

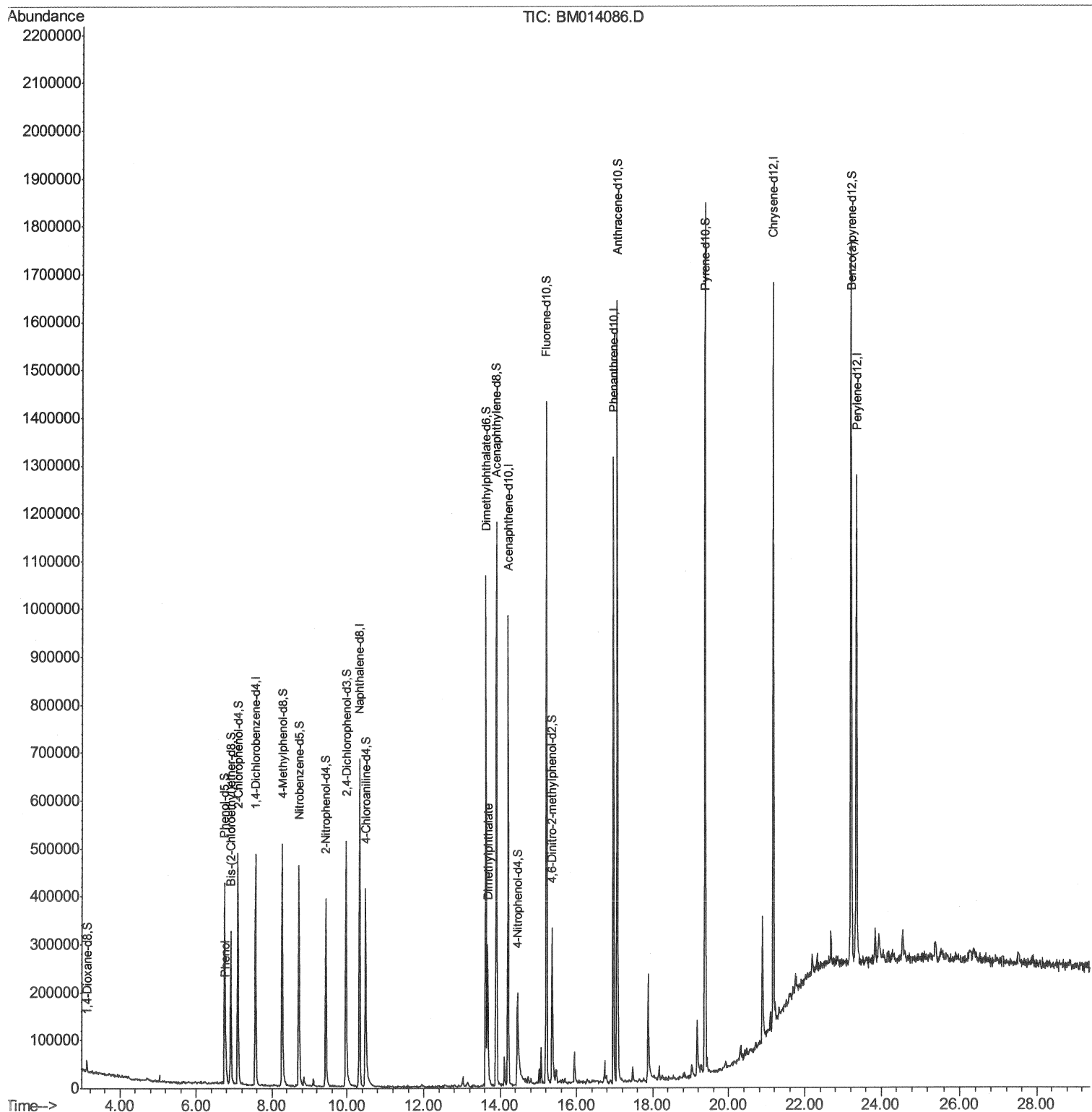
Data Path : Z:\HPCHEM1\BNA M\DATA\BM020118\
Data File : BM014086.D
Acq On : 01 Feb 2018 22:46
Operator : SJ/JU
Sample : J1317-12
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F6D08

Manual Integrations
APPROVED

Sohil
2/2/2018 4:39:30 PM

Quant Time: Feb 02 02:56:12 2018
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 02 01:03:24 2018
Response via : Initial Calibration



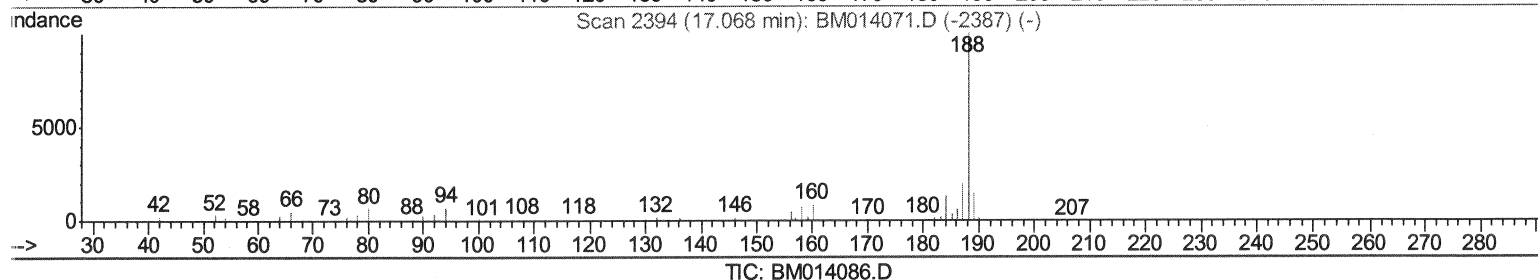
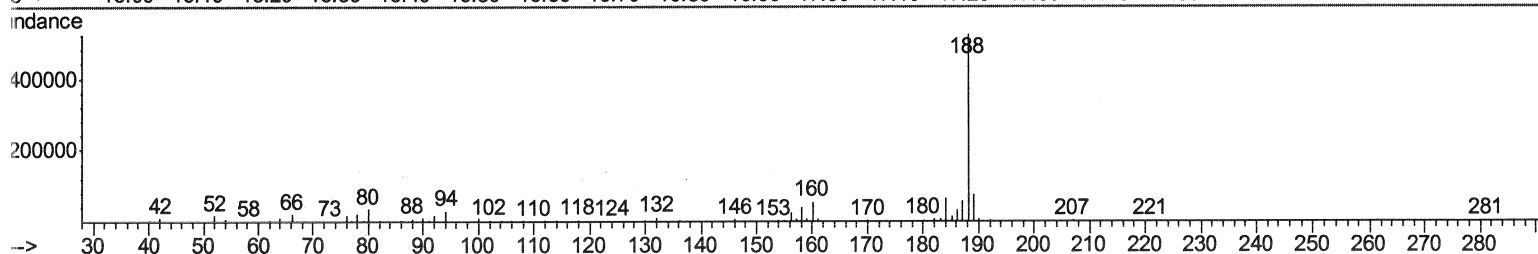
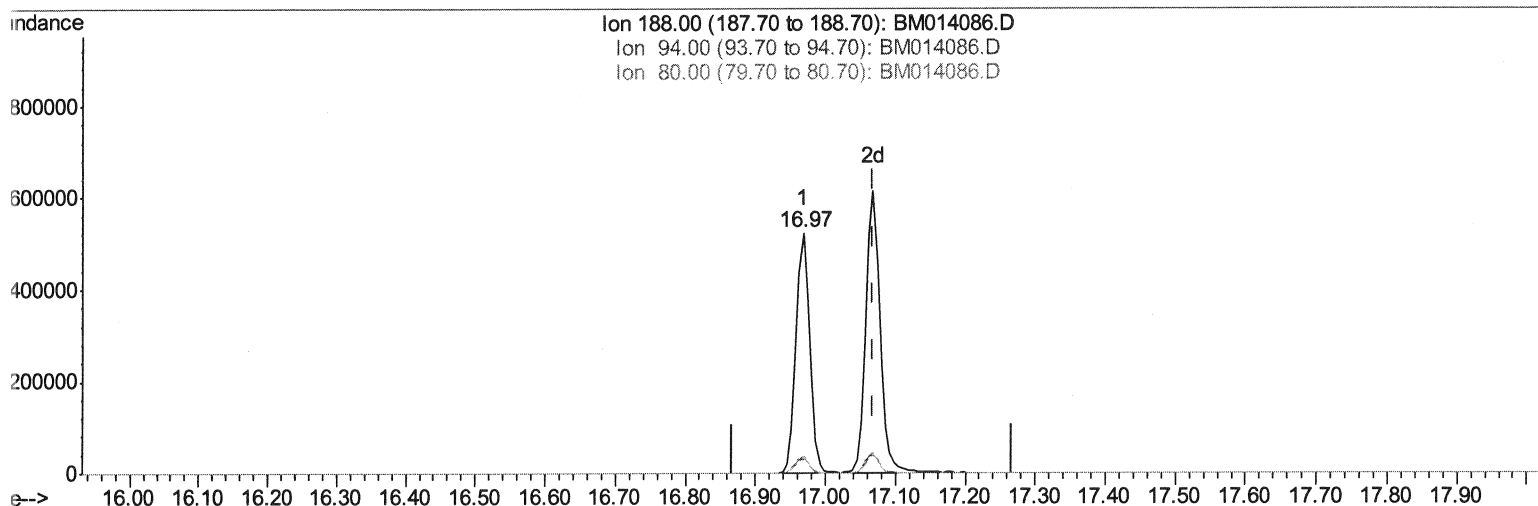
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TIC: BM014086.D

(70) Anthracene-d10 (S)

16.968min (-0.100) 20.60ng/ul

response 715616

Ion	Exp%	Act%
188.00	100	100
94.00	5.70	5.84
80.00	5.90	6.83
0.00	0.00	0.00

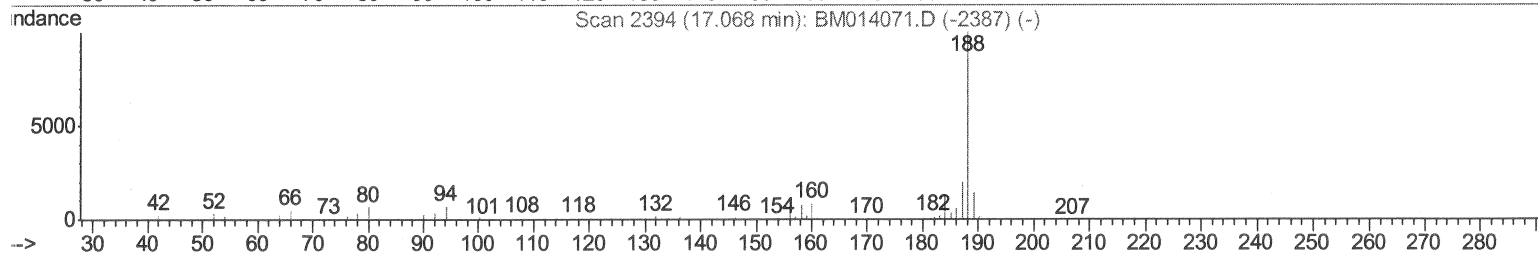
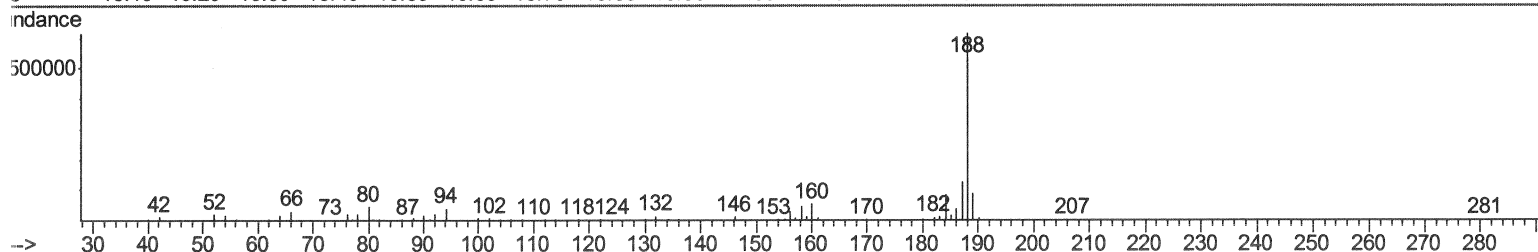
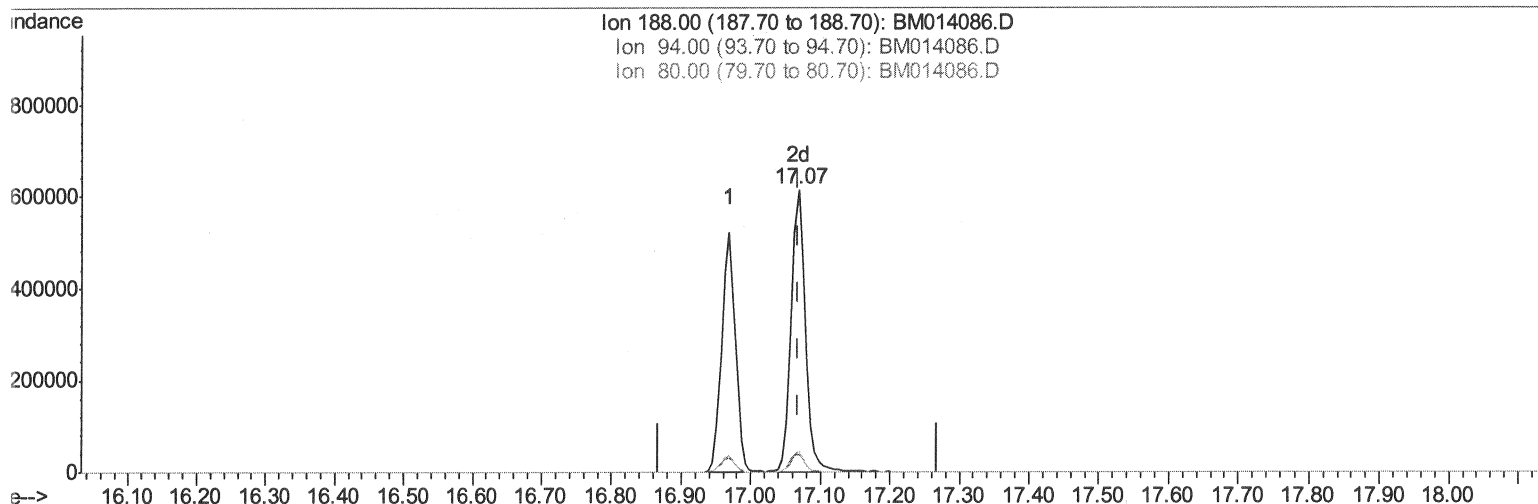
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Manual Integrations
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 Response via : Initial Calibration



TIC: BM014086.D

(70) Anthracene-d10 (S)

17.068min (+0.000) 25.08ng/ul m

response 871406

Ion	Exp%	Act%
188.00	100	100
94.00	5.70	6.42
80.00	5.90	7.40#
0.00	0.00	0.00

ata Path : Z:\HPCHEM1\BNA M\DATA\BM020118\
 ata File : BM014086.D
 ca On : 01 Feb 2018 22:46
 perator : SJ/JU
 ample : J1317-12
 isc :
 LS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F6D08

Manual Integrations
 APPROVED

Sohil
 2/2/2018 4:39:30 PM

uant Time: Feb 02 02:56:12 2018
 uant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M
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 Last Update : Fri Feb 02 01:03:24 2018
 esponse via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	118524	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	477737	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.21	164	301986	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.97	188	715616	20.00	ng/ul	0.00
75) Chrysene-d12	21.17	240	765247	20.00	ng/ul	0.00
83) Perylene-d12	23.36	264	771521	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.12	96	10472	4.05	ng/uL	0.00
5) Phenol-d5	6.75	99	208491	23.40	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.90	67	128056	23.83	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	180379	25.57	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	176635	23.68	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	88395	25.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	99720	26.13	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.97	165	187149	23.31	ng/ul	0.00
29) 4-Chloroaniline-d4	10.48	131	211434	30.38	ng/ul	0.00
43) Dimethylphthalate-d6	13.63	166	625771	24.97	ng/ul	0.00
46) Acenaphthylene-d8	13.90	160	790208	25.99	ng/ul	0.00
51) 4-Nitrophenol-d4	14.46	143	57062	16.52	ng/ul	0.00
57) Fluorene-d10	15.22	176	557473	24.85	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.36	200	73893	16.82	ng/ul	0.00
70) Anthracene-d10	17.07	188	871406m	25.08	ng/ul	0.00
76) Pyrene-d10	19.37	212	978832	26.56	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.21	264	962351	25.41	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.77	94	32347	3.630 ng/ul	96
44) Dimethylphthalate	13.68	163	173345	7.175 ng/ul	99

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