

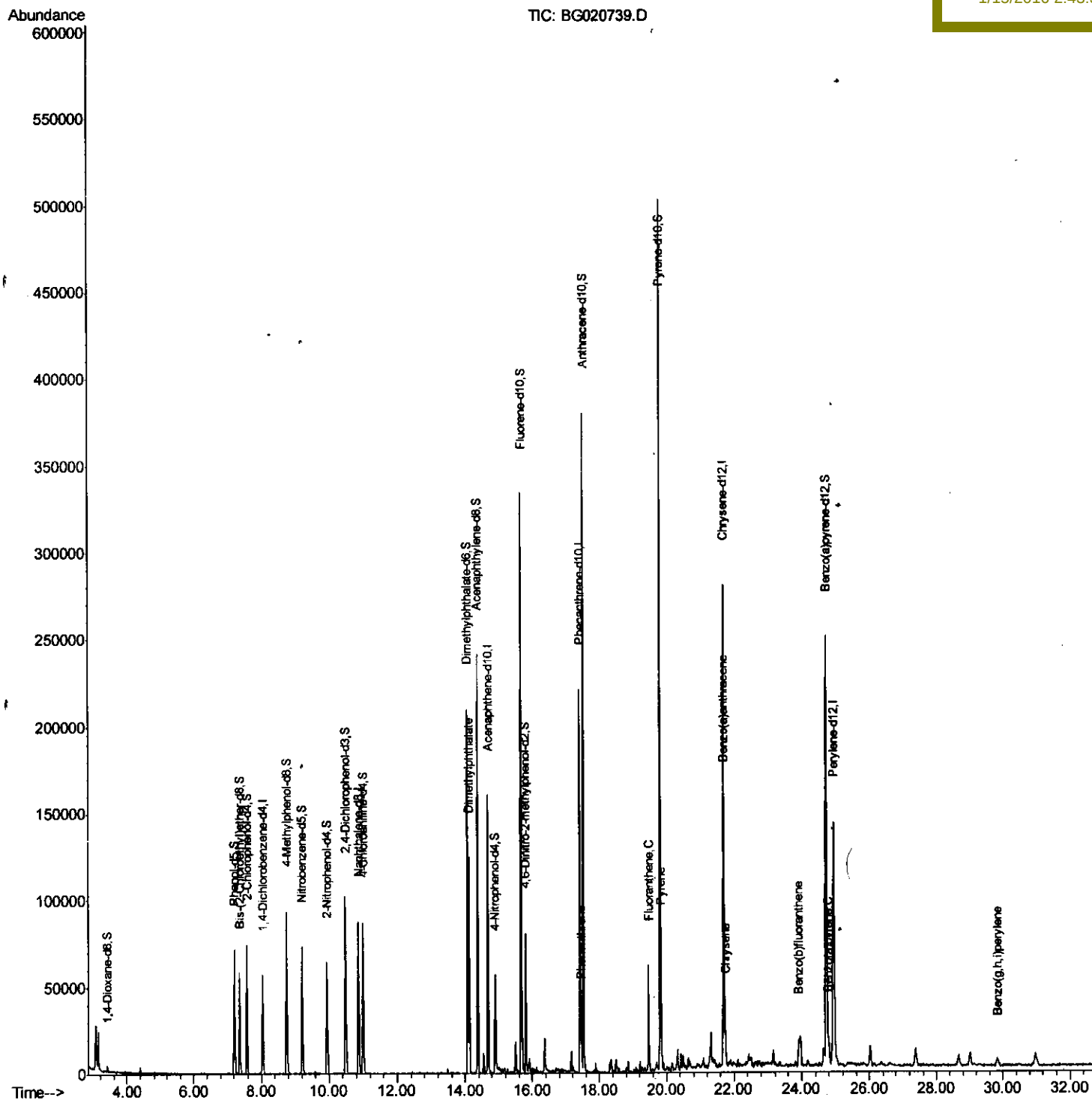
Data Path : Z:\HPCHEM1\BNA_G\Data\BG011416\
Data File : BG020739.D
Acq On : 15 Jan 2016 1:06
Operator : SJ/IZ
Sample : H1101-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Quant Time: Jan 15 02:12:51 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011216.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jan 14 23:48:02 2016
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
BC9G9

Manual Integrations
APPROVED

mohammad
1/15/2016 2:43:36 PM



Quantitation Report (Qedit)

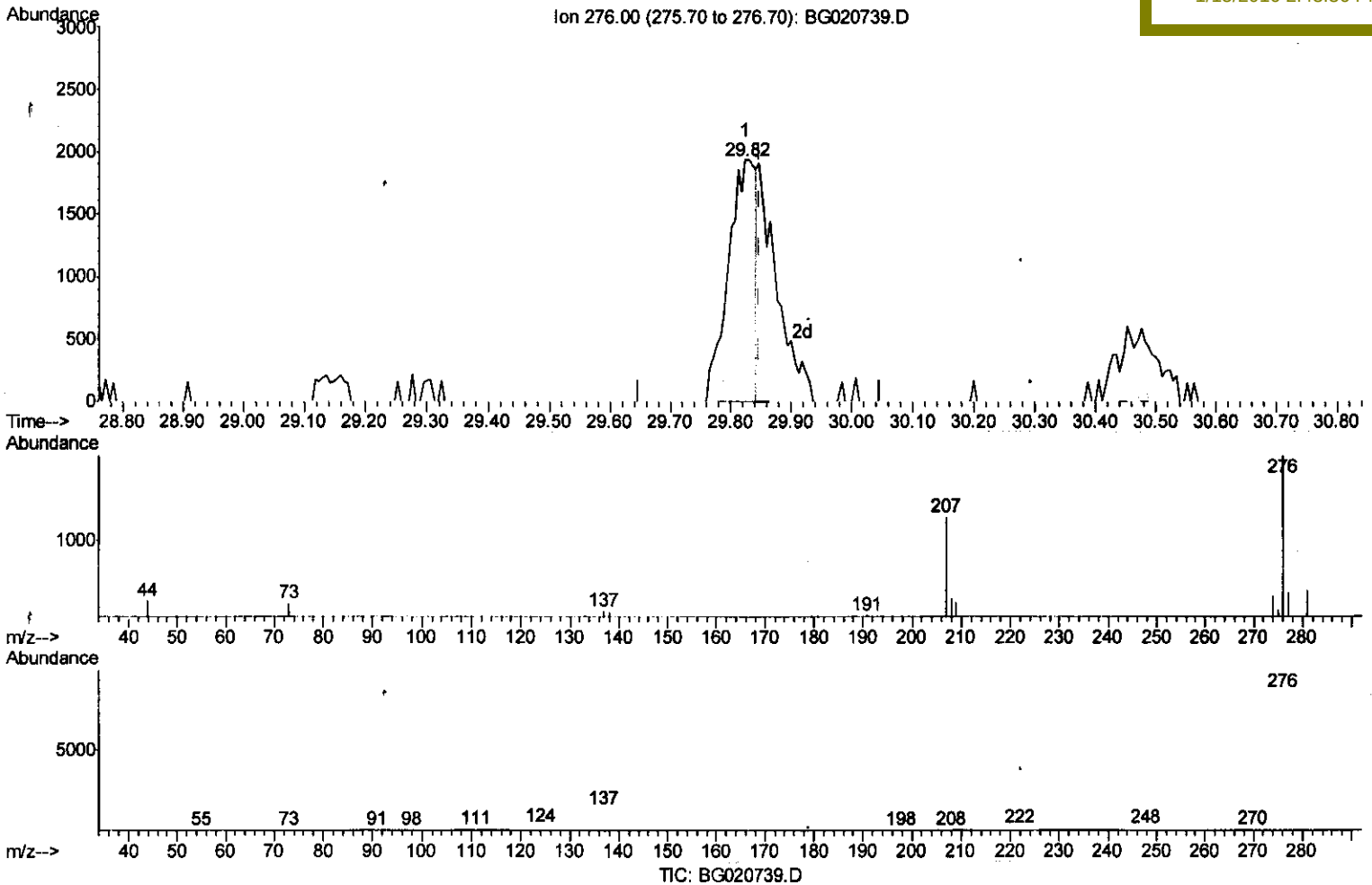
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Quant Time: Jan 15 02:07:52 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 14 23:48:02 2016
 Response via : Initial Calibration

Manual Integrations
 APPROVED

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(94) Benzo(g,h,i)perylene

29.824min (-0.023) 0.64ng/ul

response 6102

Ion	Exp%	Act%
276.00	100	100
138.00	14.00	10.58#
277.00	23.80	21.57
0.00	0.00	0.00

Quantitation Report (Qedit)

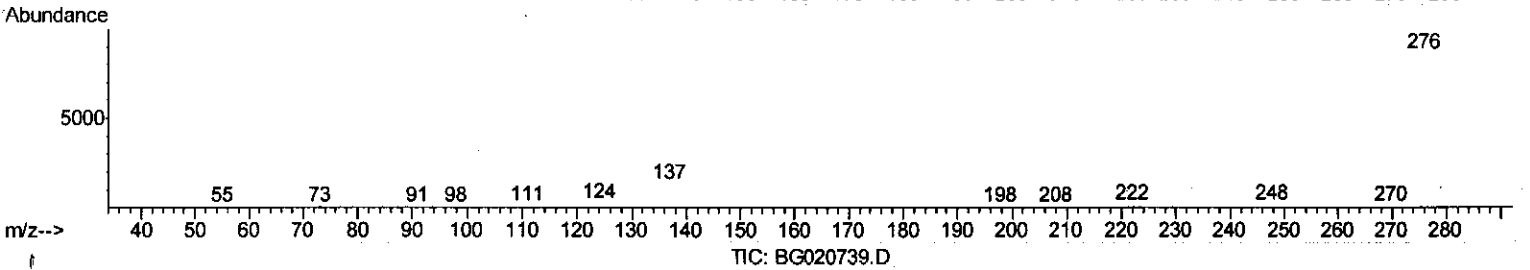
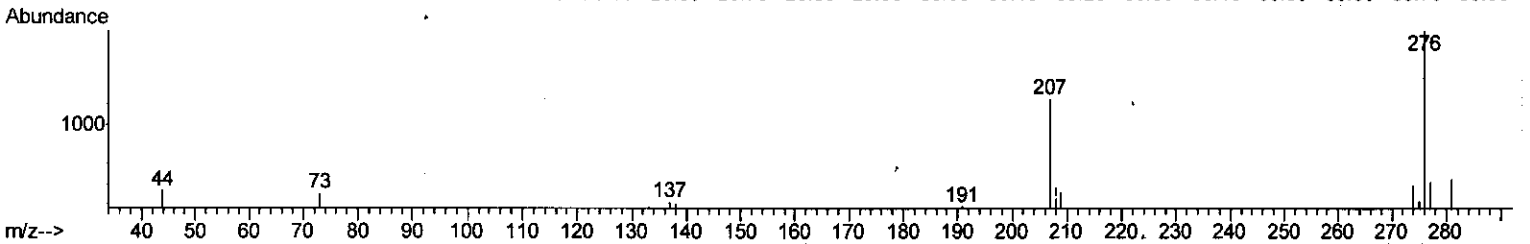
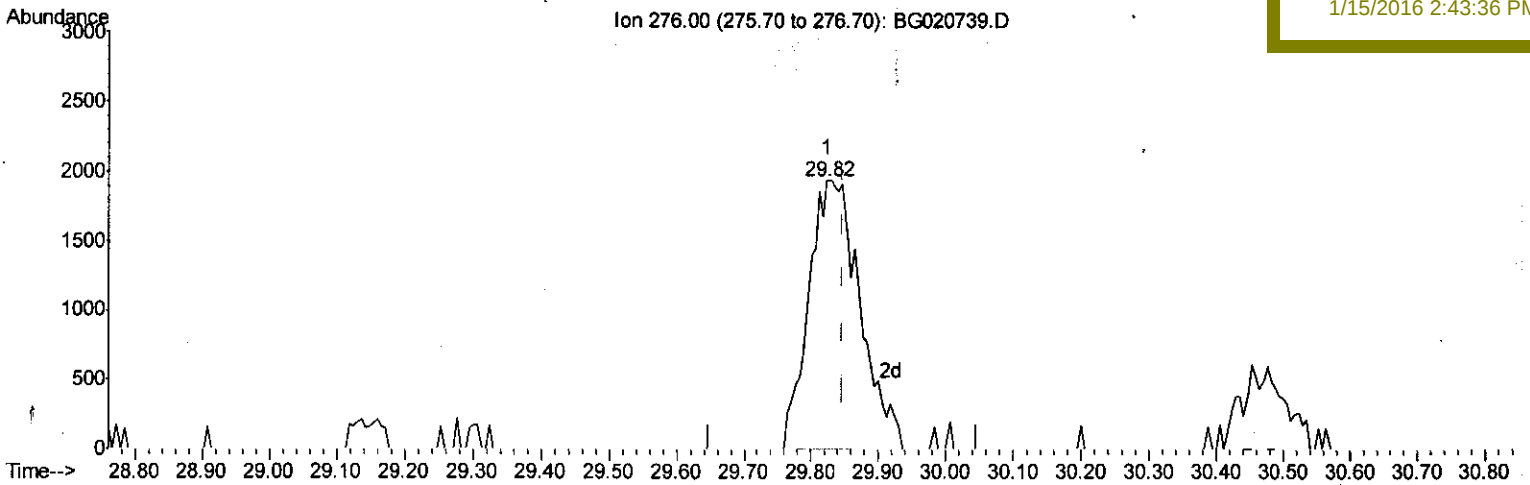
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(94) Benzo(g,h,i)perylene

29.824min (-0.023) 1.07ng/ul m

response 10244

Ion	Exp%	Act%
276.00	100	100
138.00	14.00	10.58#
277.00	23.80	21.57
0.00	0.00	0.00

S.J
 1/16/16

Data Path : Z:\HPCHEM1\BNA_G\Data\BG011416\
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Manual Integrations
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mohammad
 1/15/2016 2:43:36 PM

Quant Time: Jan 15 02:12:51 2016
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 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 14 23:48:02 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	17655	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	75727	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.68	164	54675	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	141509	20.00	ng/ul	0.00
78) Chrysene-d12	21.70	240	170180	20.00	ng/ul	0.00
86) Perylene-d12	24.95	264	158727	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.40	96	1992	5.51	ng/uL	0.00
5) Phenol-d5	7.20	99	45310	29.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.36	67	28678	31.52	ng/ul	0.00
9) 2-Chlorophenol-d4	7.57	132	37158	31.31	ng/ul	0.00
13) 4-Methylphenol-d8	8.75	113	39692	31.10	ng/ul	0.00
19) Nitrobenzene-d5	9.21	128	19895	33.30	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	23345	33.63	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.48	165	42336	32.03	ng/ul	0.00
29) 4-Chloroaniline-d4	10.99	131	53031	40.54	ng/ul	0.00
44) Dimethylphthalate-d6	14.08	166	152128	31.89	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	181469	33.87	ng/ul	0.00
52) 4-Nitrophenol-d4	14.89	143	18763	29.28	ng/ul	0.00
58) Fluorene-d10	15.67	176	145024	34.10	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.80	200	22646	29.42	ng/ul	0.00
71) Anthracene-d10	17.53	188	229325	33.06	ng/ul	0.00
79) Pyrene-d10	19.81	212	269625	33.30	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.73	264	289918	34.27	ng/ul	0.00

Target Compounds

					Qvalue
45) Dimethylphthalate	14.12	163	85692	16.72 ng/ul#	98
70) Phenanthrene	17.47	178	18936	2.25 ng/ul	100
77) Fluoranthene	19.48	202	39649	3.87 ng/ul	99
80) Pyrene	19.84	202	36677	3.40 ng/ul	99
83) Benzo(a)anthracene	21.68	228	21564	2.03 ng/ul	99
85) Chrysene	21.74	228	20947	2.09 ng/ul	99
88) Benzo(b)fluoranthene	23.92	252	25479	2.49 ng/ul	97
91) Benzo(a)pyrene	24.80	252	18295	1.86 ng/ul	98
94) Benzo(g,h,i)perylene	29.82	276	10244m	1.07 ng/ul	

S.J
 1/16/16

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