

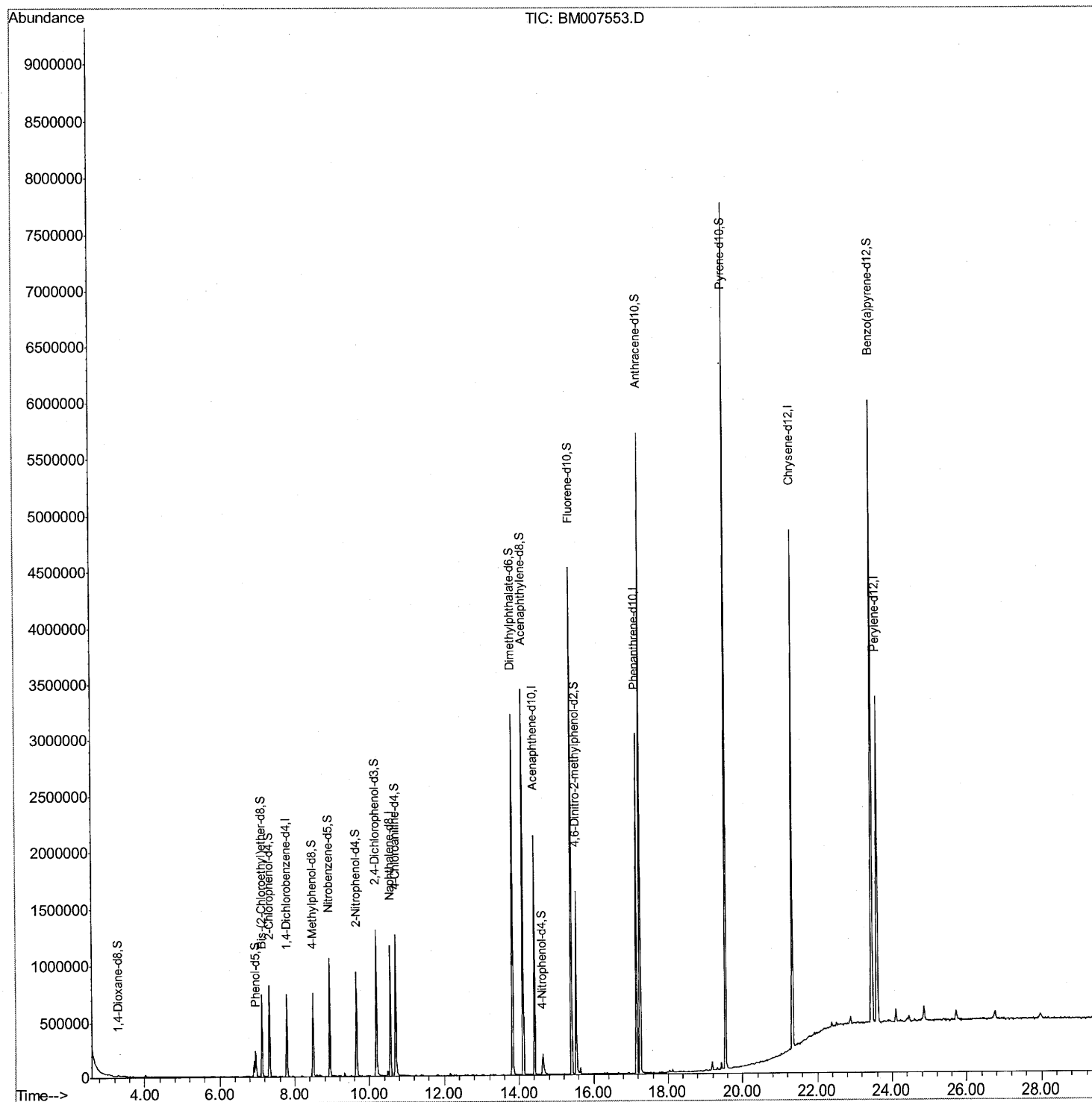
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092316\
Data File : BM007553.D
Acq On : 24 Sep 2016 09:00
Operator : UM/SJ
Sample : H4866-03
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
E43S4

Manual Integrations
APPROVED

sohil
9/26/2016 6:35:42 PM

Quant Time: Sep 26 03:20:14 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM092316.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 24 04:39:16 2016
Response via : Initial Calibration



Quantitation Report (Qedit)

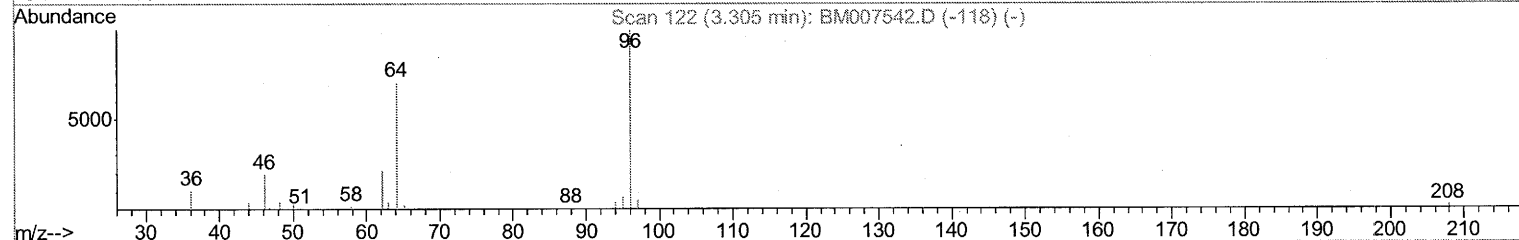
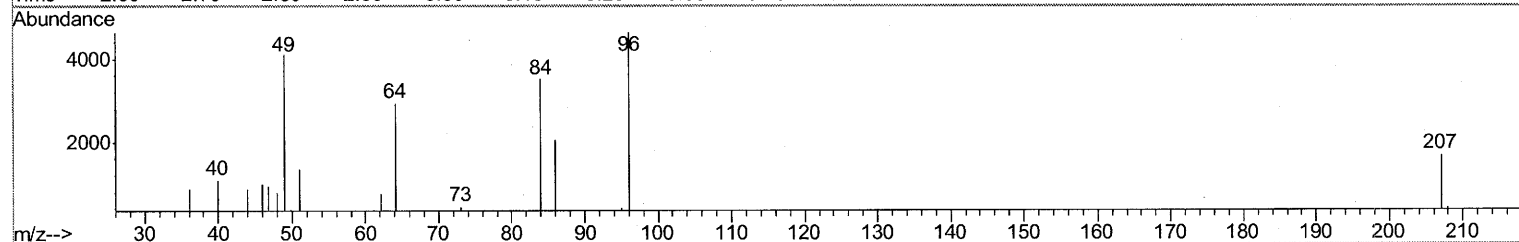
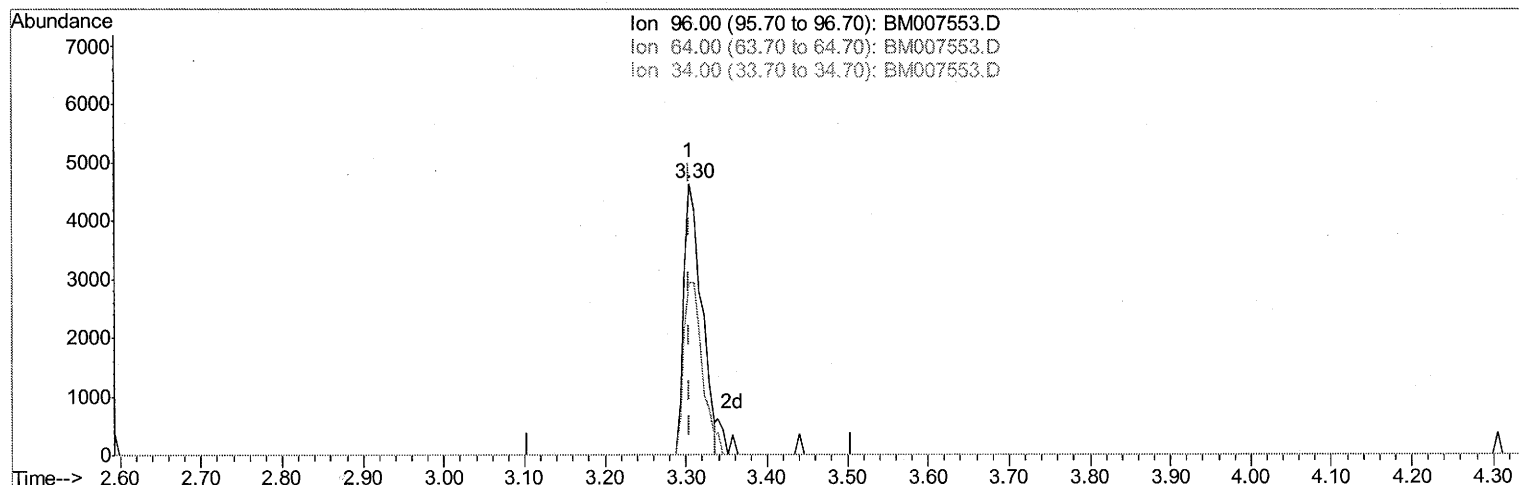
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Quant Time: Sep 26 03:09:05 2016
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 Response via : Initial Calibration



TIC: BM007553.D

(3) 1,4-Dioxane-d8 (S)

3.305min (+0.000) 1.71ng/uL

response 6967

Ion	Exp%	Act%
96.00	100	100
64.00	68.70	67.35
34.00	0.00	0.00
0.00	0.00	0.00

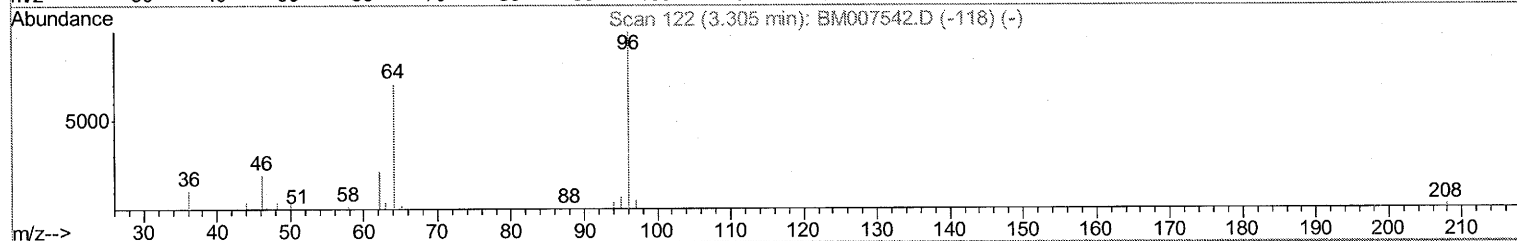
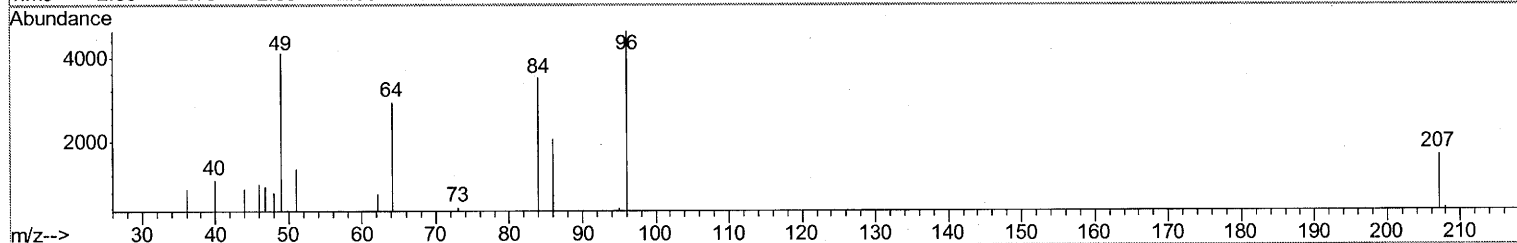
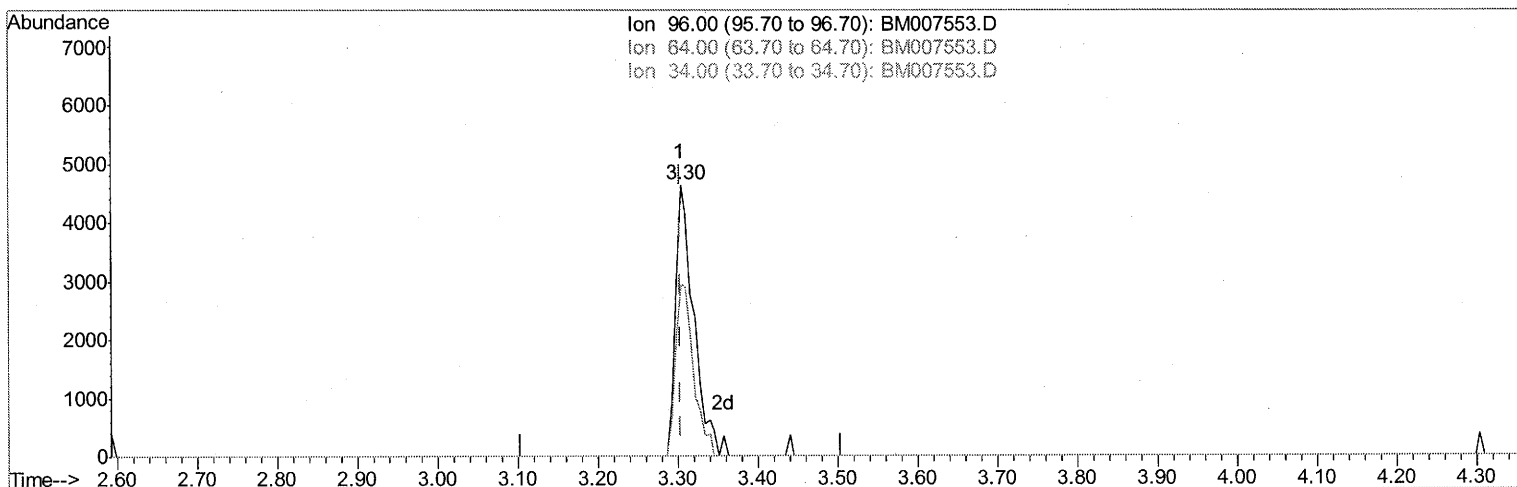
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Manual Integrations
APPROVED

sohil
9/26/2016 6:35:42 PM

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BM007553.D

(3) 1,4-Dioxane-d8 (S)

3.305min (+0.000) 1.80ng/uL m

response 7342

Ion	Exp%	Act%
96.00	100	100
64.00	68.70	63.91
34.00	0.00	0.00
0.00	0.00	0.00

UM
09/26/16

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092316\
 Data File : BM007553.D
 Acq On : 24 Sep 2016 09:00
 Operator : UM/SJ
 Sample : H4866-03
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 E43S4

Manual Integrations
 APPROVED

sohil
 9/26/2016 6:35:42 PM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	200522	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	929008	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.41	164	747613	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	1732918	20.00	ng/ul	0.00
75) Chrysene-d12	21.34	240	2377107	20.00	ng/ul	0.00
83) Perylene-d12	23.60	264	2257907	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.30	96	7342m	1.80	ng/uL	0.00
5) Phenol-d5	6.95	99	141991	7.98	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	333028	34.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	360959	27.93	ng/ul	0.00
13) 4-Methylphenol-d8	8.48	113	283619	18.67	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	244417	36.80	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	289969	38.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	499515	33.23	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	672519	42.61	ng/ul	0.00
43) Dimethylphthalate-d6	13.83	166	2158802	39.70	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	2377020	37.67	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	75403	7.86	ng/ul	0.00
57) Fluorene-d10	15.40	176	1852654	38.85	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	373890	35.36	ng/ul	0.00
70) Anthracene-d10	17.26	188	3207537	40.28	ng/ul	0.00
76) Pyrene-d10	19.55	212	4042483	42.85	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.46	264	4135672	41.40	ng/ul	0.00

UM
 09/26/16

Target Compounds Qvalue

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