

Quantitation Report (Qedit)

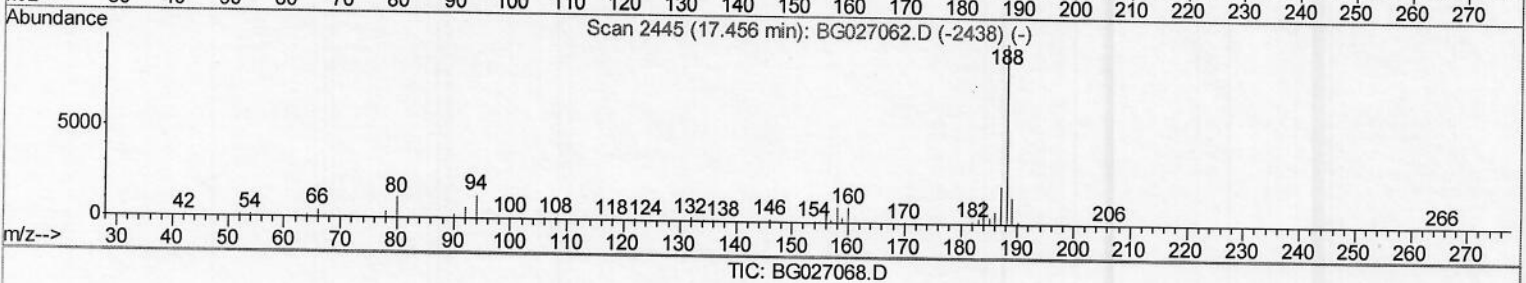
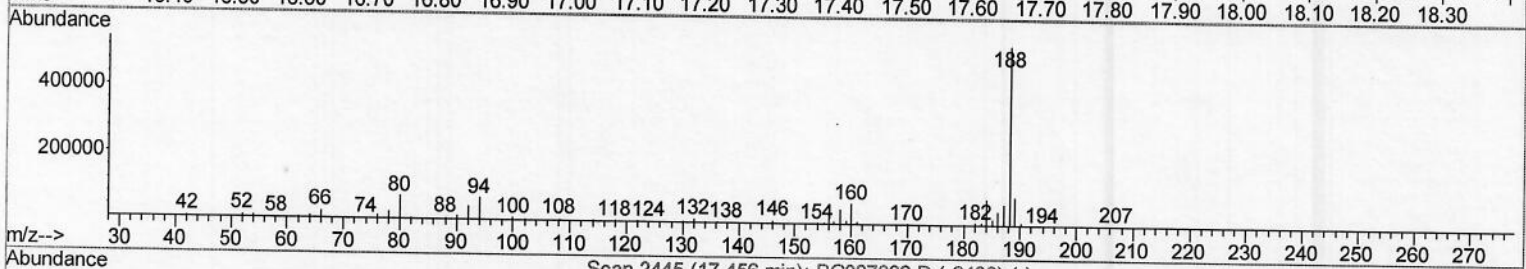
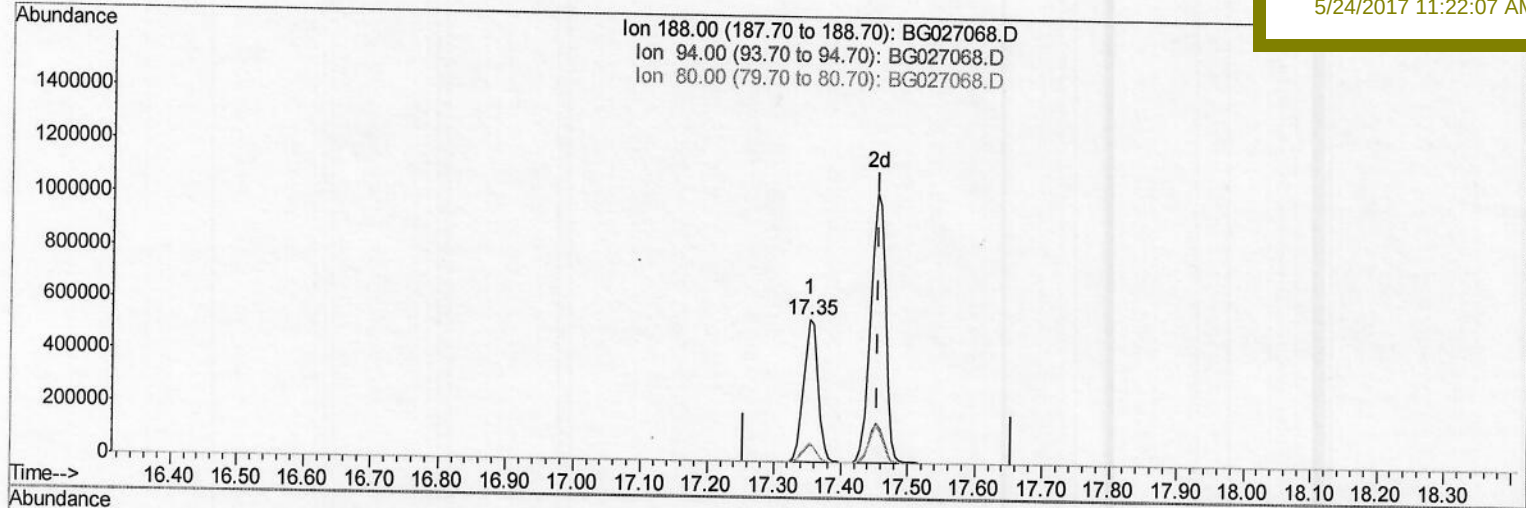
Data Path : Z:\HPCHEM1\BNA G\DATA\BG052317\
 Data File : BG027068.D
 Acq On : 22 May 2017 21:34
 Operator : SJ/MA
 Sample : PB99151BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: May 23 02:45:24 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG051817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 23 02:44:34 2017
 Response via : Initial Calibration

Instrument :
 BNA_G
ClientSampleId :
 SBLK51

Manual Integrations
APPROVED

mohammad
 5/24/2017 11:22:07 AM



(14) Anthracene-d10 (S)
 17.352min (-0.103) 20.51ng/ui
 response 816597

Ion	Exp%	Act%
188.00	100	100
94.00	11.10	12.24
80.00	11.80	12.93
0.00	0.00	0.00

Quantitation Report (Qedit)

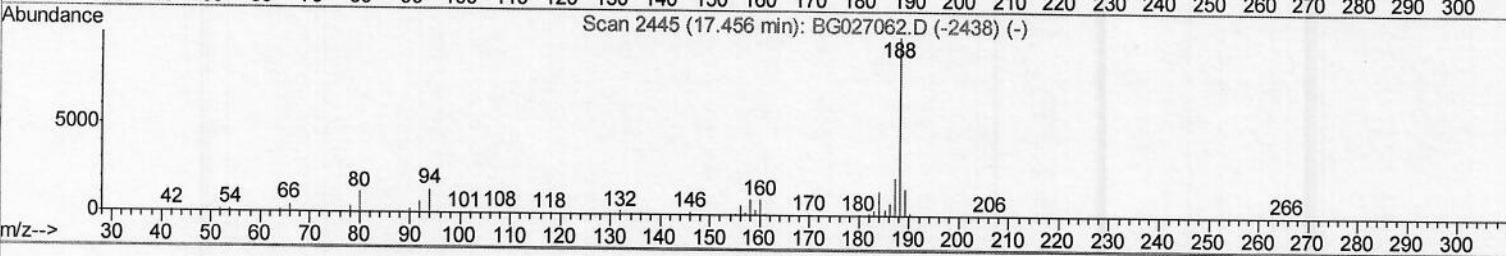
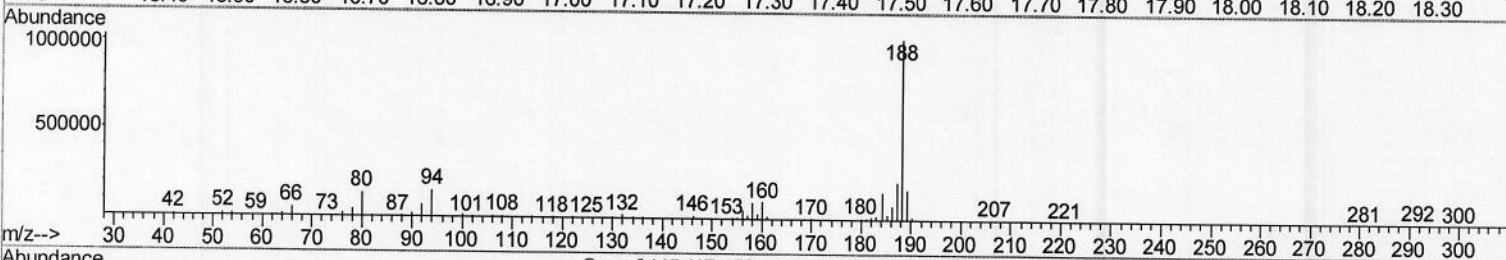
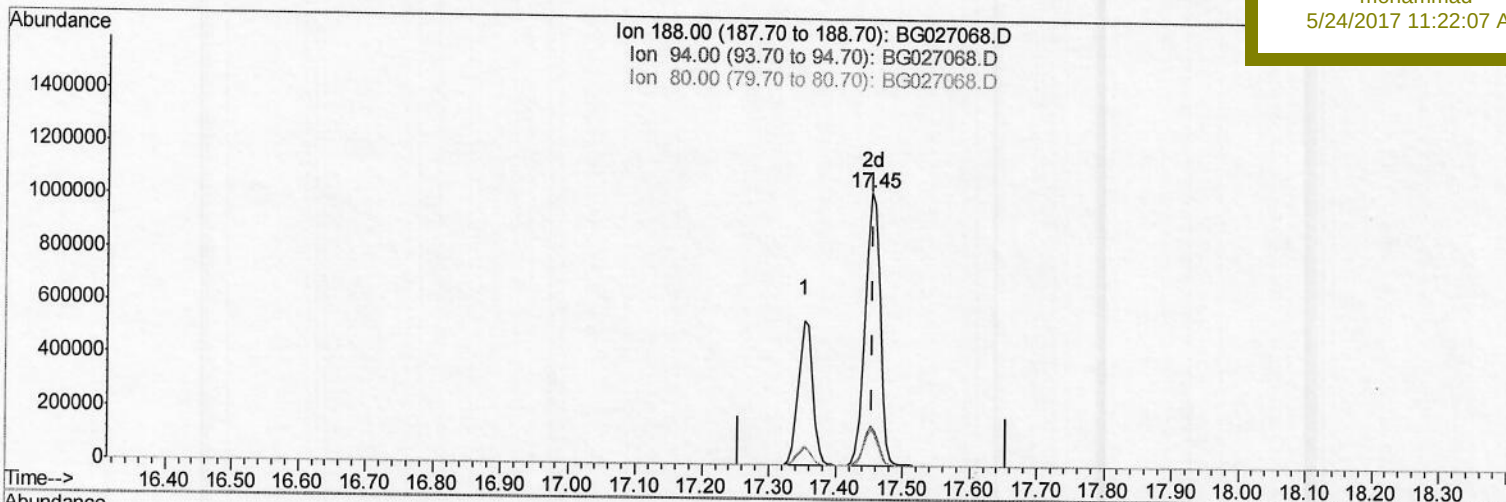
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 BNA_G
 ClientSampled :
 SBLK51

Manual Integrations
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TIC: BG027068.D

(14) Anthracene-d10 (S)

17.452min (-0.003) 40.29ng/ul m>

response 1604165

Ion	Exp%	Act%
188.00	100	100
94.00	11.10	14.69#
80.00	11.80	13.23
0.00	0.00	0.00

S.J
 05/30/17

Data Path : Z:\HPCHEM1\BNA G\DATA\BG052317\
 Data File : BG027068.D
 Acq On : 22 May 2017 21:34
 Operator : SJ/MA
 Sample : PB99151BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: May 23 03:01:25 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG051817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 23 02:44:34 2017
 Response via : Initial Calibration

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	176719	20.00	ng/ul	0.00
2) Naphthalene-d8	10.81	136	850877	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.62	164	440216	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.35	188	816597	20.00	ng/ul	0.00
17) Chrysene-d12	21.60	240	736210	20.00	ng/ul	0.00
22) Perylene-d12	24.77	264	625559	20.00	ng/ul	0.00

System Monitoring Compounds

7) Acenaphthylene-d8	14.31	160	1787544	39.59	ng/ul	0.00
10) Fluorene-d10	15.61	176	1139955	37.84	ng/ul	0.00
14) Anthracene-d10	17.45	188	1604165m	40.29	ng/ul	0.00
18) Pvrene-d10	19.74	212	1535209	39.94	ng/ul	0.00
25) Benzo(a)pyrene-d12	24.56	264	1354182	43.57	ng/ul	0.00

Target Compounds

Ovalue

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

> 9.7
 05/30/17

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Manual Integrations
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