

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

Quantitation Report (Qedit)

Data File : BG026750.D

Acq On : 28 Apr 2017 17:06

Operator : SJ/MA

Sample : I2816-06

Misc :

ALS Vial : 9 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

JHSP7

Quant Time: Apr 29 00:39:43 2017

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

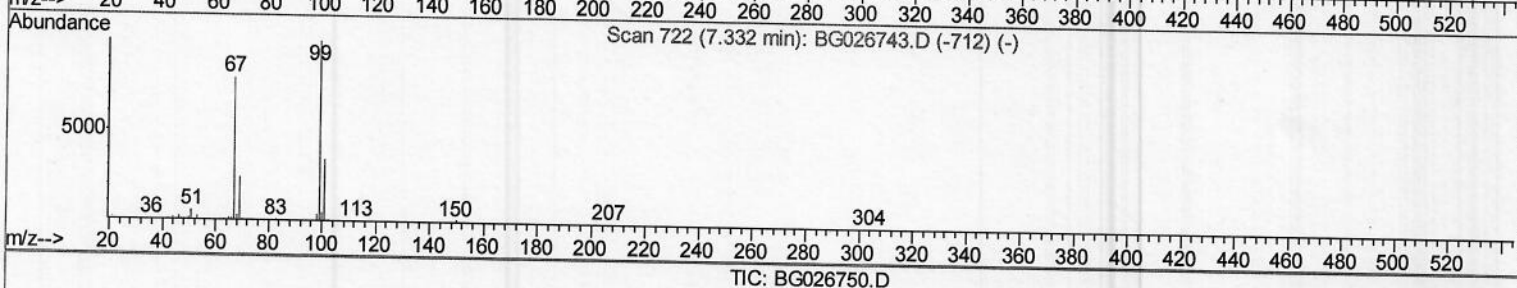
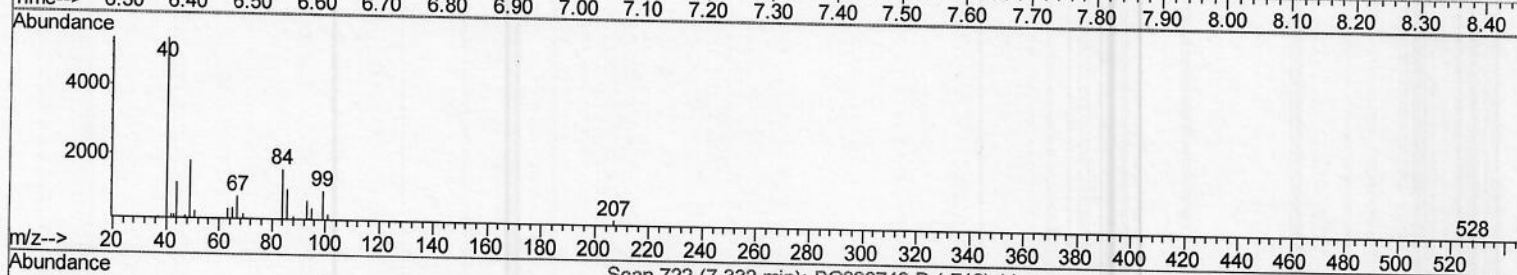
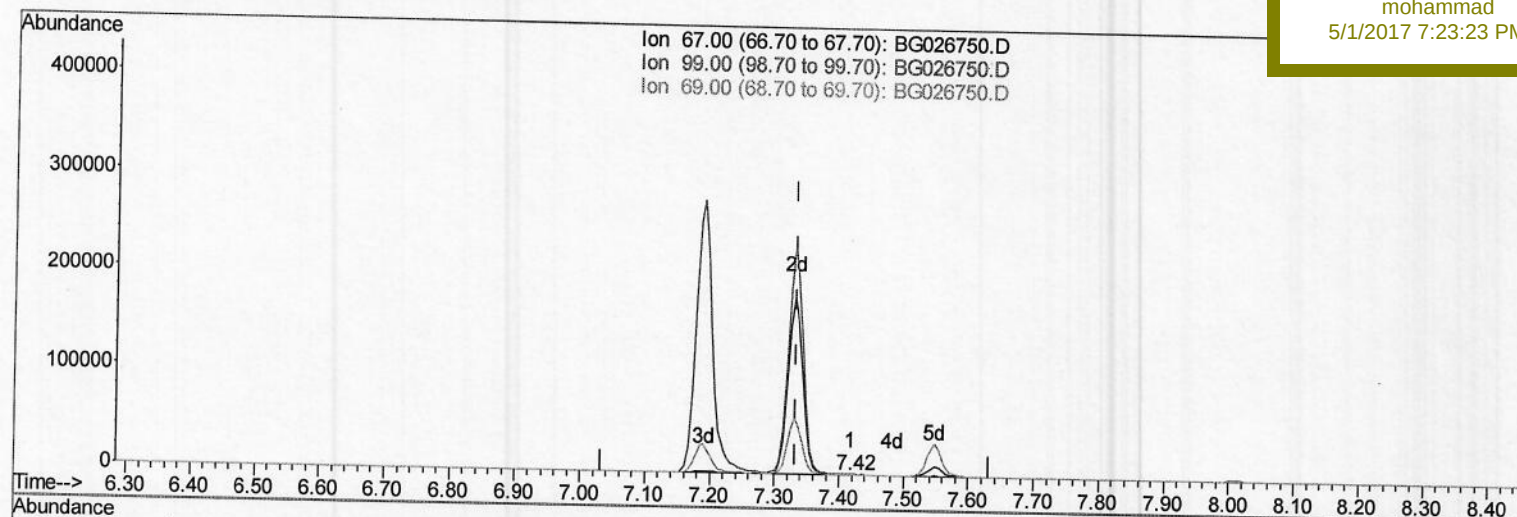
Quant Title : SVOA CALIBRATION

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

Response via : Initial Calibration

Manual Integrations
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5/1/2017 7:23:23 PM



(7) Bis-(2-Chloroethyl)ether-d8 (S)

7.419min (+0.086) 0.14ng/ul

response 1055

Ion	Exp%	Act%
67.00	100	100
99.00	102.80	121.99
69.00	31.30	36.65
0.00	0.00	0.00

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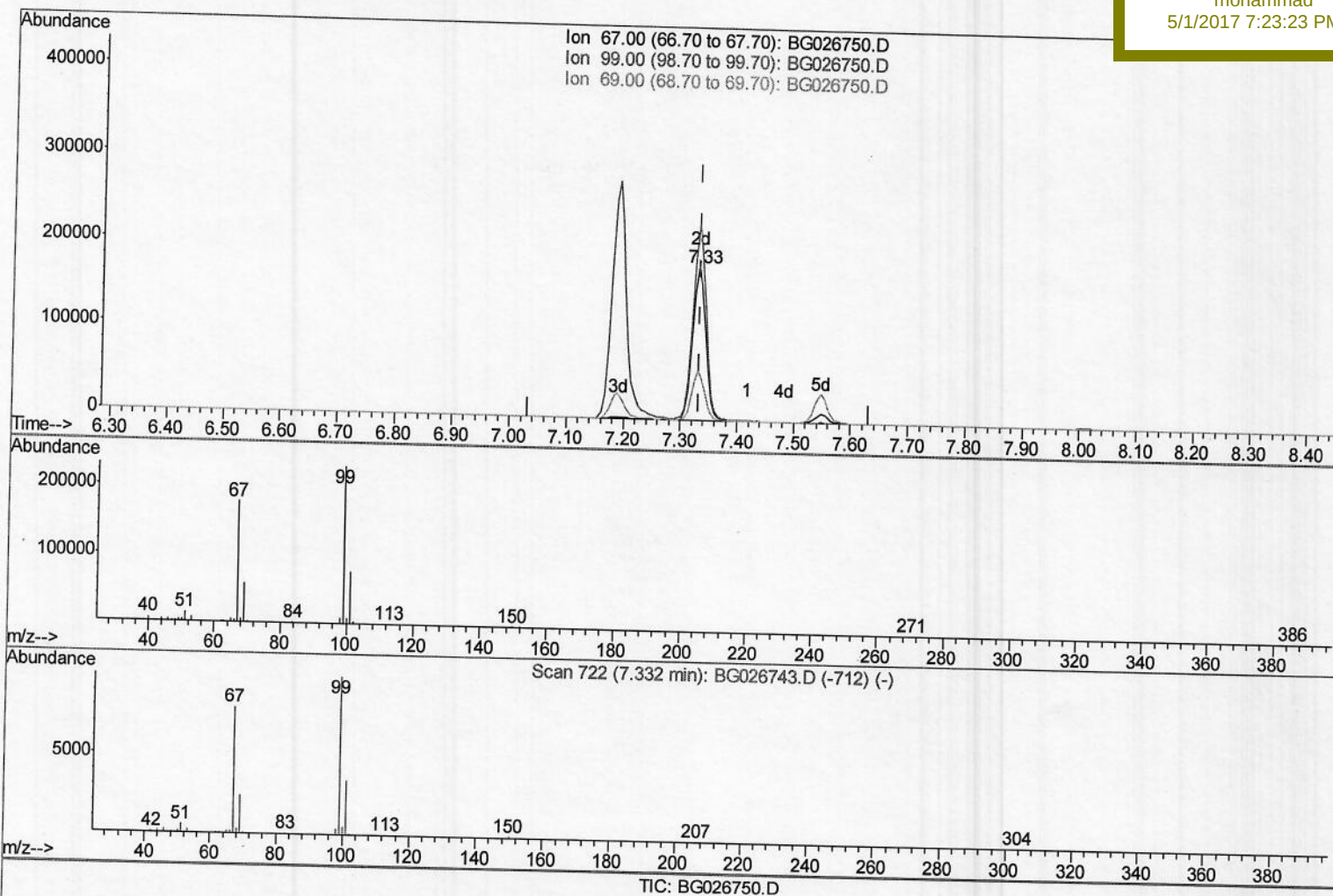
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(7) Bis-(2-Chloroethyl)ether-d8 (S)

7.330min (-0.002) 37.96ng/ul m

response 285199

Ion	Exp%	Act%
67.00	100	100
99.00	102.80	129.67#
69.00	31.30	31.71
0.00	0.00	0.00

> S.J
05/04/17

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Operator : SJ/MA

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Misc :

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	143165	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	723539	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.62	164	419454	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.35	188	825054	20.00	ng/ul	0.00
75) Chrysene-d12	21.60	240	737035	20.00	ng/ul	0.00
83) Perylene-d12	24.76	264	759307	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	22319	7.02	ng/uL	0.00
5) Phenol-d5	7.19	99	493669	38.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.33	67	285199m	37.96	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	404449	37.29	ng/ul	0.00
13) 4-Methylphenol-d8	8.72	113	425951	40.17	ng/ul	0.00
19) Nitrobenzene-d5	9.17	128	217947	36.04	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	251156	38.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.44	165	388723	37.19	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	591138	42.25	ng/ul	0.00
43) Dimethylphthalate-d6	14.02	166	1237417	36.65	ng/ul	0.00
46) Acenaphthylene-d8	14.32	160	1578193	37.21	ng/ul	0.00
51) 4-Nitrophenol-d4	14.84	143	199825	32.84	ng/ul	0.00
57) Fluorene-d10	15.61	176	1080641	36.67	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.73	200	157088	31.36	ng/ul	0.00
70) Anthracene-d10	17.45	188	1499052	37.44	ng/ul	0.00
76) Pyrene-d10	19.73	212	1428707	36.73	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.55	264	1393844	38.83	ng/ul	0.00

> S.J
05/04/17

Target Compounds

					Qvalue
6) Phenol	7.21	94	14960	1.16	ng/ul 93
44) Dimethylphthalate	14.07	163	214109	6.48	ng/ul 99

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

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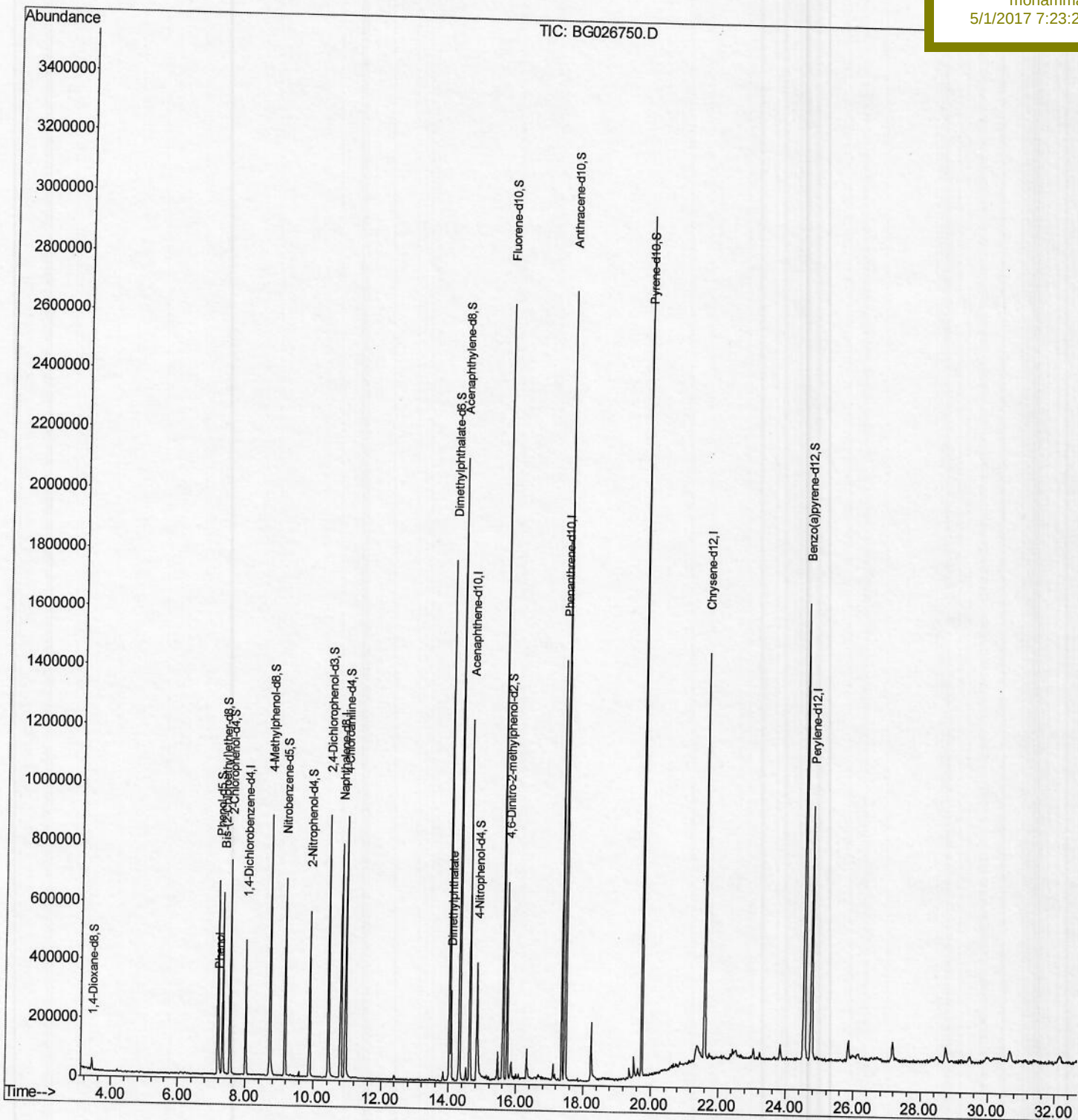
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Manual Integrations

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SOM-EPA-BG042717.M Sat Apr 29 00:58:04 2017

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