

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061917\
Quantitation Report (Qedit)

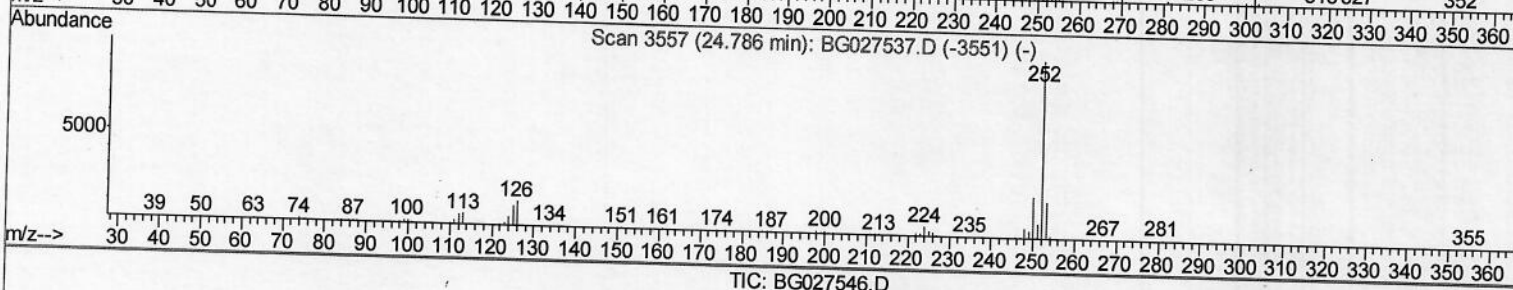
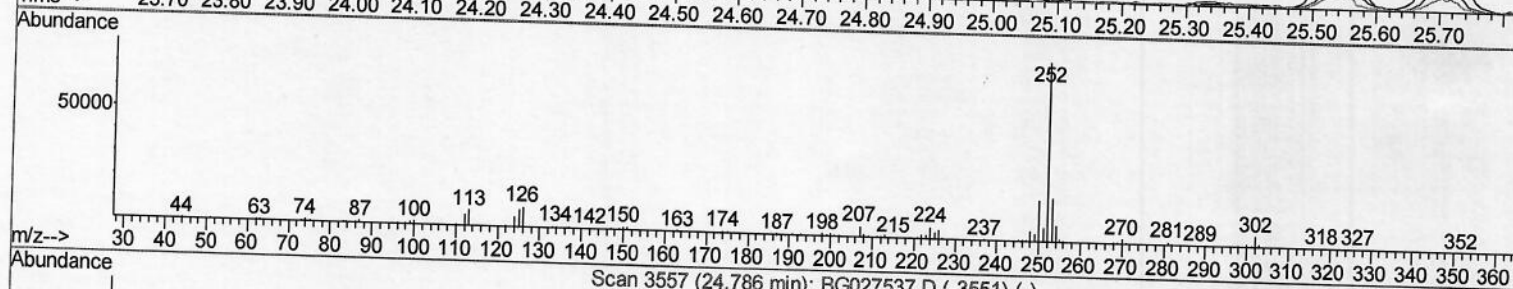
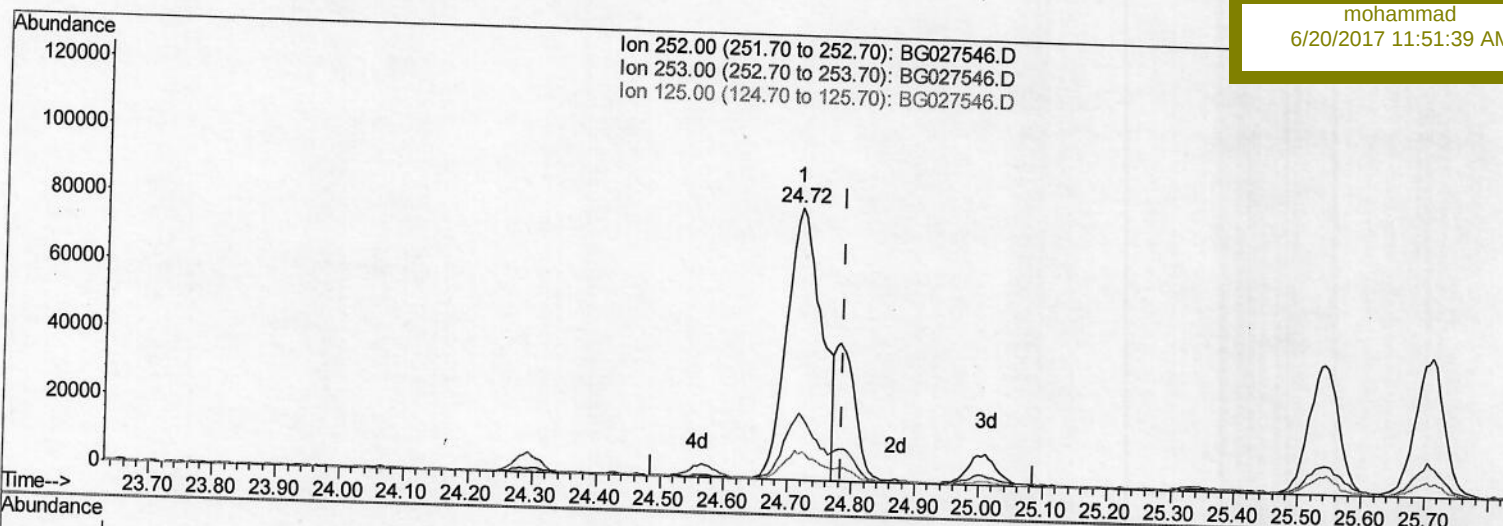
Data File : BG027546.D
Acq On : 20 Jun 2017 2:16
Operator : SJ/MA
Sample : I3627-15 25X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jun 20 03:30:00 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 20 01:50:01 2017
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
F5A88

Manual Integrations
APPROVED

mohammad
6/20/2017 11:51:39 AM



(86) Benzo(k)fluoranthene

24.717min (-0.069) 19.61ng/ul

response 304688

Ion	Exp%	Act%
252.00	100	100
253.00	23.10	24.42
125.00	5.10	9.72#
0.00	0.00	0.00

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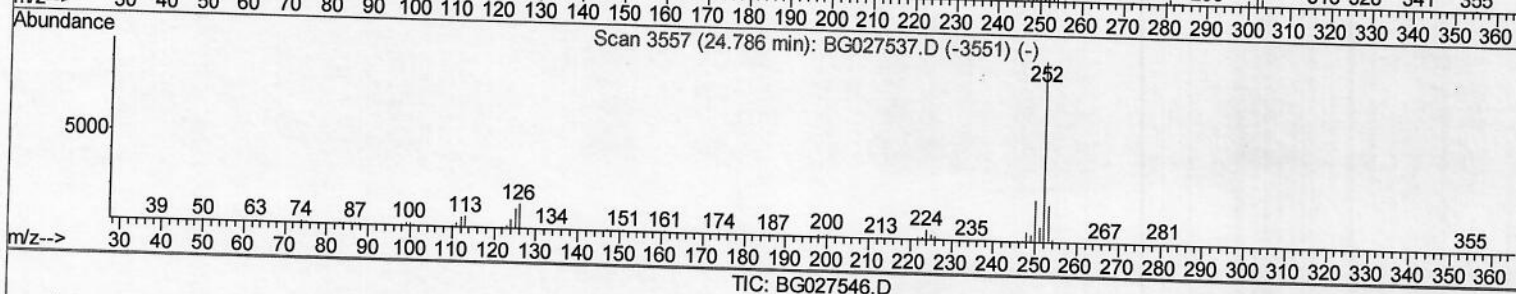
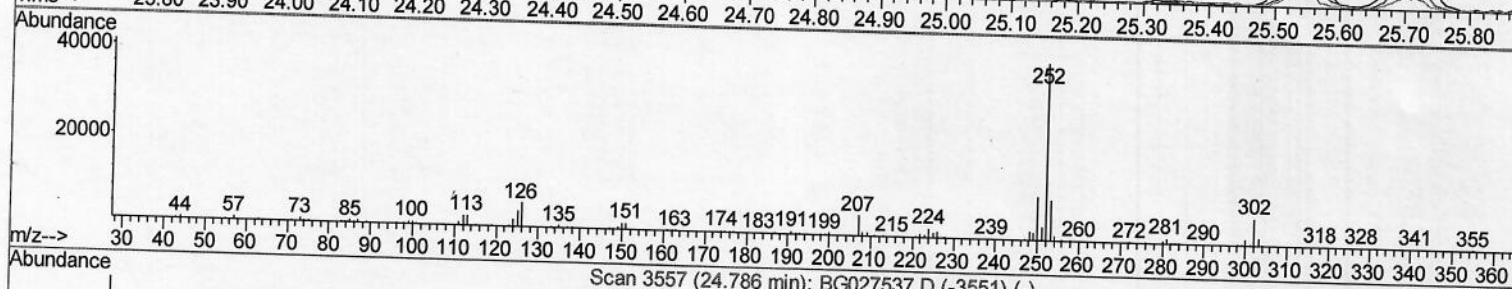
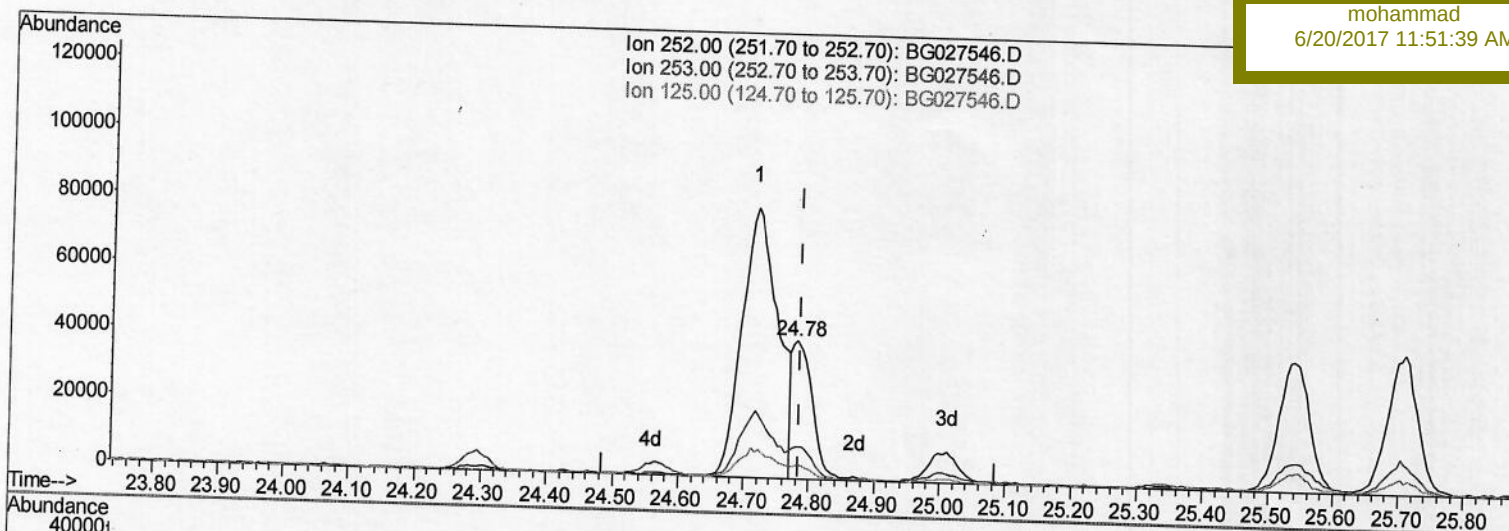
F5A88

Manual Integrations

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(86) Benzo(k)fluoranthene

24.781min (-0.004) 5.60ng/ul m

response 86922

Ion Exp% Act%

252.00 100 100

253.00 23.10 22.87

125.00 5.10 9.59#

0.00 0.00 0.00

S.J
06/30/2017

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 Operator : SJ/MA
 Sample : I3627-15 25X
 Misc :
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Instrument :
 BNA_G
 Client Sample ID :
 F5A88

Quant Time: Jun 20 03:43:20 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 20 01:50:01 2017
 Response via : Initial Calibration

Manual Integrations
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 6/20/2017 11:51:39 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	8.52	152	32443	20.00	ng/ul	0.00
18) Naphthalene-d8	11.34	136	167666	20.00	ng/ul	0.00
35) Acenaphthene-d10	15.10	164	111673	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.85	188	257649	20.00	ng/ul	0.00
75) Chrysene-d12	22.23	240	286408	20.00	ng/ul	0.00
83) Perylene-d12	25.89	264	271368	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)
3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.66	99	2676	0.76	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.84	67	1723	0.72	ng/ul	0.00
9) 2-Chlorophenol-d4	8.06	132	2098	0.90	ng/ul	0.00
13) 4-Methylphenol-d8	9.19	113	2450	0.89	ng/ul	0.00
19) Nitrobenzene-d5	9.68	128	985	0.81	ng/ul	0.00
22) 2-Nitrophenol-d4	10.40	143	941	0.73	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.94	165	2213	0.91	ng/ul	0.00
29) 4-Chloroaniline-d4	11.47	131	1653	0.53	ng/ul	0.00
43) Dimethylphthalate-d6	14.49	166	8823	0.94	ng/ul	0.00
46) Acenaphthylene-d8	14.80	160	9616	0.89	ng/ul	0.00
51) 4-Nitrophenol-d4	15.29	143	586	0.32	ng/ul	0.01
57) Fluorene-d10	16.09	176	8737	1.00	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
70) Anthracene-d10	17.95	188	13221	1.10	ng/ul	0.00
76) Pyrene-d10	20.22	212	15256	1.09	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.62	264	12986	1.06	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
69) Phenanthrene	17.90	178	48450	3.50	ng/ul	96
71) Anthracene	17.99	178	98118	6.97	ng/ul	97
72) Carbazole	18.25	167	12173	1.01	ng/ul	99
74) Fluoranthene	19.89	202	669237	42.96	ng/ul#	88
77) Pyrene	20.25	202	640211	37.82	ng/ul#	87
80) Benzo(a)anthracene	22.21	228	267857	16.79	ng/ul	98
82) Chrysene	22.28	228	286274	19.87	ng/ul	99
85) Benzo(b)fluoranthene	24.72	252	304688	18.85	ng/ul#	93
86) Benzo(k)fluoranthene	24.78	252	86922m	5.60	ng/ul	
88) Benzo(a)pyrene	25.71	252	131054	8.45	ng/ul#	97
89) Indeno(1,2,3-cd)pyrene	30.08	276	71595	4.04	ng/ul#	89
90) Dibenzo(a,h)anthracene	30.13	278	22470	1.48	ng/ul#	92
91) Benzo(g,h,i)perylene	31.40	276	47172	3.25	ng/ul#	95

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

> 5.5
 06/30/2017

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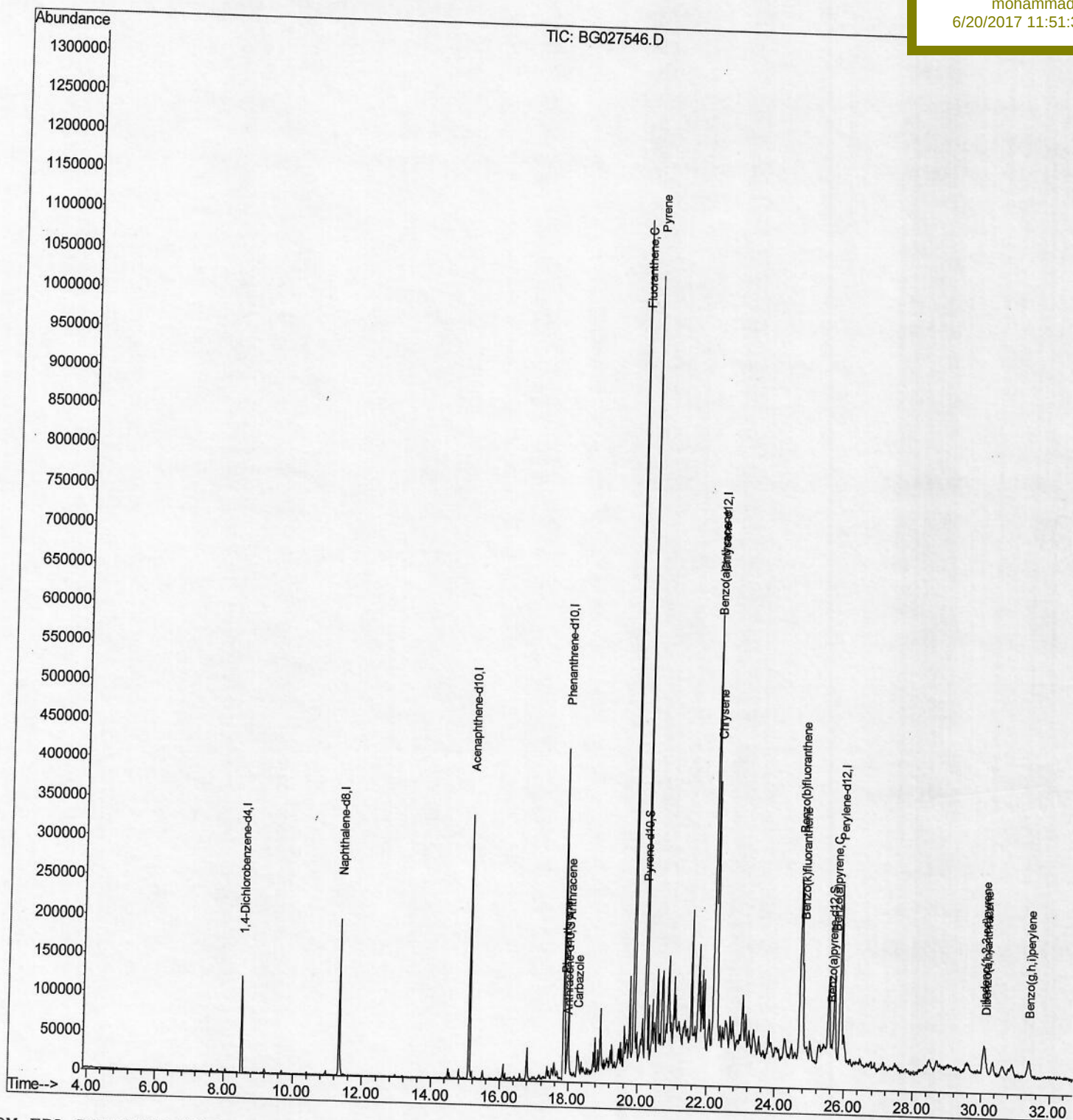
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SOM-EPA-BG061617.M Tue Jun 20 03:44:12 2017

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