

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG042817\

Quantitation Report (Qedit)

Data File : BG026752.D

Acq On : 28 Apr 2017 18:24

Operator : SJ/MA

Sample : I2816-08

Misc :

ALS Vial : 11 Sample Multiplier: 1

Quant Time: Apr 29 00:40:00 2017

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG042717.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat Apr 29 00:39:01 2017

Response via : Initial Calibration

Instrument :

BNA_G

ClientSampleId :

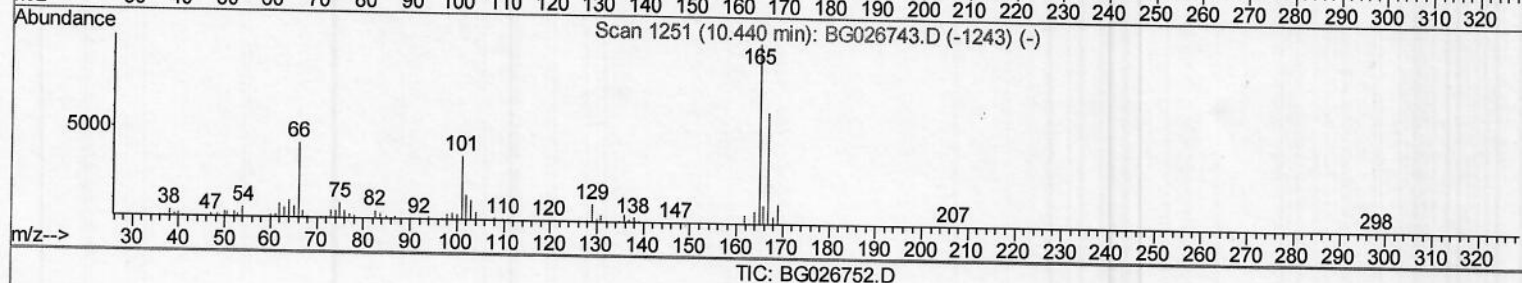
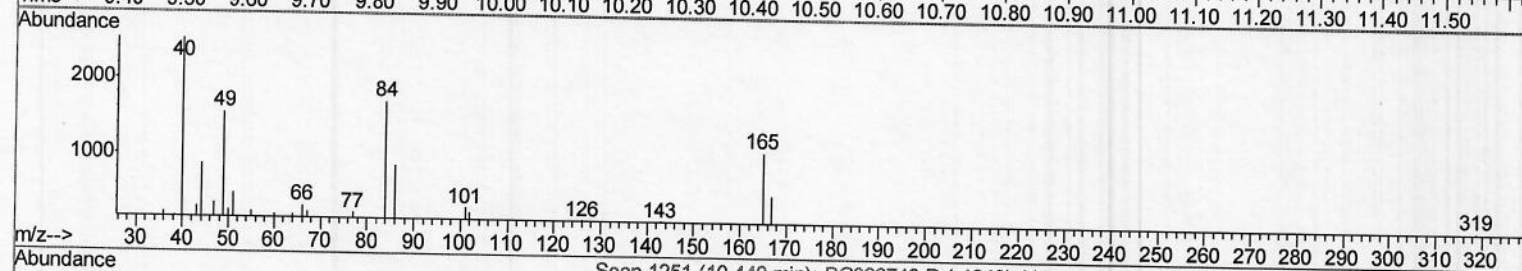
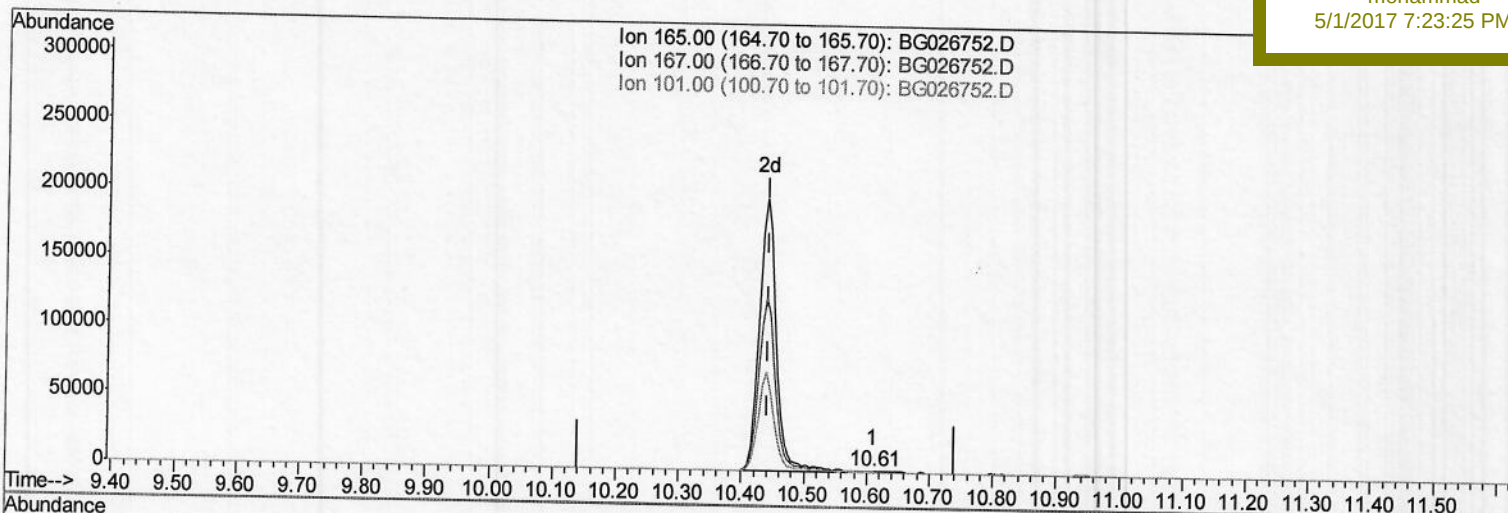
JHT20

Manual Integrations

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mohammad

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(26) 2,4-Dichlorophenol-d3 (S)

10.609min (+0.169) 0.04ng/ul

response 400

Ion	Exp%	Act%
165.00	100	100
167.00	58.70	48.55
101.00	29.80	28.36
0.00	0.00	0.00

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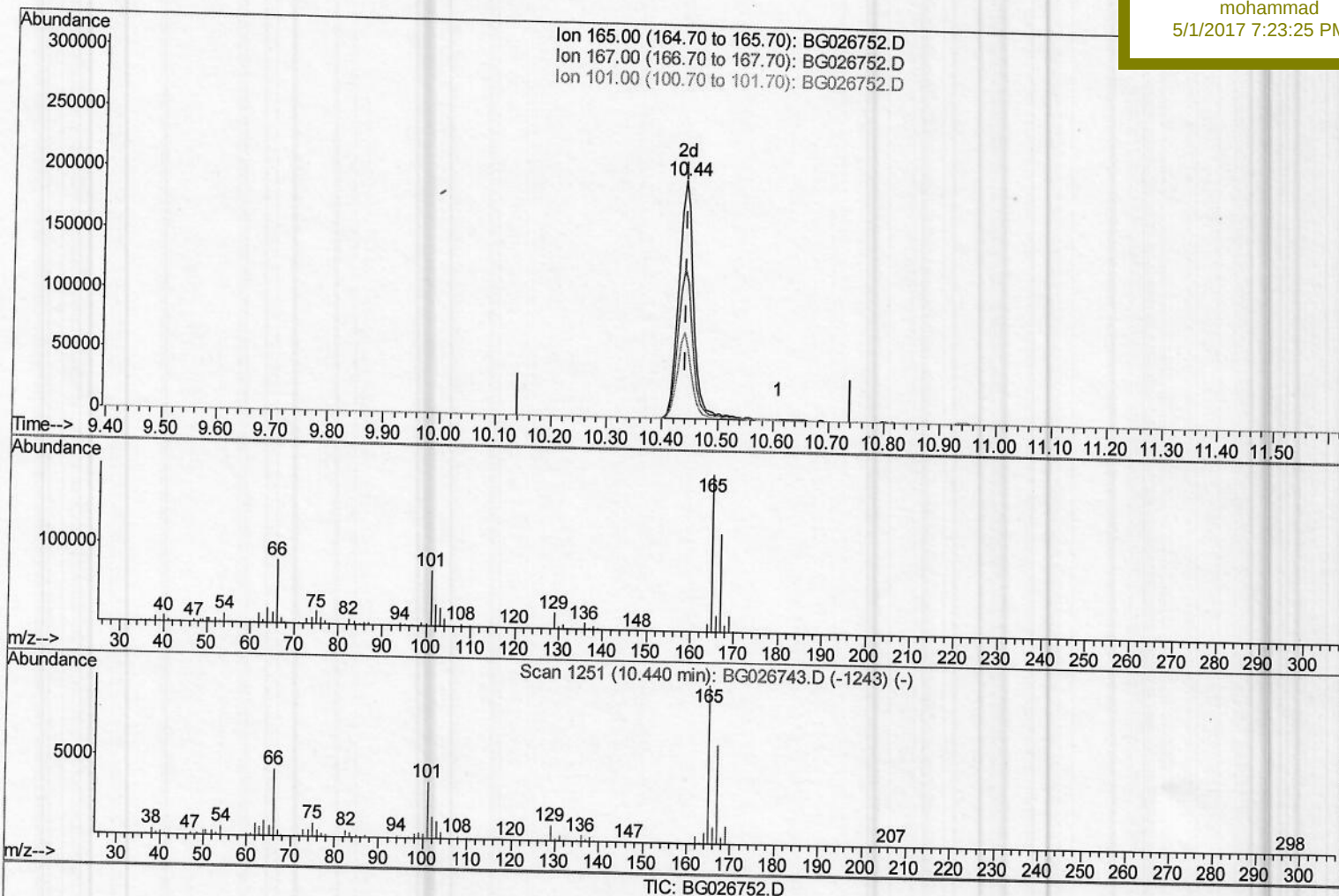
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(26) 2,4-Dichlorophenol-d3 (S)

10.439min (-0.001) 36.18ng/ul m

response 366926

Ion Exp% Act%

165.00 100 100

167.00 58.70 63.19

101.00 29.80 36.01#

0.00 0.00 0.00

> S.J
05/04/17

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Quant Time: Apr 29 01:06:29 2017

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	137914	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	702085	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.62	164	412677	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.35	188	813041	20.00	ng/ul	0.00
75) Chrysene-d12	21.60	240	731924	20.00	ng/ul	0.00
83) Perylene-d12	24.76	264	752480	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	20493	6.69	ng/uL	0.00
5) Phenol-d5	7.19	99	457904	37.20	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.33	67	262242	36.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	366594	35.09	ng/ul	0.00
13) 4-Methylphenol-d8	8.72	113	397885	38.95	ng/ul	0.00
19) Nitrobenzene-d5	9.16	128	207306	35.32	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	233774	36.63	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.44	165	366926m	36.18	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	591691	43.59	ng/ul	0.00
43) Dimethylphthalate-d6	14.02	166	1222555	36.81	ng/ul	0.00
46) Acenaphthylene-d8	14.32	160	1523634	36.52	ng/ul	0.00
51) 4-Nitrophenol-d4	14.84	143	198407	33.14	ng/ul	0.00
57) Fluorene-d10	15.61	176	1046801	36.11	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.73	200	153656	31.12	ng/ul	0.00
70) Anthracene-d10	17.45	188	1472408	37.32	ng/ul	0.00
76) Pyrene-d10	19.73	212	1400541	36.25	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.54	264	1396455	39.25	ng/ul	0.00

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05/04/17

Target Compounds

					Qvalue
6) Phenol	7.21	94	14113	1.13	ng/ul 89
44) Dimethylphthalate	14.07	163	205582	6.32	ng/ul 98

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

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Client Sampled :

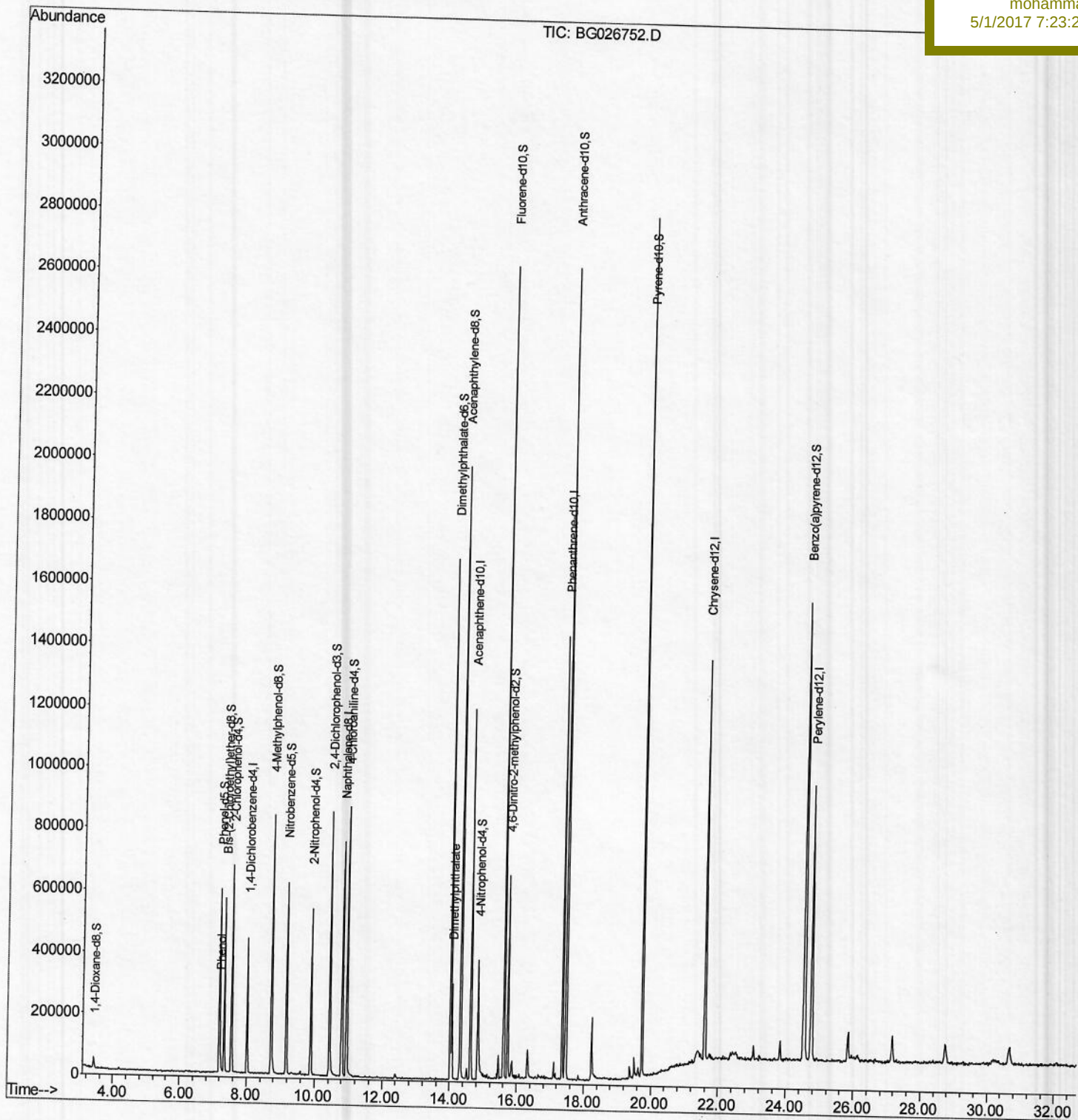
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SOM-EPA-BG042717.M Sat Apr 29 01:07:53 2017

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