

Quantitation Report (QT Reviewed)

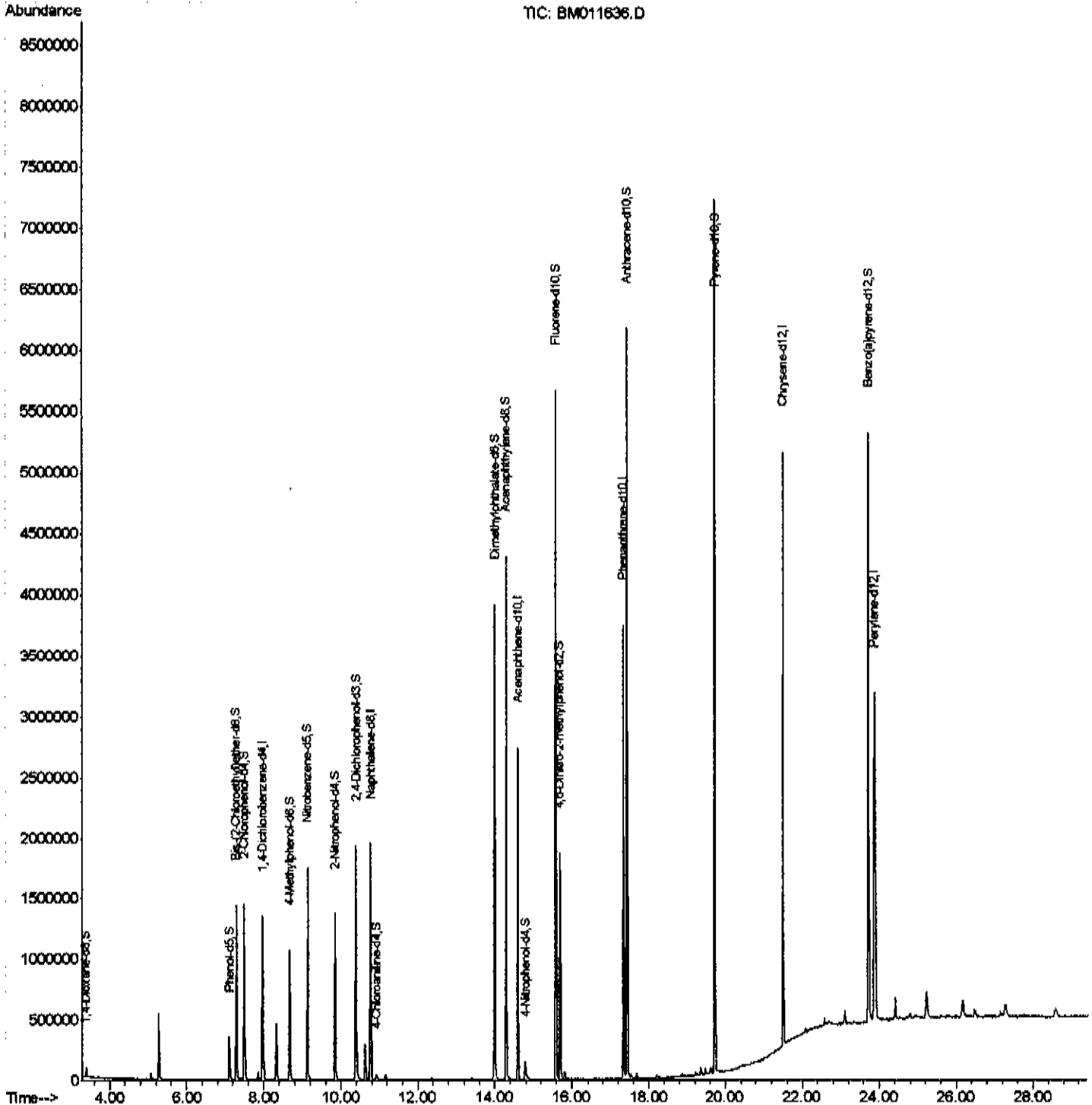
Data Path : Z:\HPCHEM1\BNA M\DATA\BM092017\
 Data File : BM011636.D
 Acq On : 20 Sep 2017 14:19
 Operator : SJ/JU
 Sample : I5273-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 COAL3

Manual Integrations
 APPROVED

Sohil
 9/22/2017 4:05:14 PM

Quant Time: Sep 21 02:42:53 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 OLast Update : Thu Sep 21 01:44:48 2017
 Response via : Initial Calibration



Instrument :
BNA_M
ClientSampleId :
C0AL3

Manual Integrations

Sohil
9/22/2017 4:05:14 PM

Abundance

Ion 131.00 (130.70 to 131.70): BM011636.D
File: BM011636.D, Date: 10-01-2011, Time: 15:35:11
Scan range: 1000000 to 1000000, Scan rate: 1000000

20000

15000

10000

5000

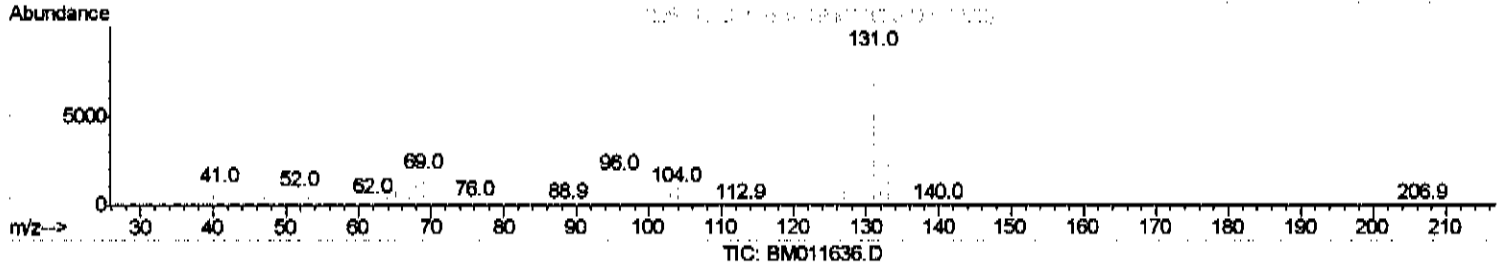
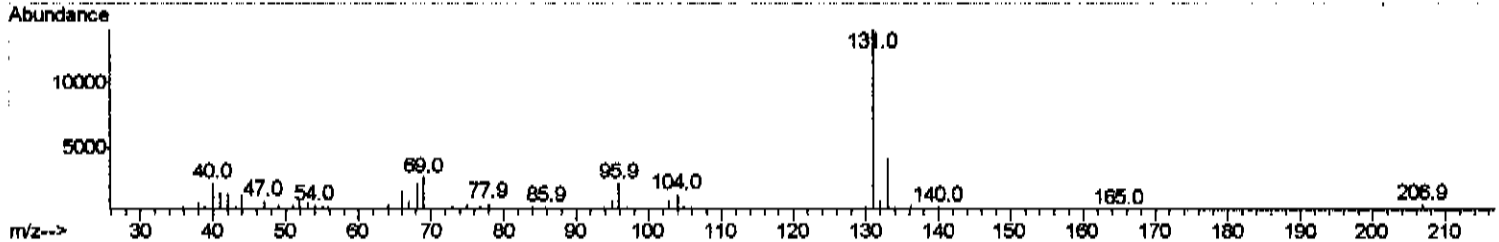
0

Time--> 9.90 10.00 10.10 10.20 10.30 10.40 10.50 10.60 10.70 10.80 10.90 11.00 11.10 11.20 11.30 11.40 11.50 11.60 11.70 11.80 11.90

1
10.92

3d

2d



гепропна 33343

Ion	Exp%	Act%
131.00	100	100
133.00	33.90	30.60
69.00	18.60	19.67
0.00	0.00	0.00

Quantitation Report (Qedit)

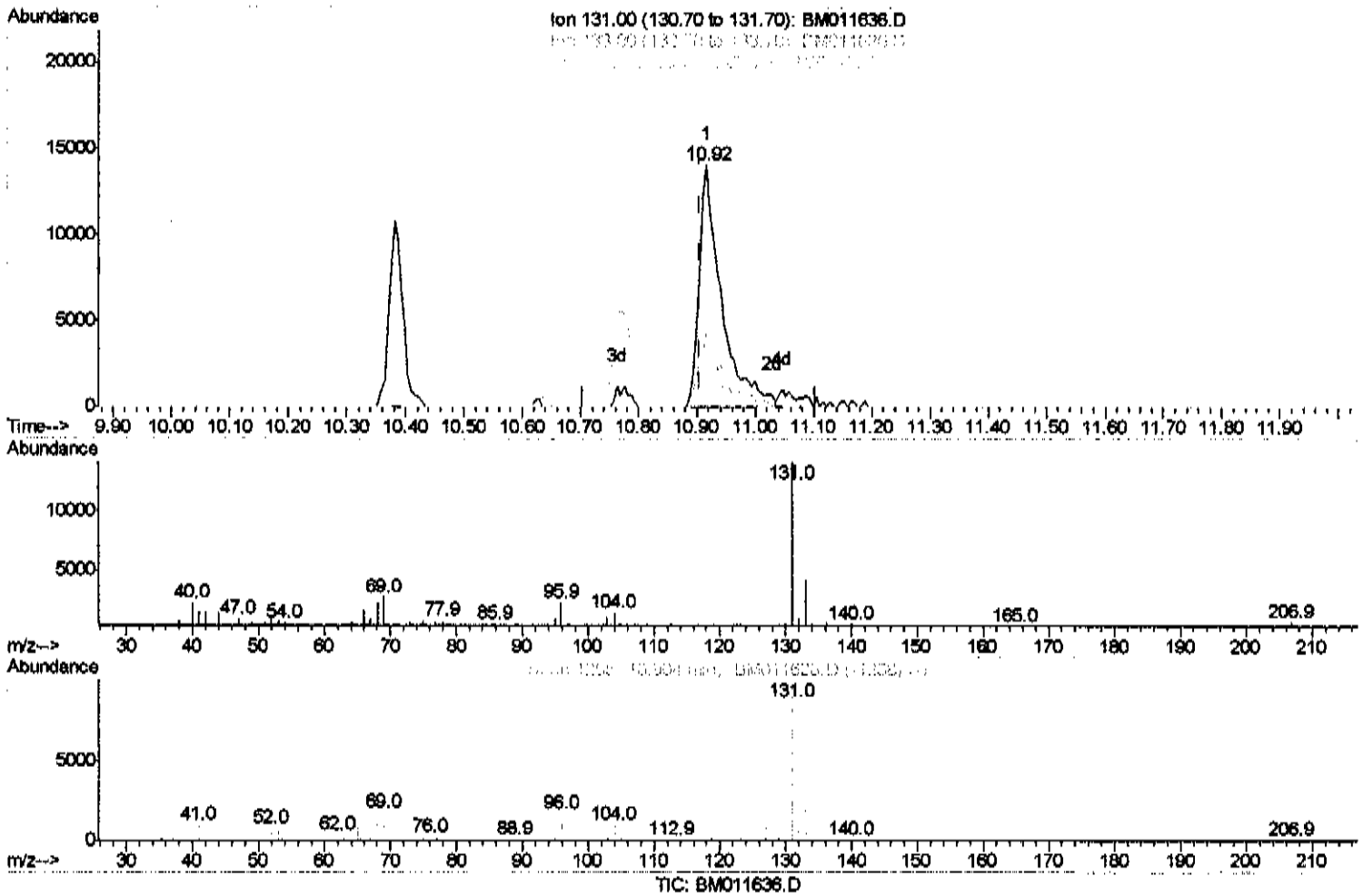
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Manual Integrations
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(29) 4-Chloroaniline-d4 (S)

10.916min (+0.012) 1.34ng/ul m

response 36893

Ion	Exp%	Act%
131.00	100	100
133.00	33.90	30.60
69.00	18.60	19.67
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Instrument :
BNA_M
ClientSampled :
C0AL3

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Manual Integrations
APPROVED

Sohil
9/22/2017 4:05:14 PM

Quant Time: Sep 21 02:42:53 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 21 01:44:48 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	342786	20.00	ng/ul	0.00
18) Naphthalene-d8	10.77	136	1565870	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.59	164	916138	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.33	188	2055011	20.00	ng/ul	0.00
78) Chrysene-d12	21.50	240	2184187	20.00	ng/ul	0.00
86) Perylene-d12	23.87	264	2015302	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.39	96	36519	4.79	ng/uL	0.00
5) Phenol-d5	7.11	99	202497	6.15	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.29	67	701221	31.61	ng/ul	0.00
9) 2-Chlorophenol-d4	7.49	132	598057	24.11	ng/ul	0.00
13) 4-Methylphenol-d8	8.66	113	382131	14.87	ng/ul	0.00
19) Nitrobenzene-d5	9.13	128	373791	31.45	ng/ul	0.00
22) 2-Nitrophenol-d4	9.85	143	396316	32.08	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.38	165	689839	27.73	ng/ul	0.00
29) 4-Chloroaniline-d4	10.92	131	36893m	1.34	ng/ul	0.01
44) Dimethylphthalate-d6	14.00	166	2439777	31.50	ng/ul	0.00
47) Acenaphthylene-d8	14.29	160	2952257	31.52	ng/ul	0.00
52) 4-Nitrophenol-d4	14.79	143	45373	3.12	ng/ul	0.00
58) Fluorene-d10	15.59	176	2139764	31.90	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.70	200	312065	26.10	ng/ul	0.00
71) Anthracene-d10	17.43	188	3352412	33.13	ng/ul	0.00
79) Pyrene-d10	19.72	212	3704064	36.22	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.72	264	3183618	33.15	ng/ul	0.00

Target Compounds Ovalue

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