

Data Path : Z:\HPCHEM1\BNA_G\Data\BG050117\

Quantitation Report (Qedit)

Data File : BG026833.D

Acq On : 2 May 2017 8:18

Operator : SJ/MA

Sample : I2867-19

Misc :

ALS Vial : 31 Sample Multiplier: 1

Quant Time: May 06 04:40:47 2017

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

Quant Title : SVOA CALIBRATION

QLast Update : Thu May 04 17:40:44 2017

Response via : Initial Calibration

Instrument :

BNA_G

ClientSampleId :

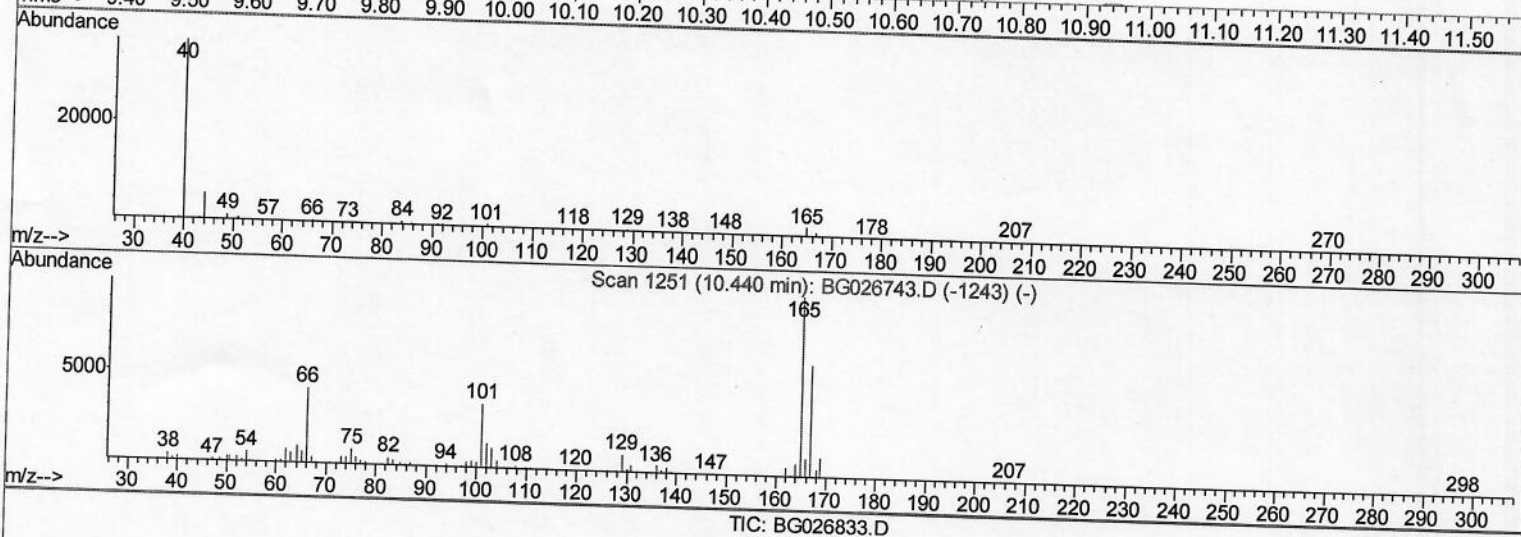
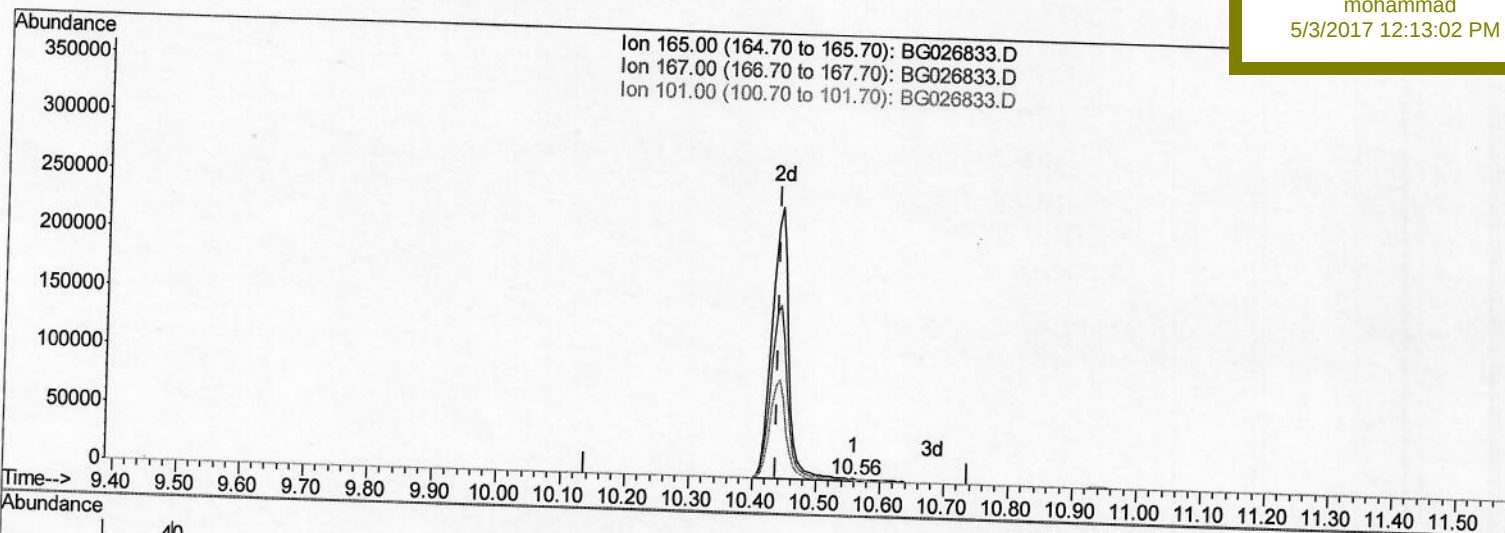
JHTB1

Manual Integrations

APPROVED

mohammad

5/3/2017 12:13:02 PM



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

10.558min (+0.121) 0.08ng/ul

response 1106

Ion	Exp%	Act%
165.00	100	100
167.00	58.70	50.43
101.00	29.80	29.31
0.00	0.00	0.00

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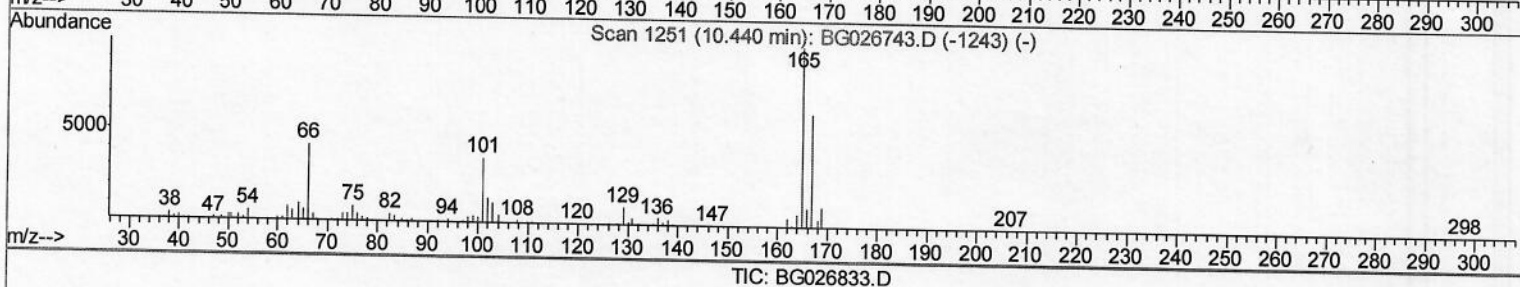
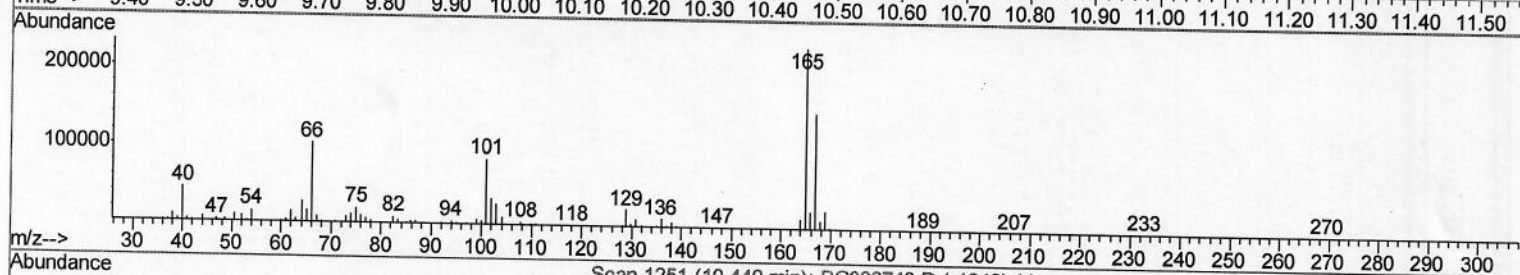
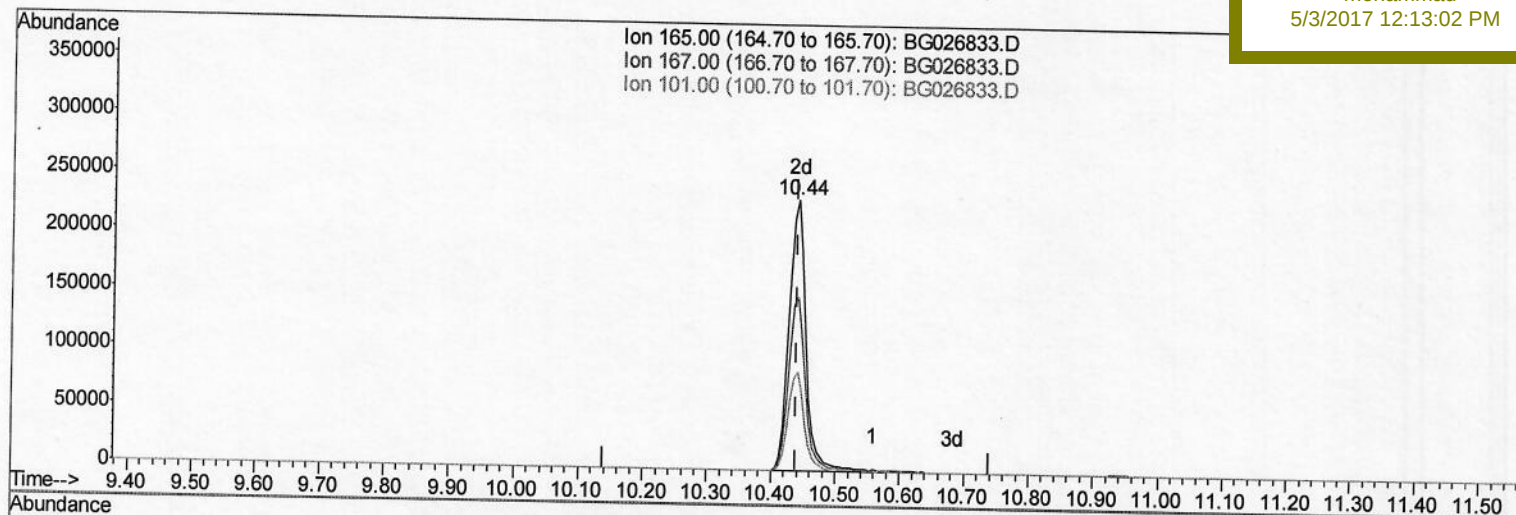
Data File : BG026833.D
Acq On : 2 May 2017 8:18
Operator : SJ/MA
Sample : I2867-19
Misc :
ALS Vial : 31 Sample Multiplier: 1

Quant Time: May 03 00:52:24 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG042717.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 03 00:50:58 2017
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
JHTB1

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

10.441min (+0.001) 33.07ng/ul m

response 434049

Ion Exp% Act%

165.00 100 100

167.00 58.70 64.15

101.00 29.80 36.53#

0.00 0.00 0.00

29.5
05/08/17

Data Path : Z:\HPCHEM1\BNA_G\Data\BG050117\

Quantitation Report

(LSC Reviewed)

Data File : BG026833.D

Acq On : 2 May 2017 8:18

Operator : SJ/MA

Sample : I2867-19

Misc :

ALS Vial : 31 Sample Multiplier: 1

Quant Time: May 03 00:52:24 2017

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

Quant Title : SVOA CALIBRATION

QLast Update : Wed May 03 00:50:58 2017

Response via : Initial Calibration

Instrument :

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	175127	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	908559	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.62	164	525120	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.36	188	975776	20.00	ng/ul	0.00
75) Chrysene-d12	21.60	240	899457	20.00	ng/ul	0.00
83) Perylene-d12	24.77	264	935111	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	22860	5.88	ng/uL	0.00
5) Phenol-d5	7.19	99	557551	35.67	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.33	67	338370	36.81	ng/ul	0.00
9) 2-Chlorophenol-d4	7.54	132	453861	34.21	ng/ul	0.00
13) 4-Methylphenol-d8	8.73	113	470812	36.29	ng/ul	0.00
19) Nitrobenzene-d5	9.17	128	250015	32.92	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	271626	32.88	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.44	165	434049m	33.07	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	627334	35.71	ng/ul	0.00
43) Dimethylphthalate-d6	14.02	166	1370402	32.42	ng/ul	0.00
46) Acenaphthylene-d8	14.32	160	1799009	33.88	ng/ul	0.00
51) 4-Nitrophenol-d4	14.84	143	212119	27.84	ng/ul	0.00
57) Fluorene-d10	15.61	176	1164484	31.57	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.73	200	157172	26.53	ng/ul	0.00
70) Anthracene-d10	17.46	188	1607651	33.95	ng/ul	0.00
76) Pyrene-d10	19.74	212	1593533	33.57	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.55	264	1536781	34.76	ng/ul	0.00

Target Compounds

						Qvalue
6) Phenol	7.22	94	31859	2.01	ng/ul#	86
44) Dimethylphthalate	14.07	163	248419	6.00	ng/ul	98

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

9.3
05/08/17

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Manual Integrations
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