

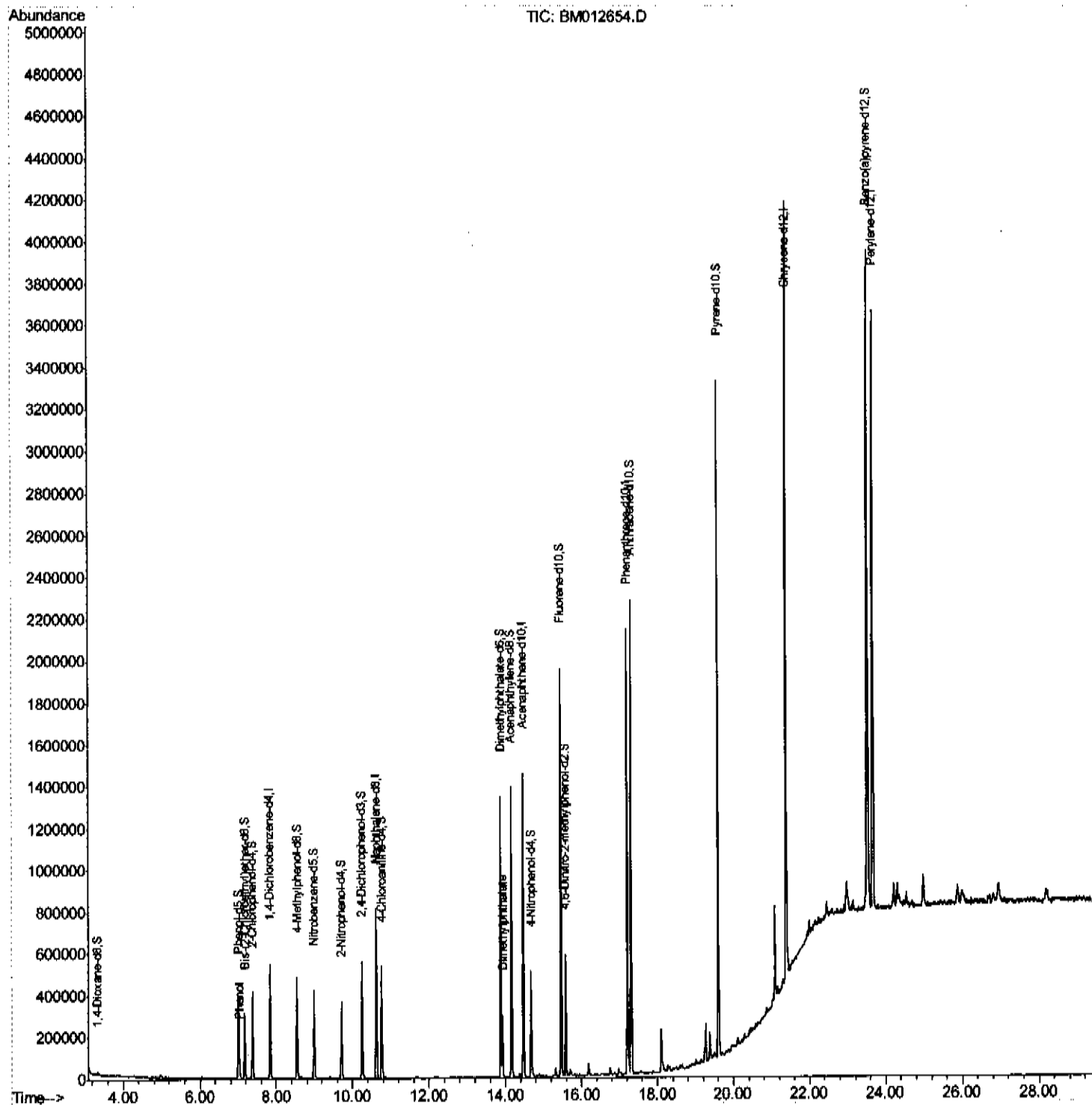
Data Path : Z:\HPCHEM1\BNA M\DATA\BM110117\
Data File : BM012654.D
Acq On : 02 Nov 2017 01:34
Operator : SJ/JU
Sample : I5937-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
B0DT0

Manual Integrations
APPROVED

Sohil
11/2/2017 5:22:29 PM

Quant Time: Nov 02 04:33:34 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 02 04:04:31 2017
Response via : Initial Calibration



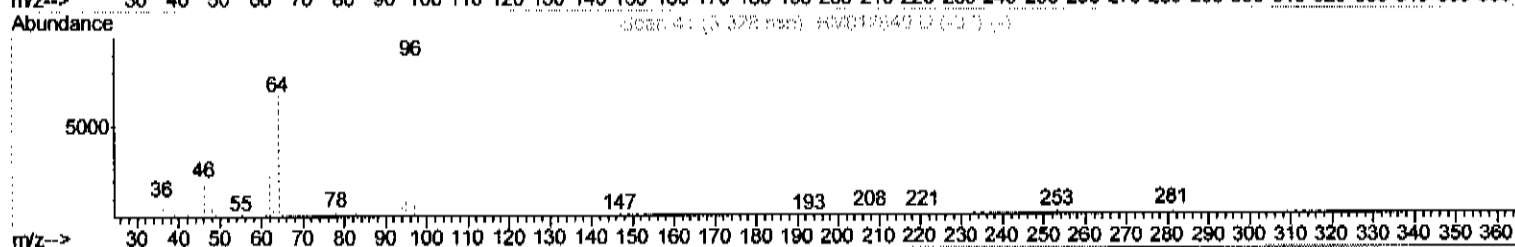
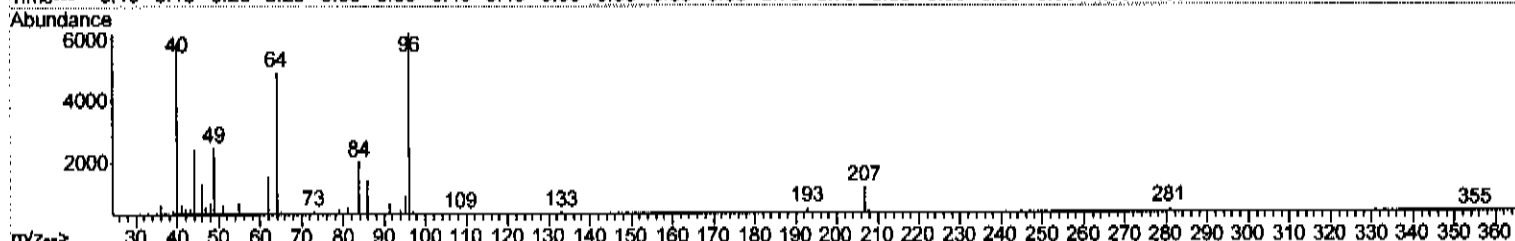
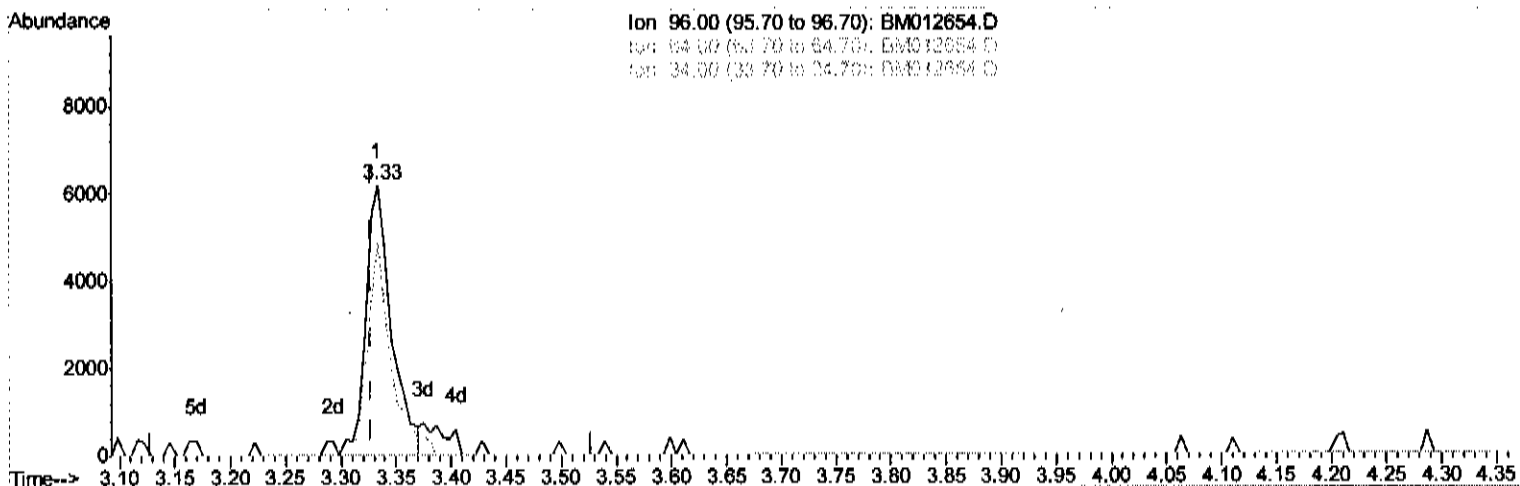
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TIC: BM012654.D

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

3.334min (+0.006) 3.45ng/uL

response 9881

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	70.23
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

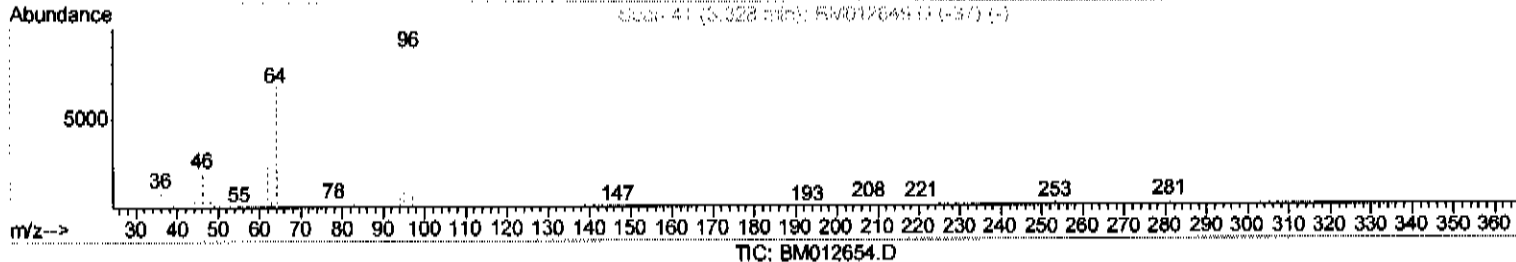
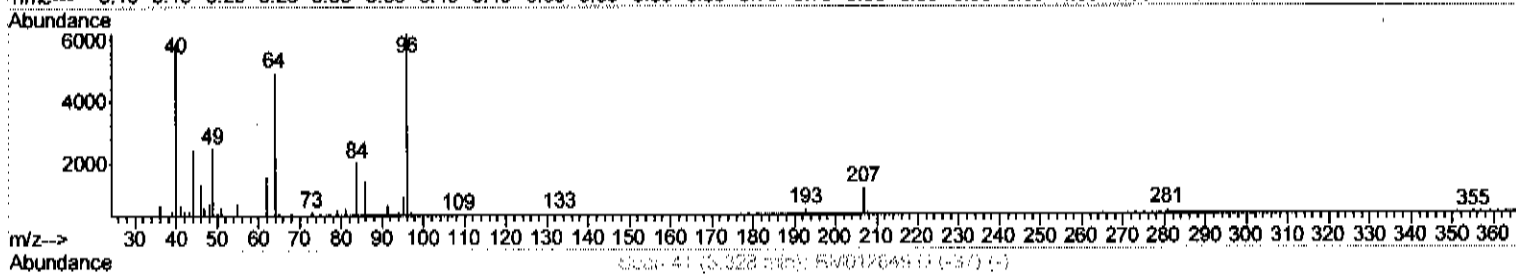
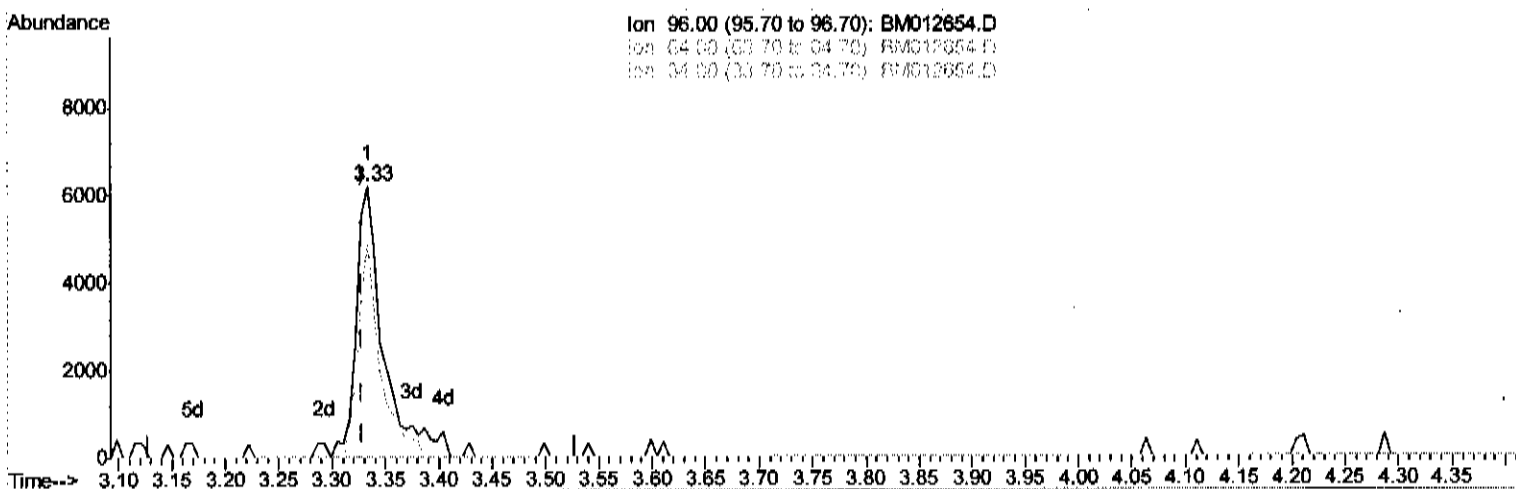
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BNA_M
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B0DT0

Manual Integrations
APPROVED

Sohil
11/2/2017 5:22:29 PM

Quant Time: Nov 02 04:16:51 2017
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Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 02 04:04:31 2017
Response via : Initial Calibration



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

3.334min (+0.006) 3.86ng/uL m

response 11033

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	62.89
34.00	0.00	0.00
0.00	0.00	0.00

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 Operator : SJ/JU
 Sample : I5937-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B0DT0

Manual Integrations
 APPROVED

Sohil
 11/2/2017 5:22:29 PM

Quant Time: Nov 02 04:33:34 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 02 04:04:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	146084	20.00	ng/ul	0.00
18) Naphthalene-d8	10.63	136	670234	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.47	164	476886	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.22	188	1249725	20.00	ng/ul	0.00
75) Chrysene-d12	21.39	240	1810729	20.00	ng/ul	0.00
83) Perylene-d12	23.69	264	2279334	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.33	96	11033m	3.86	ng/uL	0.00
5) Phenol-d5	7.01	99	224450	18.39	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.17	67	137771	19.02	ng/ul	0.00
9) 2-Chlorophenol-d4	7.37	132	185615	20.31	ng/ul	0.00
13) 4-Methylphenol-d8	8.55	113	184544	17.69	ng/ul	0.00
19) Nitrobenzene-d5	8.99	128	101098	24.08	ng/ul	0.00
22) 2-Nitrophenol-d4	9.72	143	121625	28.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.26	165	219265	19.14	ng/ul	0.00
29) 4-Chloroaniline-d4	10.77	131	284052	23.52	ng/ul	0.00
43) Dimethylphthalate-d6	13.88	166	878679	21.09	ng/ul	0.00
46) Acenaphthylene-d8	14.16	160	986963	20.03	ng/ul	0.00
51) 4-Nitrophenol-d4	14.67	143	132795	21.44	ng/ul	0.00
57) Fluorene-d10	15.46	176	766222	21.73	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.59	200	145745	24.10	ng/ul	0.00
70) Anthracene-d10	17.32	188	1292928	21.71	ng/ul	0.00
76) Pyrene-d10	19.60	212	1640505	19.75	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.54	264	2364662	21.84	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.04	94	16063	1.26 ng/ul#	82
44) Dimethylphthalate	13.93	163	189104	4.59 ng/ul	98

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