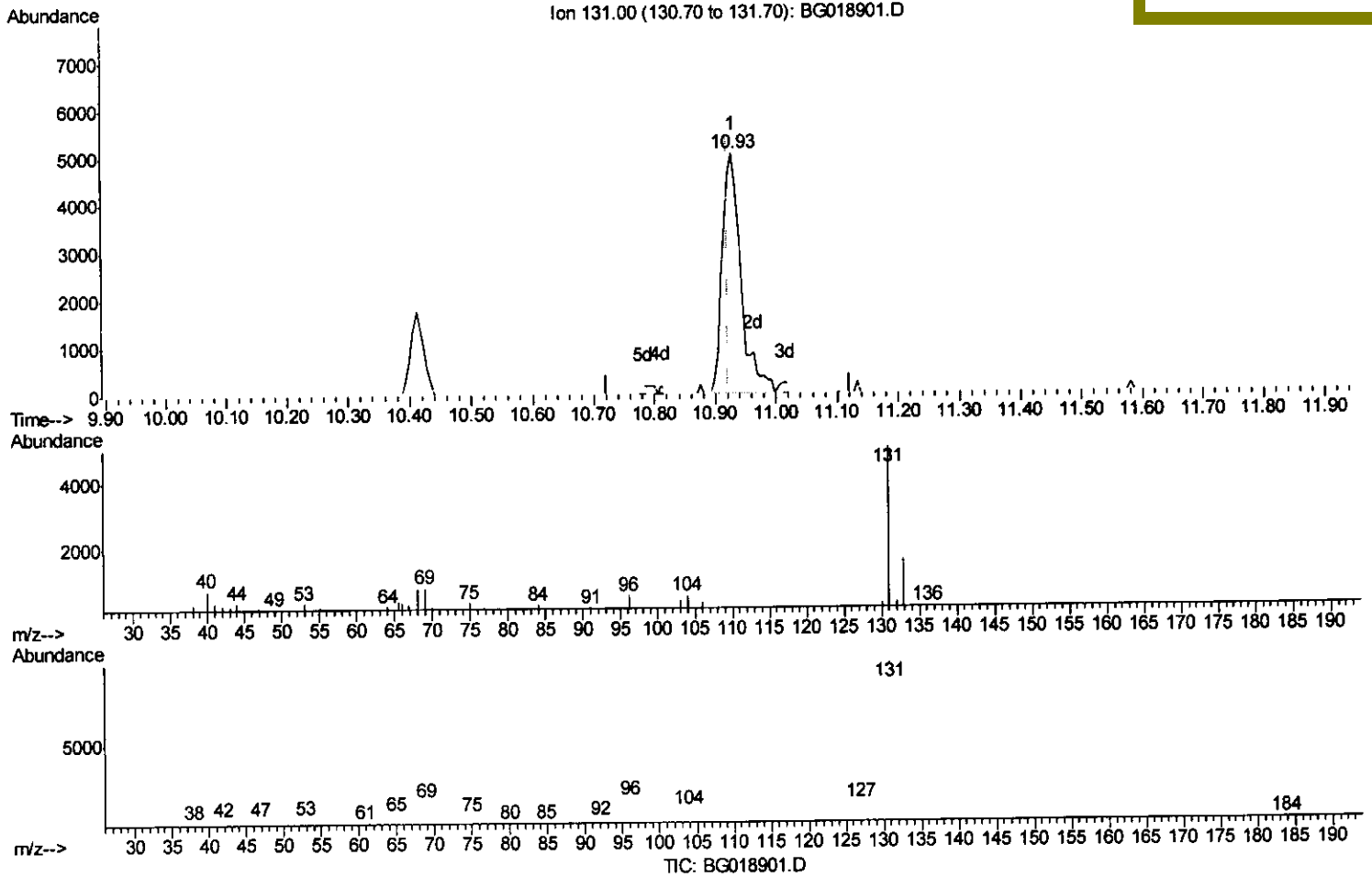
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Manual Integrations
 APPROVED

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(29) 4-Chloroaniline-d4 (S)

10.928min (+0.006) 2.93ng/ul m

response 10807

U.M
 9/29/15

Ion	Exp%	Act%
131.00	100	100
133.00	32.00	31.77
69.00	18.90	16.02
0.00	0.00	0.00

Quantitation Report (Qedit)

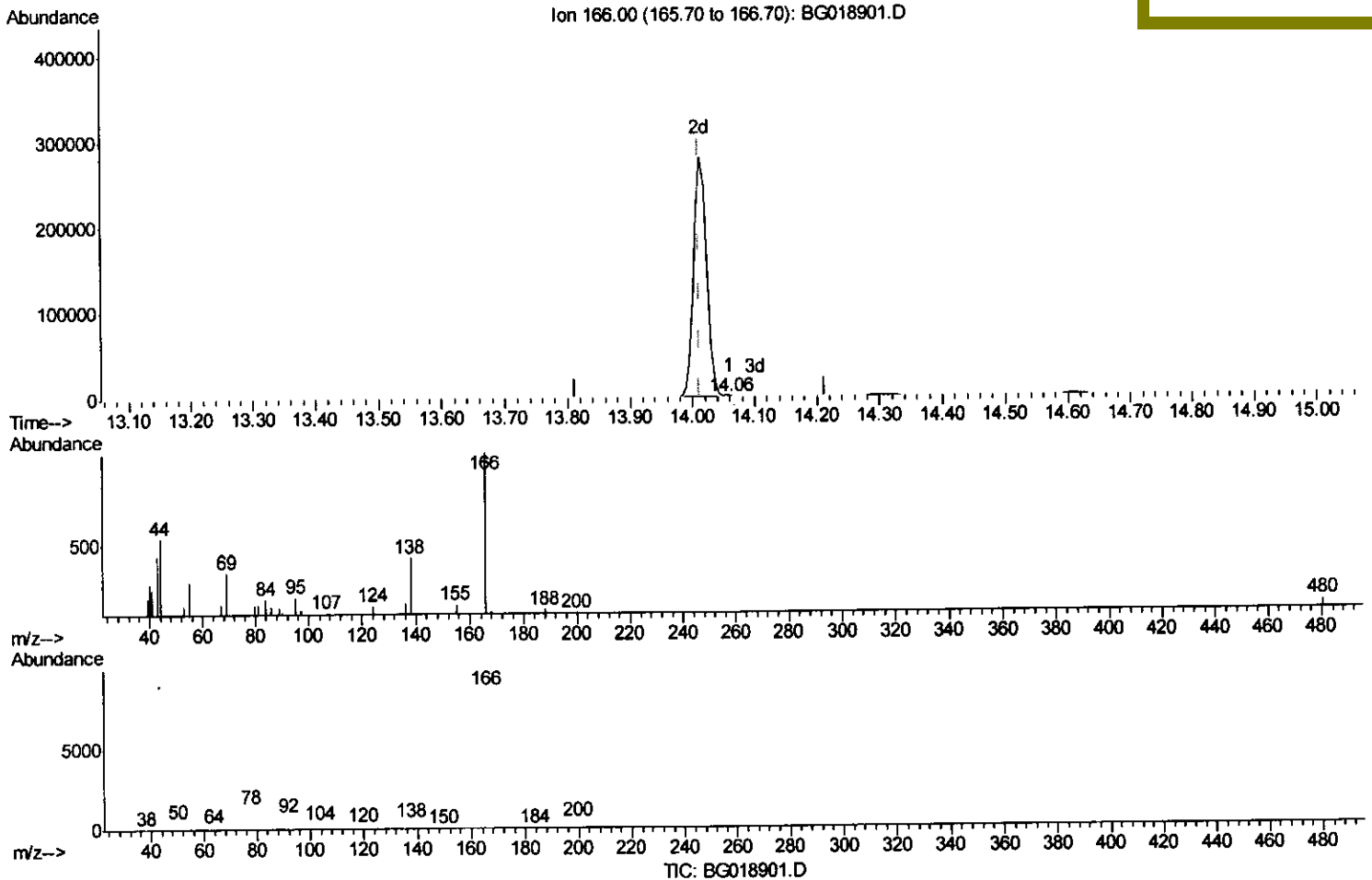
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Manual Integrations
 APPROVED

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(44) Dimethylphthalate-d6 (S)

14.059min (+0.047) 0.05ng/ul

response 582

Ion	Exp%	Act%
166.00	100	100
78.00	17.70	16.08
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

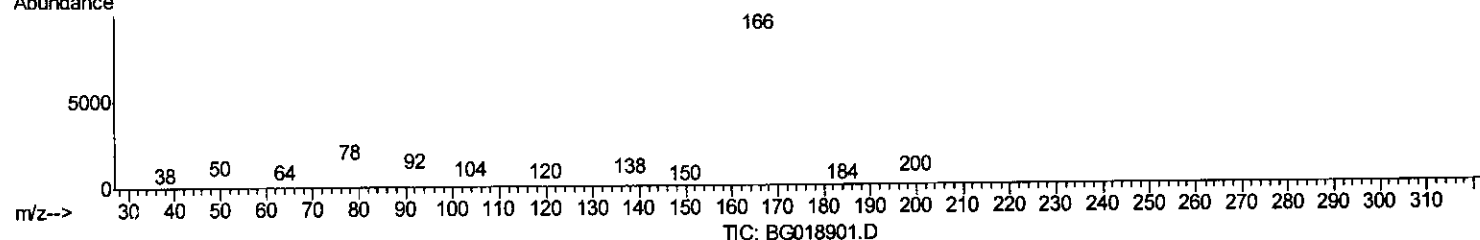
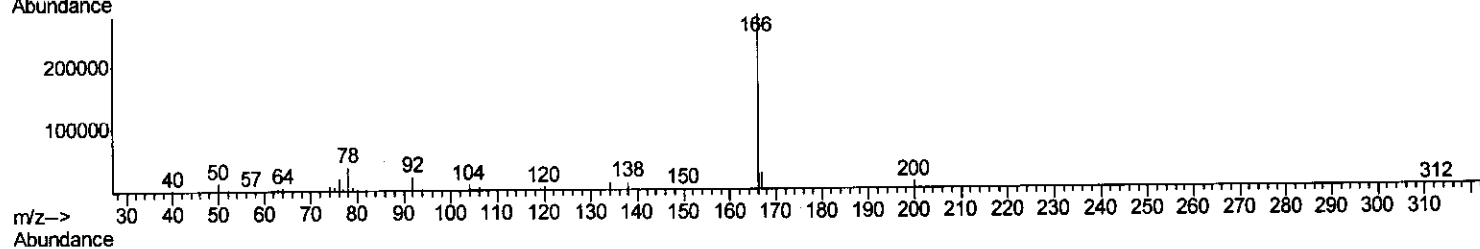
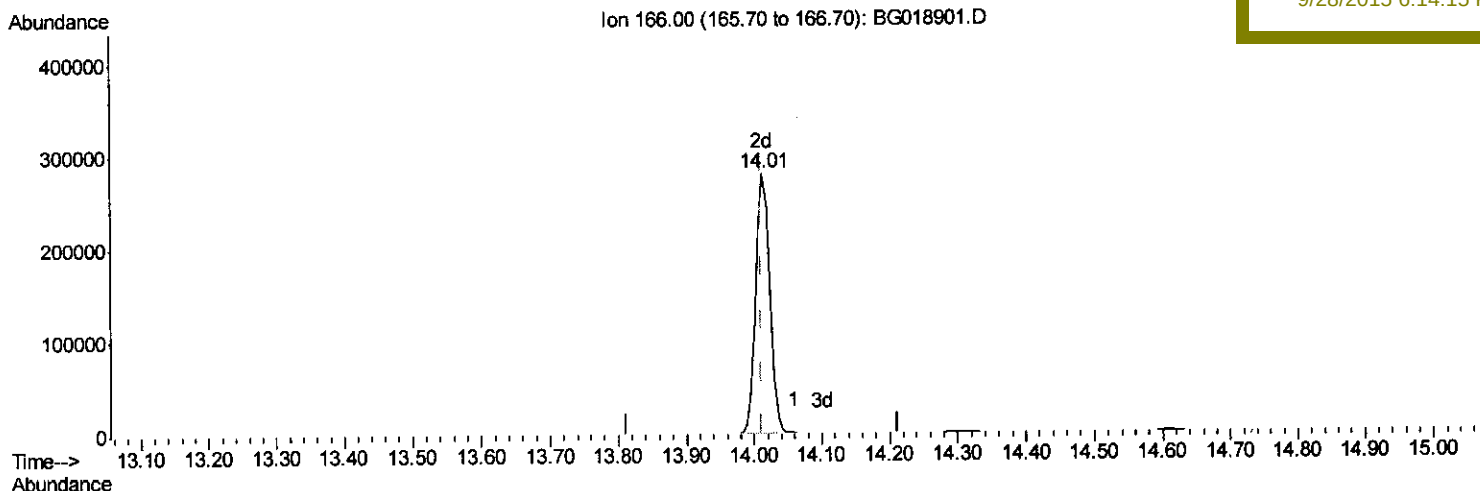
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 Operator : UM/NP
 Sample : G3800-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C03F9

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 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 26 04:18:39 2015
 Response via : Initial Calibration

Manual Integrations
 APPROVED

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(44) Dimethylphthalate-d6 (S)

14.012min (+0.000) 36.47ng/ul m U.M
 response 403908 9/29/15

Ion	Exp%	Act%
166.00	100	100
78.00	17.70	13.45#
0.00	0.00	0.00
0.00	0.00	0.00

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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	7.99	152	43209	20.00	ng/ul	0.00
18) Naphthalene-d8	10.79	136	197855	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.61	164	137614	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.36	188	377185	20.00	ng/ul	0.00
78) Chrysene-d12	21.60	240	425574	20.00	ng/ul	0.00
86) Perylene-d12	24.78	264	439021	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	946	1.05	ng/uL	0.00
5) Phenol-d5	7.16	99	26984	7.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.31	67	61865	30.09	ng/ul	0.00
9) 2-Chlorophenol-d4	7.53	132	73901	27.50	ng/ul	0.00
13) 4-Methylphenol-d8	8.69	113	55055	19.12	ng/ul	0.00
19) Nitrobenzene-d5	9.15	128	46655	31.12	ng/ul	0.00
22) 2-Nitrophenol-d4	9.87	143	53777	35.68	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.41	165	106070	33.42	ng/ul	0.00
29) 4-Chloroaniline-d4	10.93	131	10807m>	2.93	ng/ul	0.00
44) Dimethylphthalate-d6	14.01	166	403908m>	36.47	ng/ul	0.00
47) Acenaphthylene-d8	14.31	160	381802	29.09	ng/ul	0.00
52) 4-Nitrophenol-d4	14.81	143	15596	7.34	ng/ul	0.00
58) Fluorene-d10	15.60	176	250925	25.91	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.72	200	54457	30.50	ng/ul	0.00
71) Anthracene-d10	17.45	188	211736	12.75	ng/ul	0.00
79) Pyrene-d10	19.73	212	413494	21.12	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.56	264	352497	16.50	ng/ul	0.00

U.M
9/29/15

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