

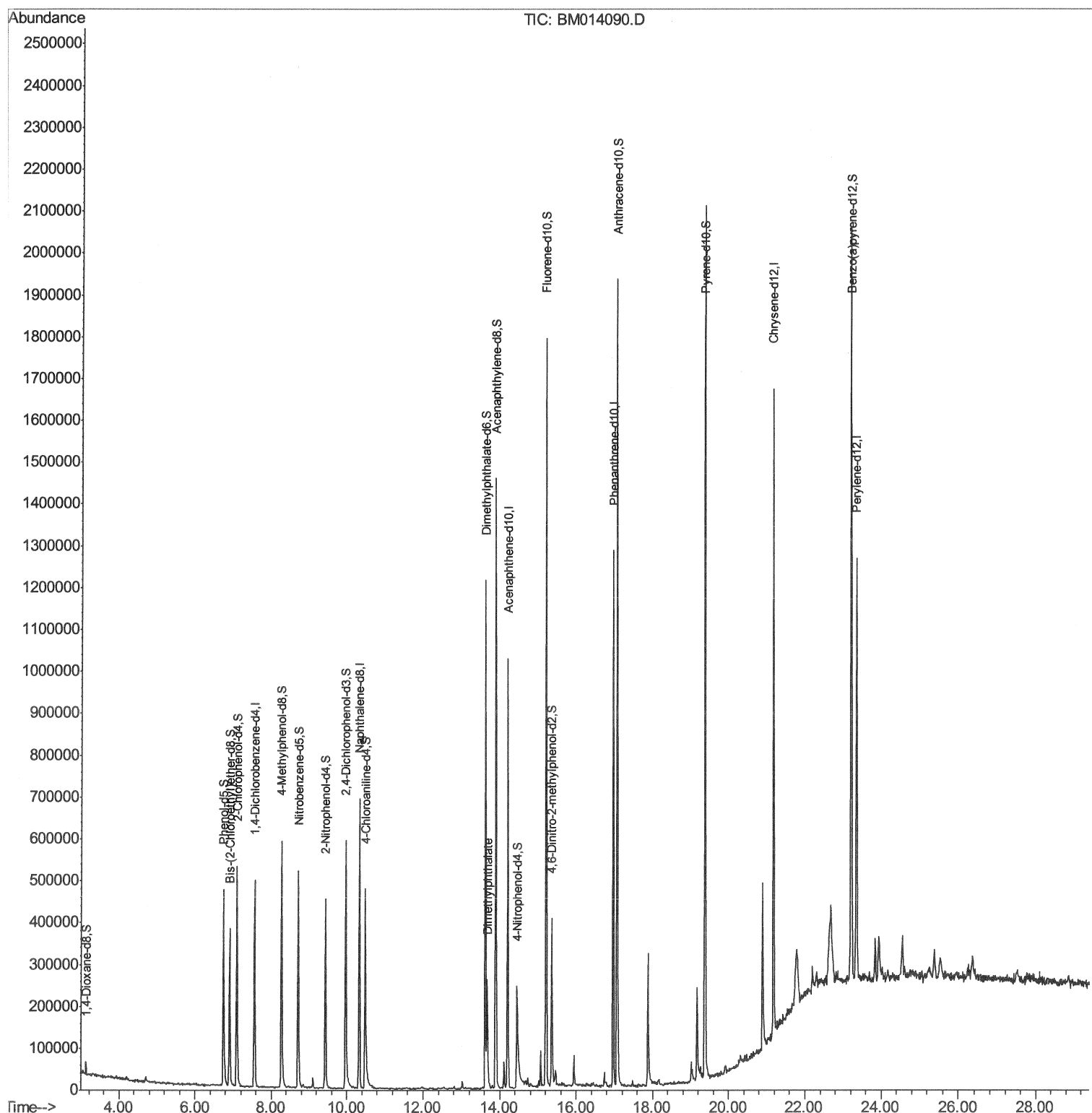
Data Path : Z:\HPCHEM1\BNA M\DATA\BM020118\
Data File : BM014090.D
Acq On : 02 Feb 2018 01:09
Operator : SJ/JU
Sample : J1316-14
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleID :
F6D37

Manual Integrations
APPROVED

Sohil
2/2/2018 4:39:42 PM

Quant Time: Feb 02 03:29:07 2018
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 02 03:15:28 2018
Response via : Initial Calibration



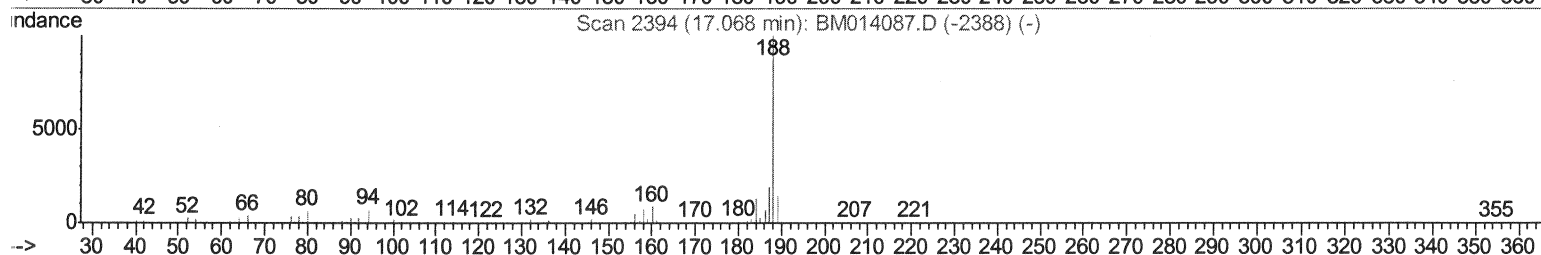
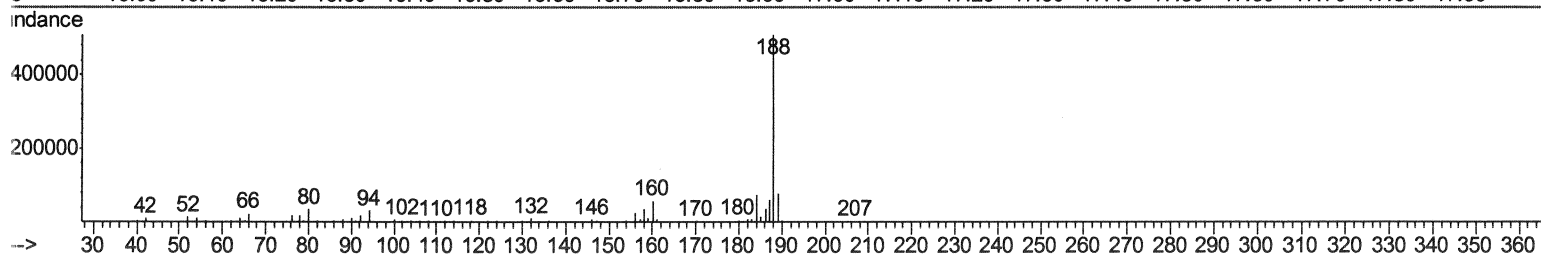
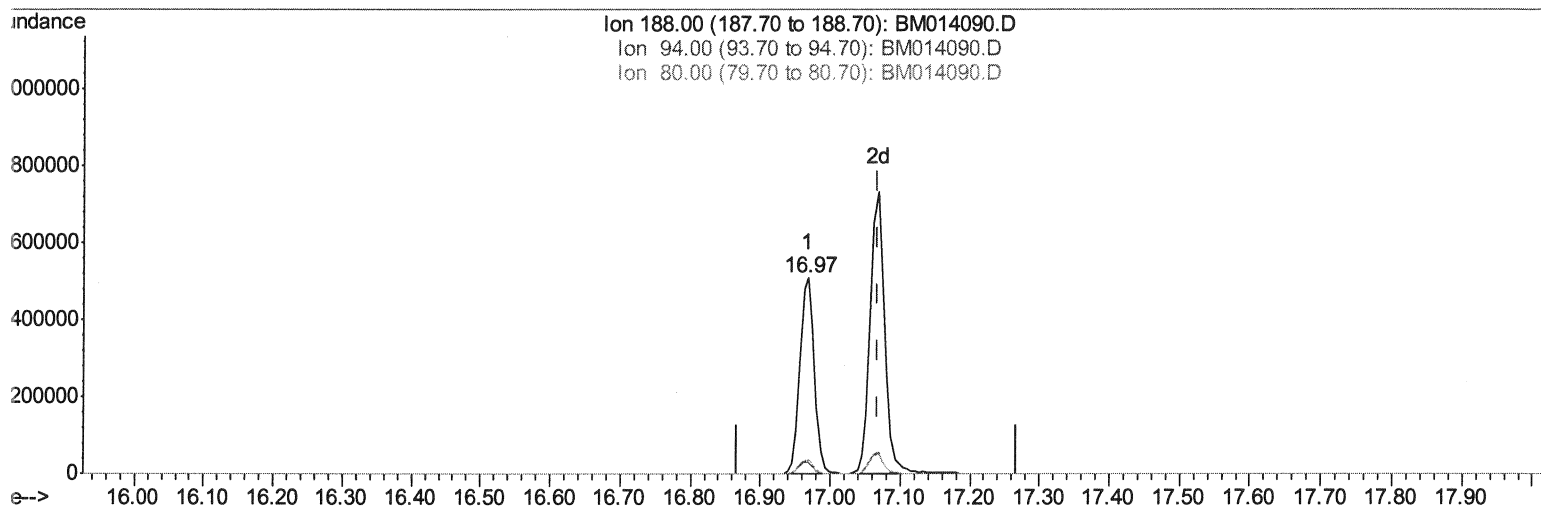
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TIC: BM014090.D

(70) Anthracene-d10 (S)

16.969min (-0.100) 20.60ng/ul

response 723562

Ion	Exp%	Act%
188.00	100	100
94.00	5.70	6.12
80.00	5.90	6.90
0.00	0.00	0.00

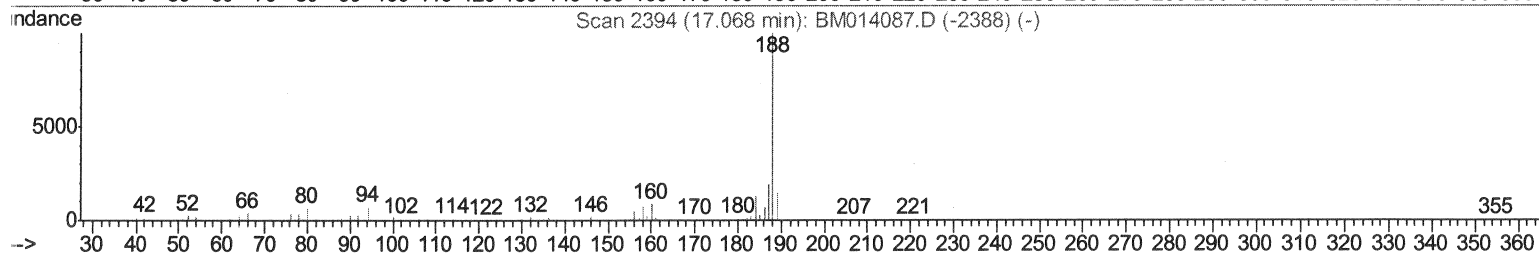
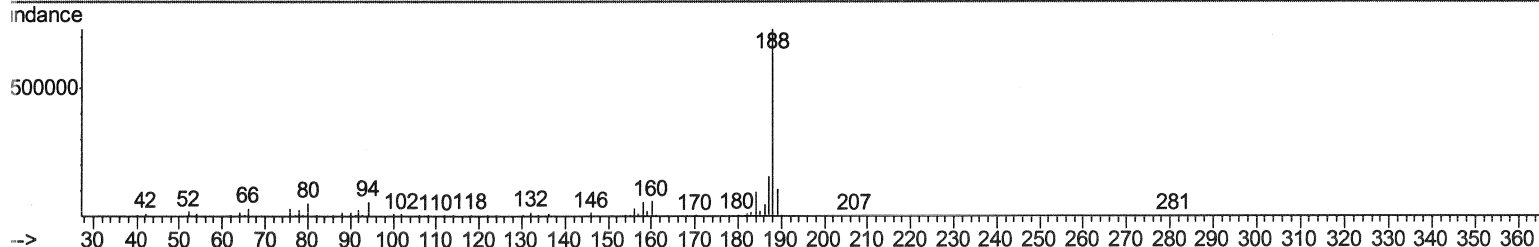
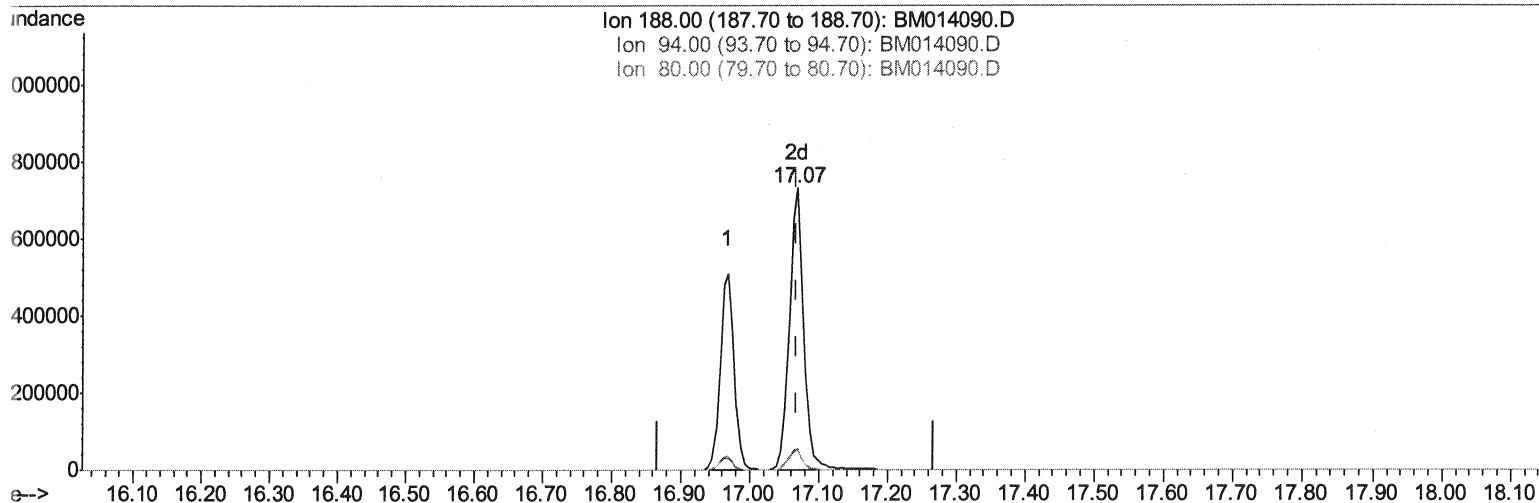
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 Operator : SJ/JU
 Sample : J1316-14
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
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Manual Integrations
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 QLast Update : Fri Feb 02 03:15:28 2018
 Response via : Initial Calibration



TIC: BM014090.D

(70) Anthracene-d10 (S)

17.069min (+0.000) 29.82ng/ul m

response 1047444

Ion	Exp%	Act%
188.00	100	100
94.00	5.70	7.64#
80.00	5.90	7.15#
0.00	0.00	0.00

ata Path : Z:\HPCHEM1\BNA M\DATA\BM020118\
 ata File : BM014090.D
 ca On : 02 Feb 2018 01:09
 operator : SJ/JU
 ample : J1316-14
 isc :
 LS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F6D37

Manual Integrations
 APPROVED

Sohil
 2/2/2018 4:39:42 PM

uant Time: Feb 02 03:29:07 2018
 uant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M
 uant Title : SVOA CALIBRATION
 Last Update : Fri Feb 02 03:15:28 2018
 esponse via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	126275	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	504952	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.21	164	311081	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.97	188	723562	20.00	ng/ul	0.00
75) Chrysene-d12	21.17	240	767997	20.00	ng/ul	0.00
83) Perylene-d12	23.35	264	773615	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.12	96	12003	4.36	ng/uL	0.00
5) Phenol-d5	6.75	99	251481	26.50	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.90	67	162068	28.31	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	208314	27.72	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	219204	27.58	ng/ul	0.00
19) Nitrobenzene-d5	8.71	128	106644	29.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	120146	29.78	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	212440	25.04	ng/ul	0.00
29) 4-Chloroaniline-d4	10.49	131	266962	36.29	ng/ul	0.00
43) Dimethylphthalate-d6	13.63	166	767892	29.74	ng/ul	0.00
46) Acenaphthylene-d8	13.90	160	959799	30.64	ng/ul	0.00
51) 4-Nitrophenol-d4	14.46	143	70511	19.82	ng/ul	0.00
57) Fluorene-d10	15.21	176	664819	28.77	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.36	200	91618	20.62	ng/ul	0.00
70) Anthracene-d10	17.07	188	1047444m	29.82	ng/ul	0.00
76) Pyrene-d10	19.37	212	1135187	30.70	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.22	264	1130223	29.77	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
44) Dimethylphthalate	13.68	163	149740	6.017	ng/ul	99

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