

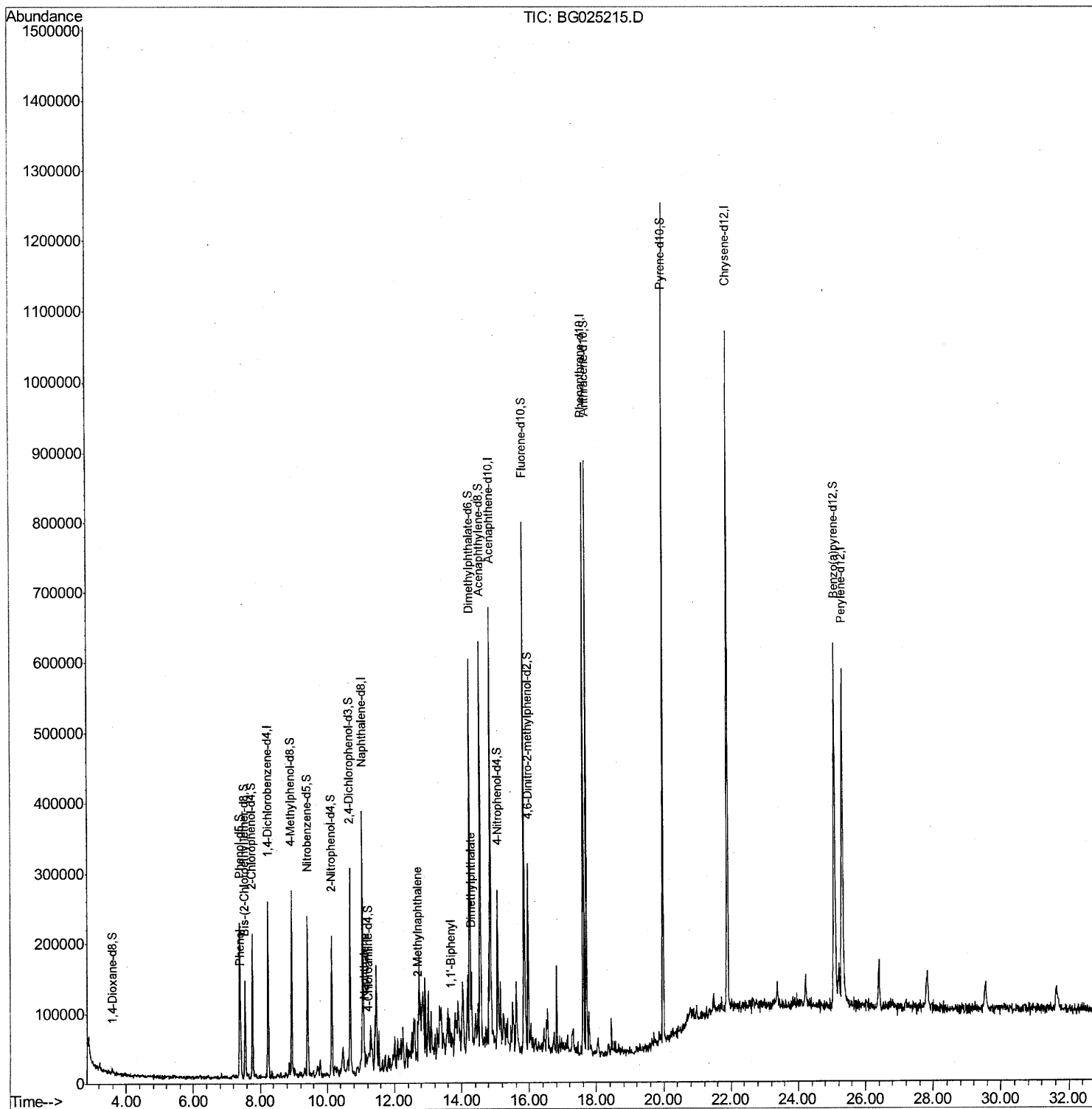
Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
DA869

Manual Integrations
APPROVED

UMANGI
12/16/2016 6:21:17 PM

Quant Time: Dec 16 05:54:34 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 16 03:23:37 2016
Response via : Initial Calibration



Quantitation Report (Qedit)

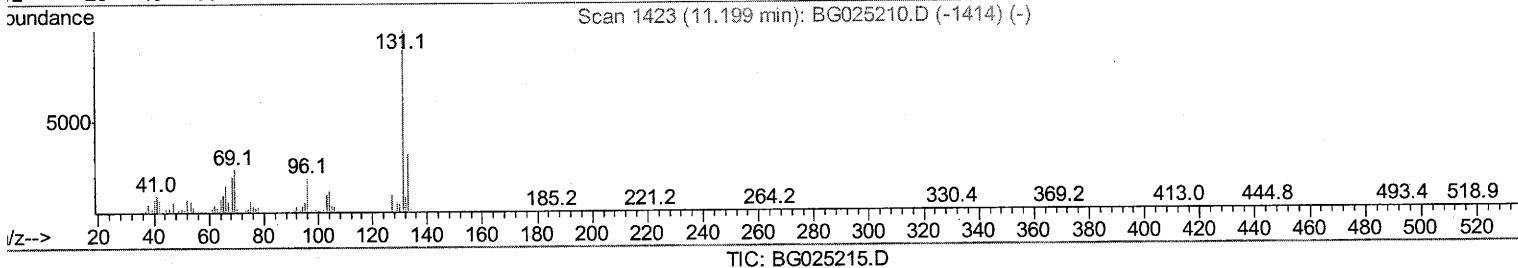
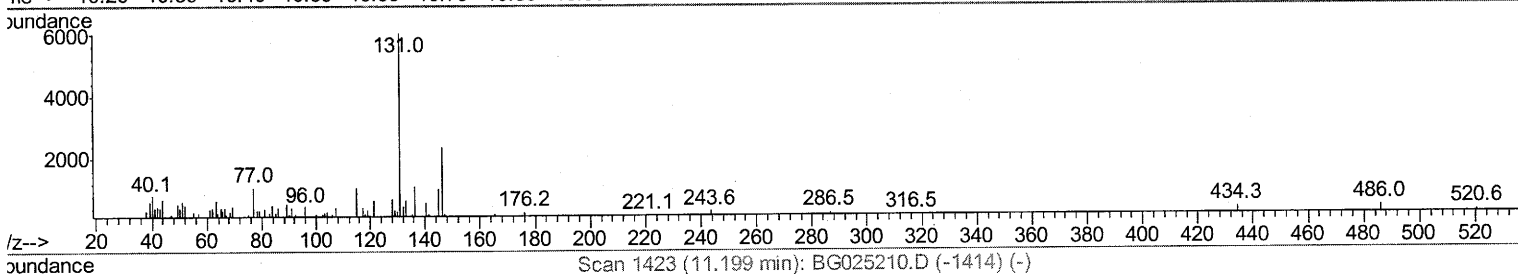
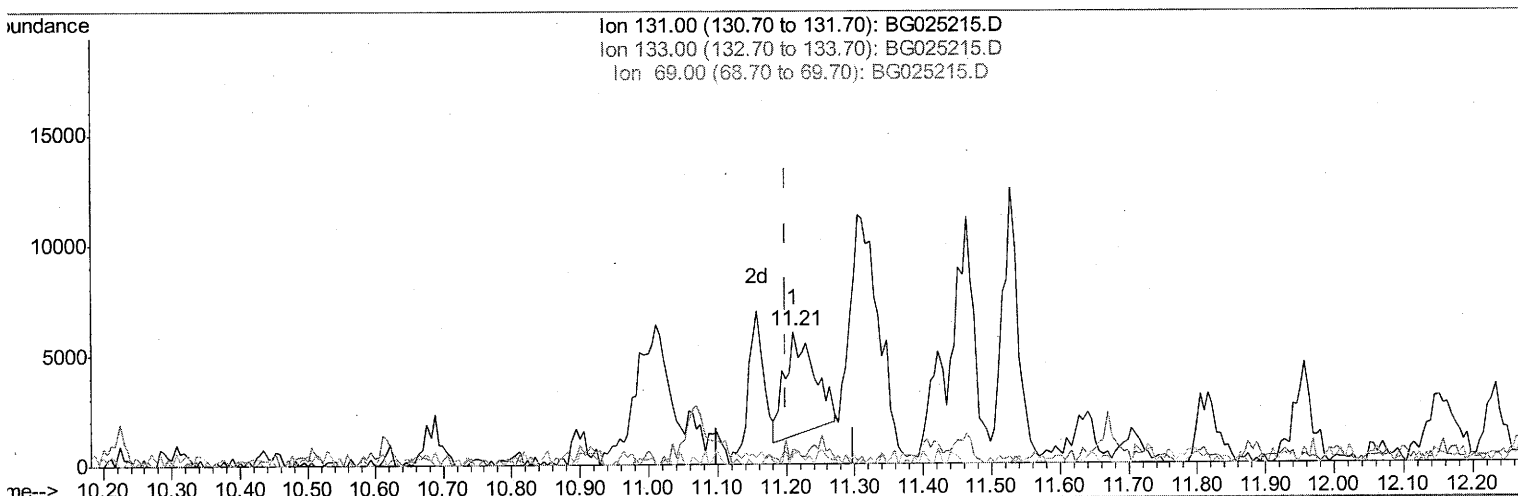
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(29) 4-Chloroaniline-d4 (S)

11.211min (+0.011) 2.94ng/ul

response 13832

Ion	Exp%	Act%
131.00	100	100
133.00	35.40	11.00#
69.00	23.80	7.68#
0.00	0.00	0.00

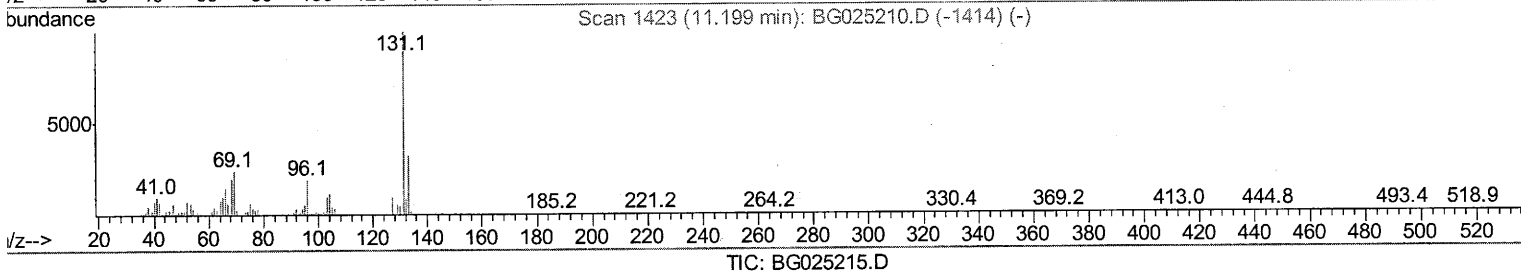
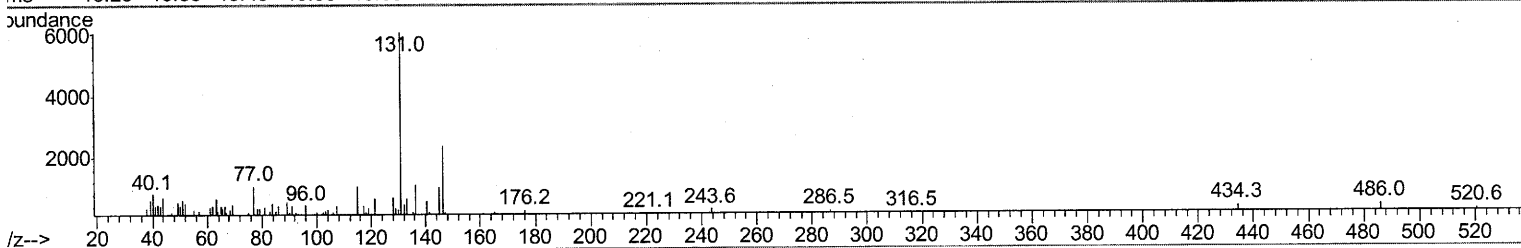
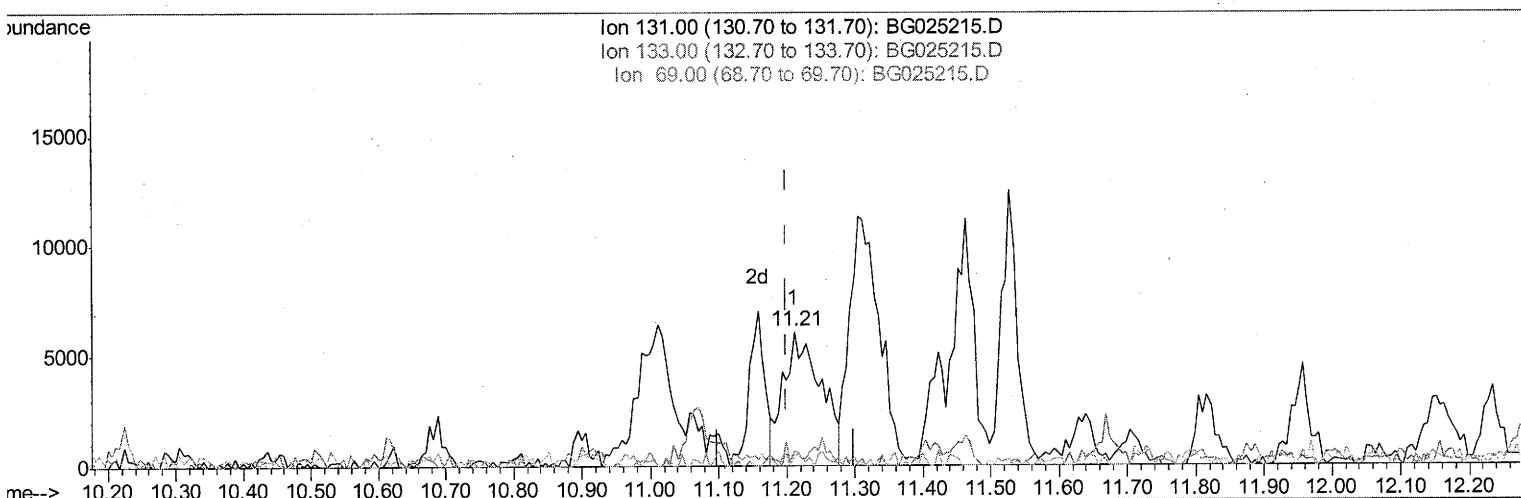
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Manual Integrations
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Response via : Initial Calibration



(29) 4-Chloroaniline-d4 (S)

11.211min (+0.011) 4.87ng/ul m

response 22913

Ion	Exp%	Act%
131.00	100	100
133.00	35.40	11.00#
69.00	23.80	7.68#
0.00	0.00	0.00

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Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
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 BNA_G
 Client Sampled :
 DA869

Manual Integrations
 APPROVED

UMANGI
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Quant Time: Dec 16 05:54:34 2016
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	63737	20.00	ng/ul	0.00
18) Naphthalene-d8	11.06	136	281769	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.86	164	234217	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.60	188	520870	20.00	ng/ul	0.00
75) Chrysene-d12	21.89	240	652037	20.00	ng/ul	0.00
83) Perylene-d12	25.31	264	626464	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.59	96	4290	3.48	ng/uL	0.00
5) Phenol-d5	7.39	99	115807	21.06	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.55	67	73947	21.74	ng/ul	0.00
9) 2-Chlorophenol-d4	7.76	132	81619	21.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.94	113	100367	21.84	ng/ul	0.00
19) Nitrobenzene-d5	9.41	128	42603	21.35	ng/ul	0.00
22) 2-Nitrophenol-d4	10.14	143	53454	21.49	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.68	165	101103	21.13	ng/ul	0.00
29) 4-Chloroaniline-d4	11.21	131	22913m	4.87	ng/ul	0.01
43) Dimethylphthalate-d6	14.25	166	345458	21.44	ng/ul	0.00
46) Acenaphthylene-d8	14.55	160	397711	22.54	ng/ul	0.00
51) 4-Nitrophenol-d4	15.07	143	52206	18.85	ng/ul	0.00
57) Fluorene-d10	15.84	176	315261	20.79	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.97	200	60390	18.75	ng/ul	0.00
70) Anthracene-d10	17.70	188	500236	22.75	ng/ul	0.00
76) Pyrene-d10	19.97	212	649479	24.18	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.08	264	607062	23.92	ng/ul	0.00

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Target Compounds

					Ovalue
6) Phenol	7.41	94	22969	4.07 ng/ul#	94
28) Naphthalene	11.11	128	24608	1.88 ng/ul#	94
34) 2-Methylnaphthalene	12.70	142	21830	2.11 ng/ul	88
40) 1,1'-Biphenyl	13.69	154	19428	1.37 ng/ul#	77
44) Dimethylphthalate	14.29	163	45922	2.90 ng/ul#	95

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