

Quantitation Report (Qedit)

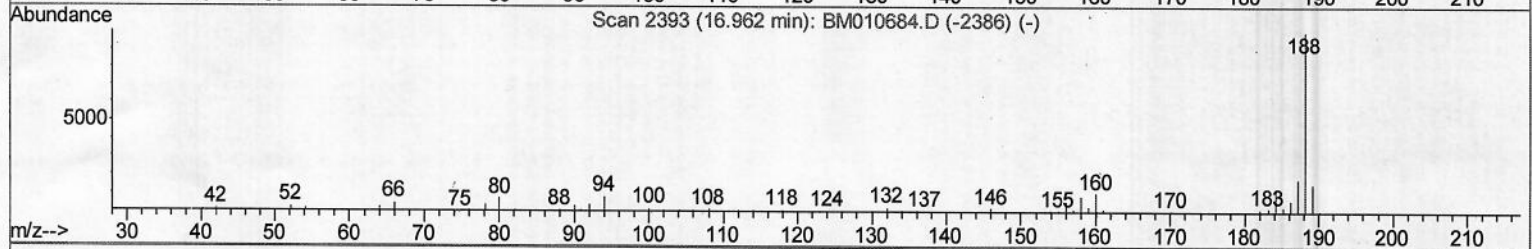
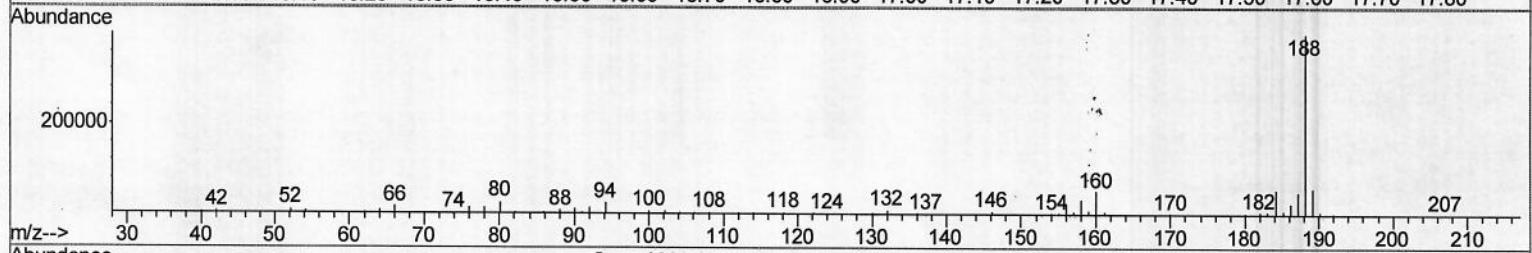
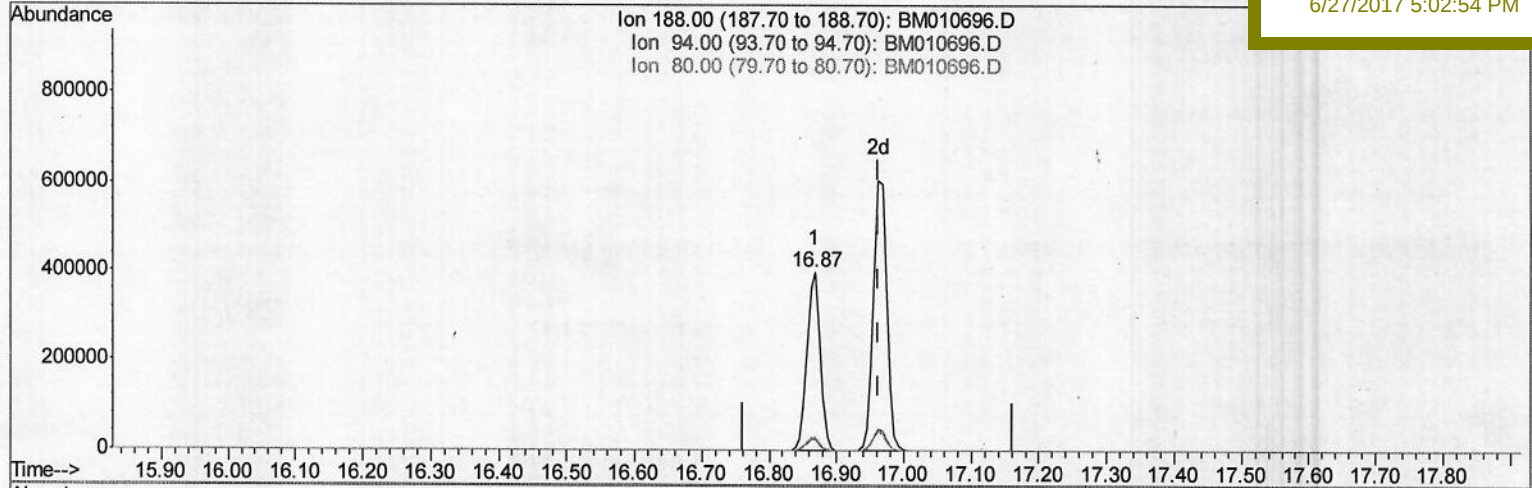
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
 Data File : BM010696.D
 Acq On : 23 Jun 2017 18:50
 Operator : SJ/JU
 Sample : PB100052BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jun 24 03:39:17 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jun 24 03:35:21 2017
 Response via : Initial Calibration

Instrument :
 BNA_M
ClientSampleId :
 SBLK52

Manual Integrations
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TIC: BM010696.D

(70) Anthracene-d10 (S)

16.868min (-0.094) 20.71ng/ul

response 539036

Ion	Exp%	Act%
188.00	100	100
94.00	5.70	6.18
80.00	5.90	6.58
0.00	0.00	0.00

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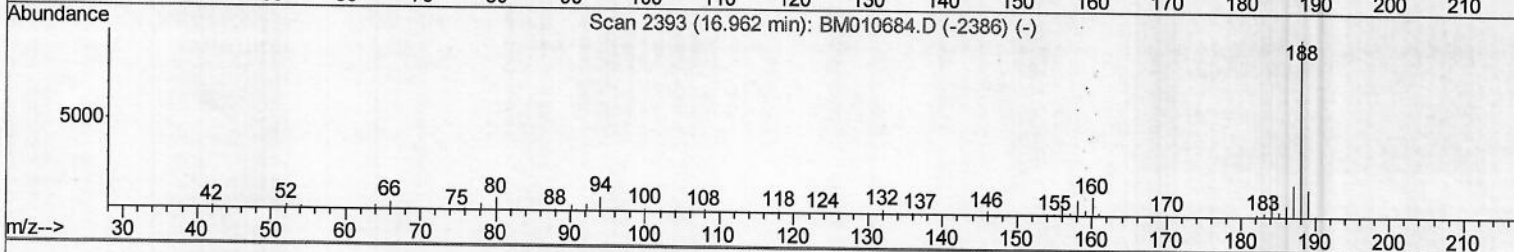
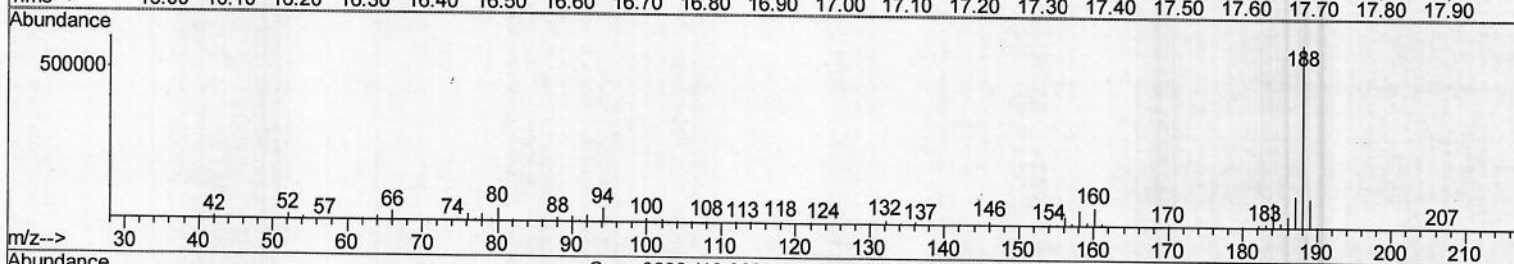
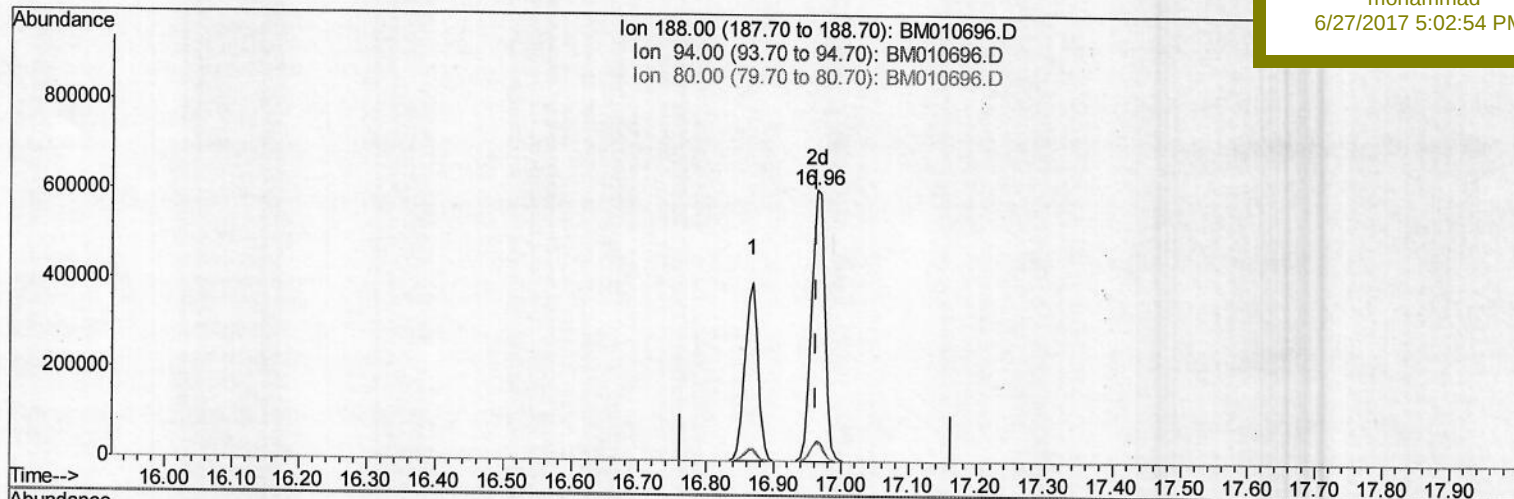
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TIC: BM010696.D

(70) Anthracene-d10 (S)

16.962min (+0.000) 32.15ng/ul m

response 836994

Ion	Exp%	Act%
188.00	100	100
94.00	5.70	7.57#
80.00	5.90	6.95
0.00	0.00	0.00

SJ
 06/24/17

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Quant Time: Jun 24 04:33:11 2017
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	64236	20.00	ng/ul	0.00
18) Naphthalene-d8	10.22	136	302082	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.11	164	216480	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.87	188	539036	20.00	ng/ul	0.00
75) Chrysene-d12	21.08	240	571967	20.00	ng/ul	0.00
83) Perylene-d12	23.22	264	403665	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.07	96	7864	7.01	ng/uL	0.00
5) Phenol-d5	6.65	99	160246	26.19	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.81	67	94138	26.95	ng/ul	0.00
9) 2-Chlorophenol-d4	7.00	132	124953	29.11	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	141745	27.94	ng/ul	0.00
19) Nitrobenzene-d5	8.60	128	64611	29.98	ng/ul	0.00
22) 2-Nitrophenol-d4	9.32	143	82530	33.24	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.85	165	151428	28.97	ng/ul	0.00
29) 4-Chloroaniline-d4	10.36	131	191771	34.82	ng/ul	0.00
43) Dimethylphthalate-d6	13.53	166	596831	31.41	ng/ul	0.00
46) Acenaphthylene-d8	13.80	160	663346	30.01	ng/ul	0.00
51) 4-Nitrophenol-d4	14.33	143	91500	27.36	ng/ul	0.00
57) Fluorene-d10	15.12	176	492816	30.00	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.24	200	98678	29.81	ng/ul	0.00
70) Anthracene-d10	16.96	188	836994m	32.15	ng/ul	0.00
76) Pyrene-d10	19.27	212	964889	36.31	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.09	264	624888	32.03	ng/ul	0.00

→ S.J
 06/24/17

Target Compounds

Qvalue

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

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