

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062017\
Quantitation Report (QT Reviewed)

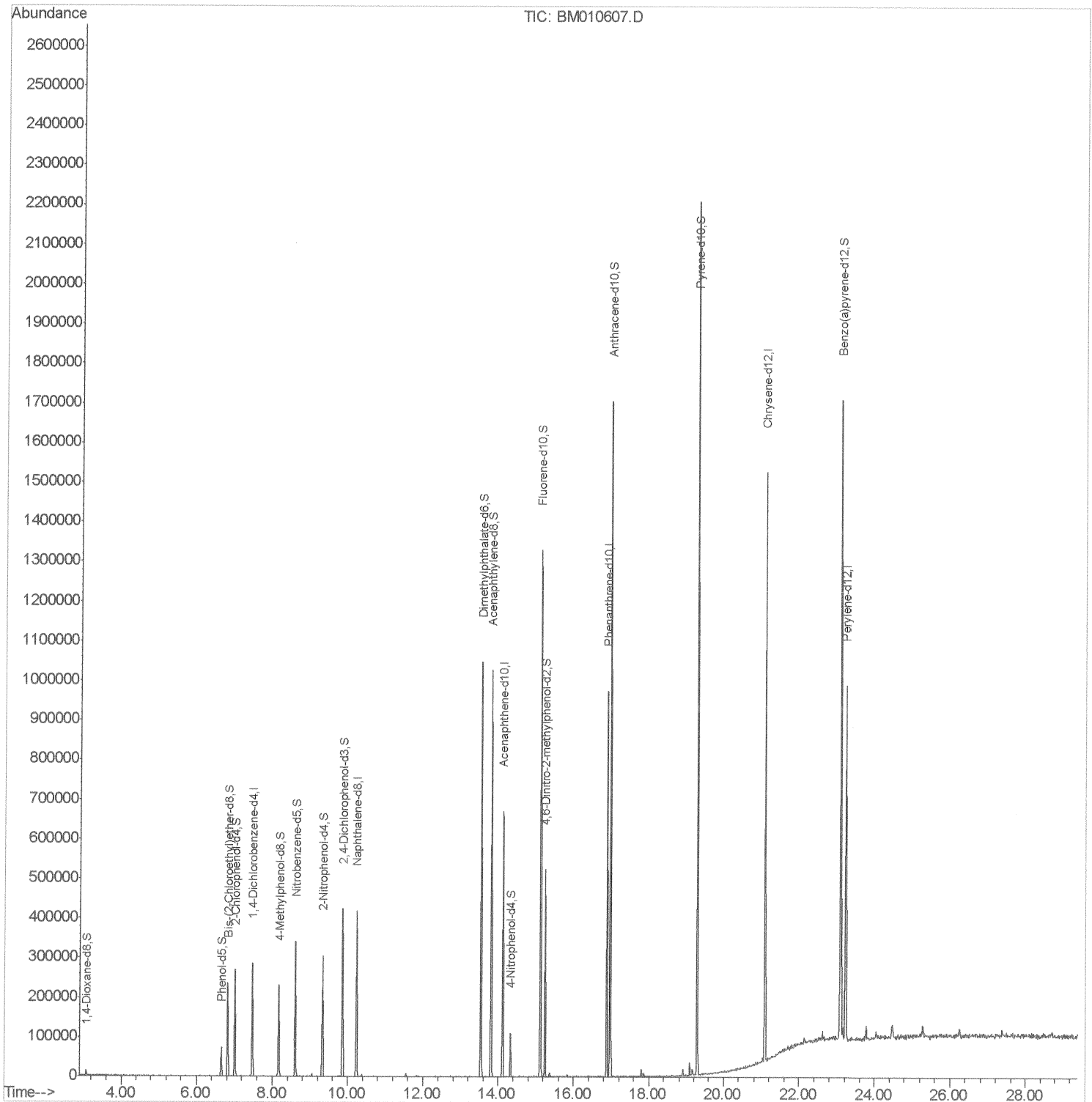
Data File : BM010607.D
Acq On : 21 Jun 2017 01:30
Operator : SJ/MA
Sample : I3790-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B77

Manual Integrations
APPROVED

mohammad
6/21/2017 4:09:07 PM

Quant Time: Jun 21 03:45:29 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 21 01:33:50 2017
Response via : Initial Calibration



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062017\

Quantitation Report (Qedit)

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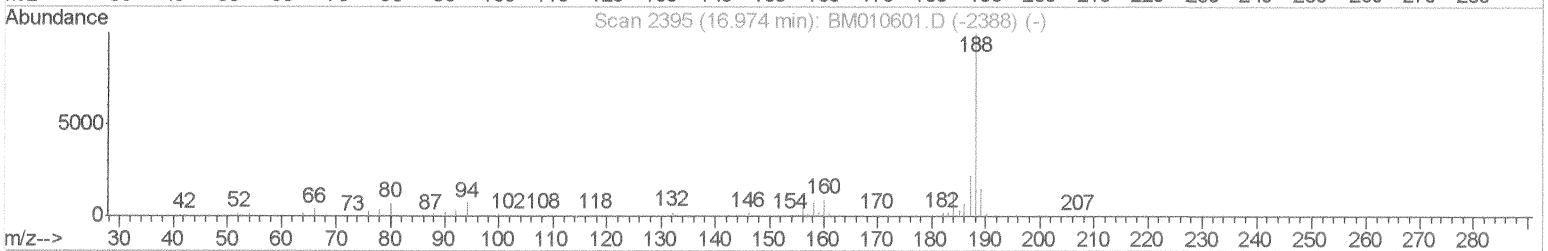
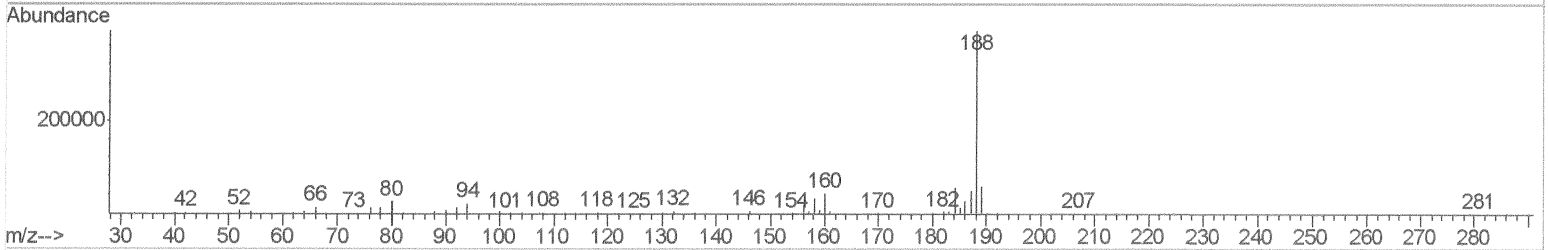
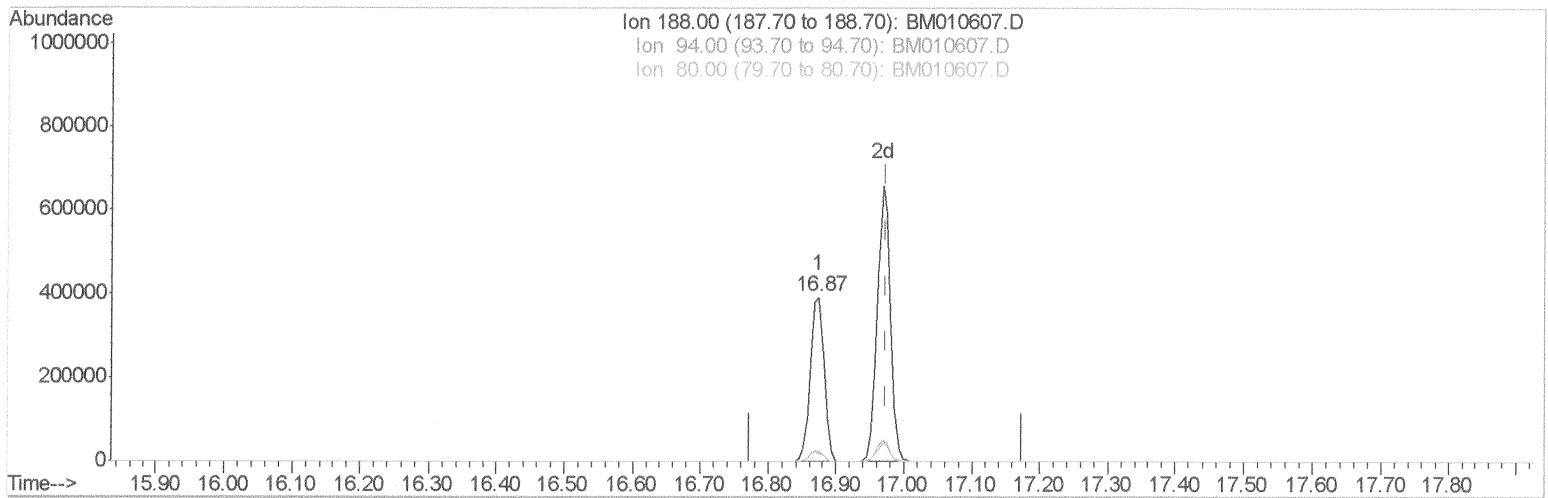
Quant Time: Jun 21 03:44:00 2017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Jun 21 01:33:50 2017

Response via : Initial Calibration



TIC: BM010607.D

(70) Anthracene-d10 (S)

16.874min (-0.100) 20.71ng/ui

response 543752

Ion	Exp%	Act%
188.00	100	100
94.00	5.70	5.73
80.00	5.90	6.90
0.00	0.00	0.00

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Manual Integrations
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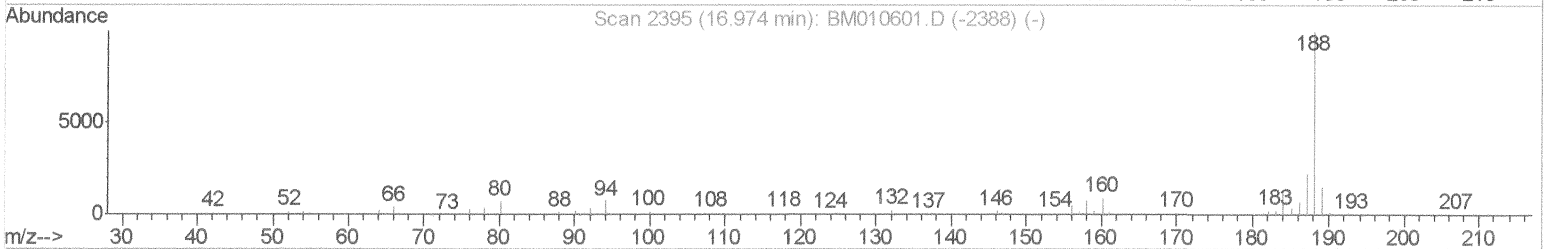
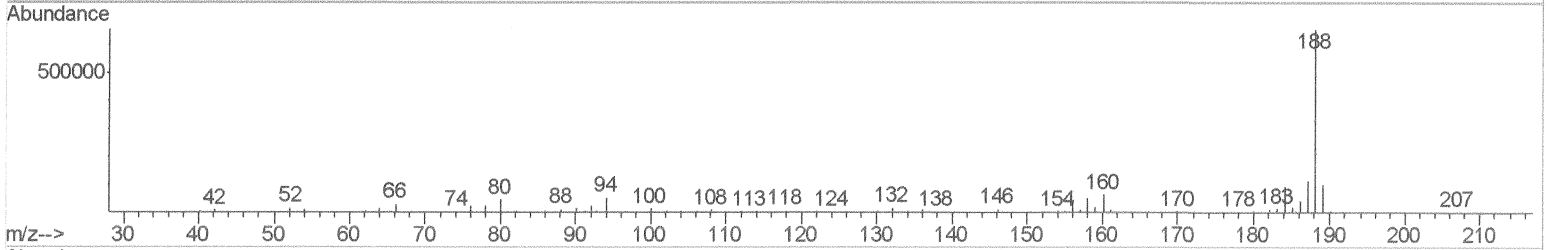
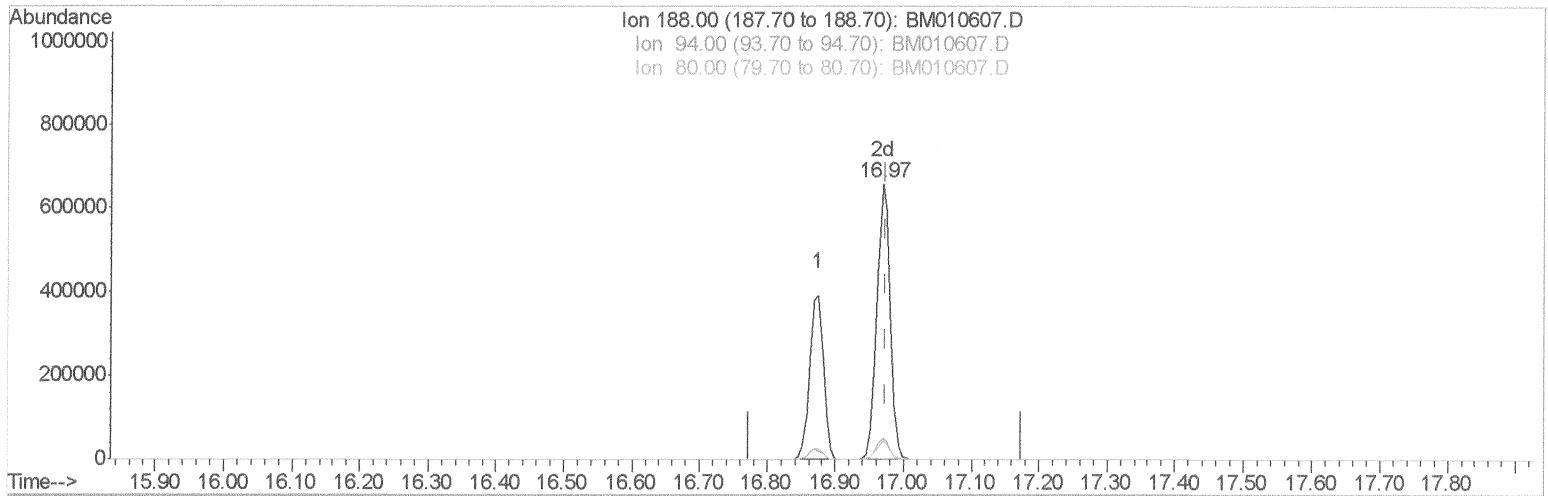
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Response via : Initial Calibration



TIC: BM010607.D

(70) Anthracene-d10 (S)

16.968min (-0.006) 33.88ng/ul m

response 889596

Ion	Exp%	Act%
188.00	100	100
94.00	5.70	7.77#
80.00	5.90	6.80
0.00	0.00	0.00

JS
06/26/16

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062017\
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Acq On : 21 Jun 2017 01:30
Operator : SJ/MA
Sample : I3790-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B77

Manual Integrations
APPROVED

mohammad
6/21/2017 4:09:07 PM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	66494	20.00	ng/ul	0.00
18) Naphthalene-d8	10.23	136	303101	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.12	164	205700	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.87	188	543752	20.00	ng/ul	0.00
75) Chrysene-d12	21.09	240	671982	20.00	ng/ul	0.00
83) Perylene-d12	23.23	264	609159	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	7117	6.13	ng/uL	0.00
5) Phenol-d5	6.65	99	35788	5.65	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.82	67	92693	25.63	ng/ul	0.00
9) 2-Chlorophenol-d4	7.00	132	97519	21.95	ng/ul	0.00
13) 4-Methylphenol-d8	8.17	113	69026	13.14	ng/ul	0.00
19) Nitrobenzene-d5	8.61	128	63231	29.25	ng/ul	0.00
22) 2-Nitrophenol-d4	9.32	143	75839	30.45	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.86	165	135142	25.77	ng/ul	0.00
29) 4-Chloroaniline-d4	10.37	131	3662	0.66	ng/ul	0.00
43) Dimethylphthalate-d6	13.54	166	587241	32.52	ng/ul	0.00
46) Acenaphthylene-d8	13.80	160	641001	30.52	ng/ul	0.00
51) 4-Nitrophenol-d4	14.32	143	19067	6.00	ng/ul	0.00
57) Fluorene-d10	15.12	176	497246	31.86	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.24	200	104911	31.41	ng/ul	0.00
70) Anthracene-d10	16.97	188	889596m	33.88	ng/ul	0.00
76) Pyrene-d10	19.28	212	1073311	34.38	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.10	264	983357	33.40	ng/ul	0.00

Target Compounds Qvalue

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

JS 06/26/17