

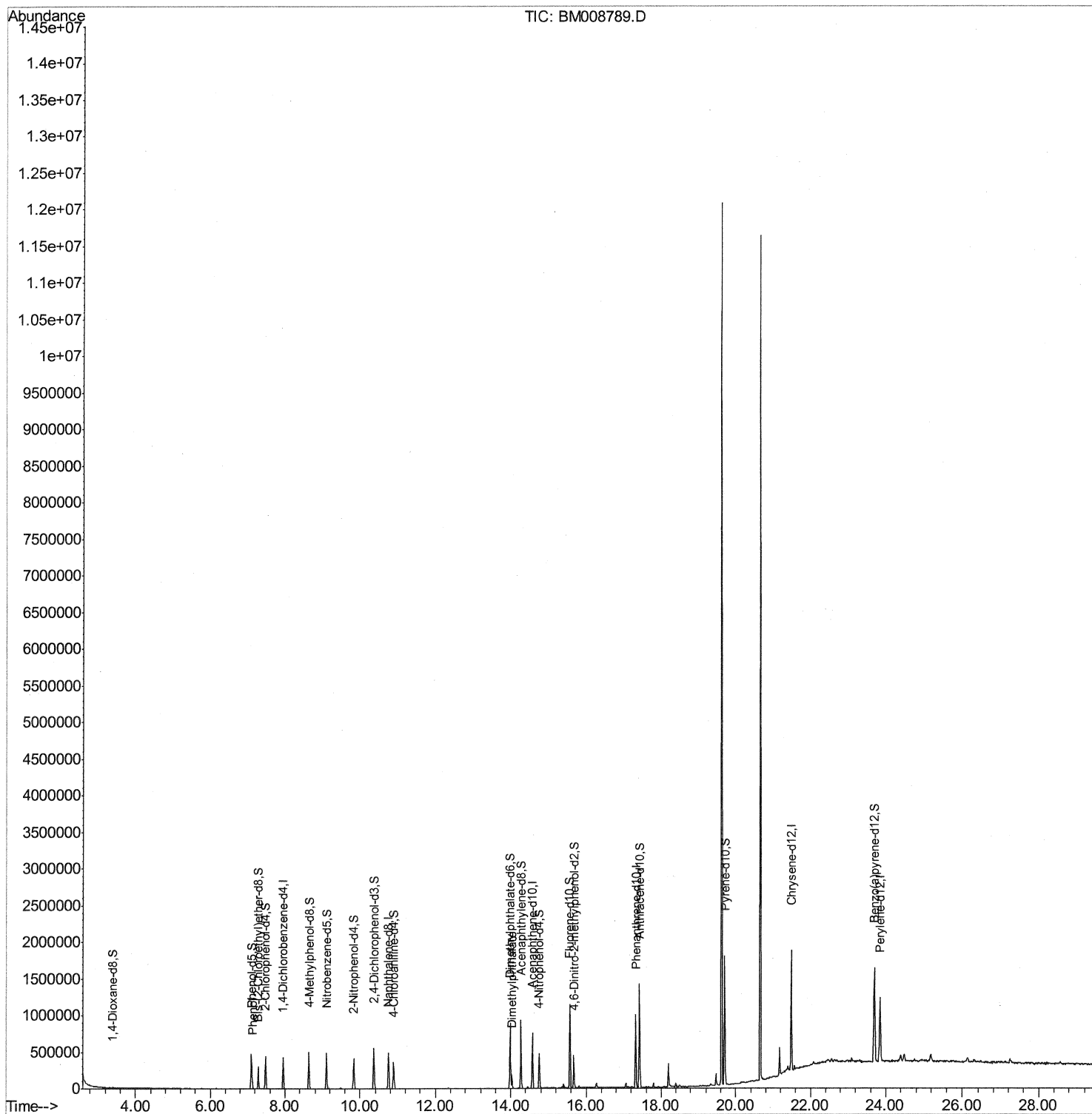
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012117\
Data File : BM008789.D
Acq On : 22 Jan 2017 02:16
Operator : UM/SJ
Sample : I1316-04
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
BDP13

Manual Integrations
APPROVED

mohammad
1/23/2017 2:49:11 PM

Quant Time: Jan 23 05:53:31 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010617.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jan 23 04:58:16 2017
Response via : Initial Calibration



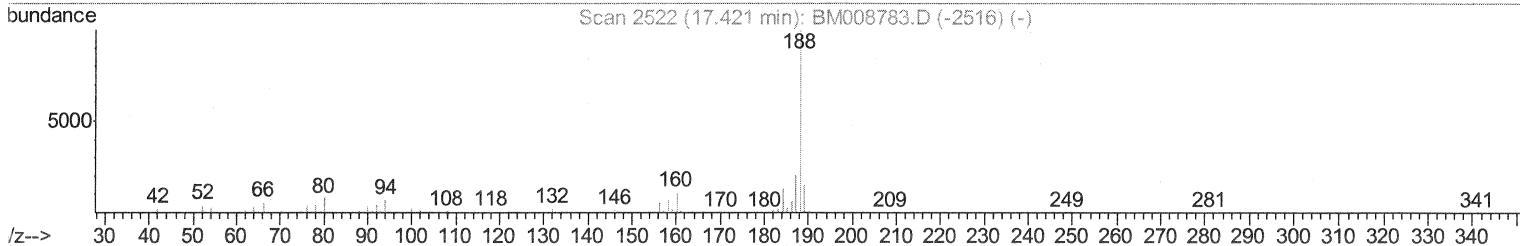
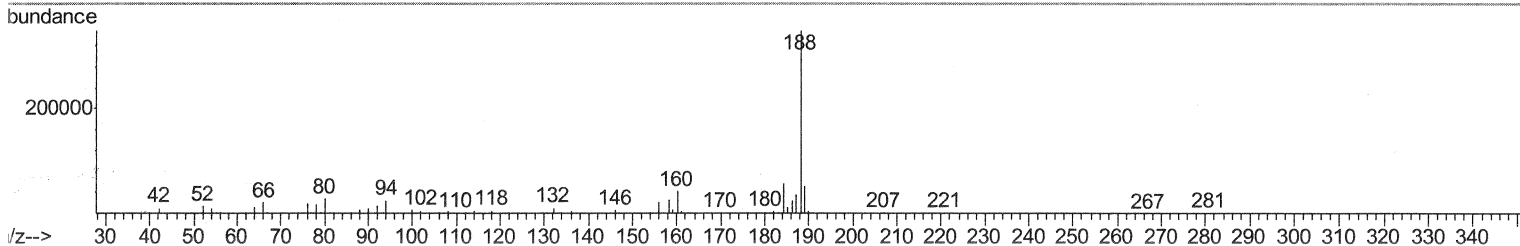
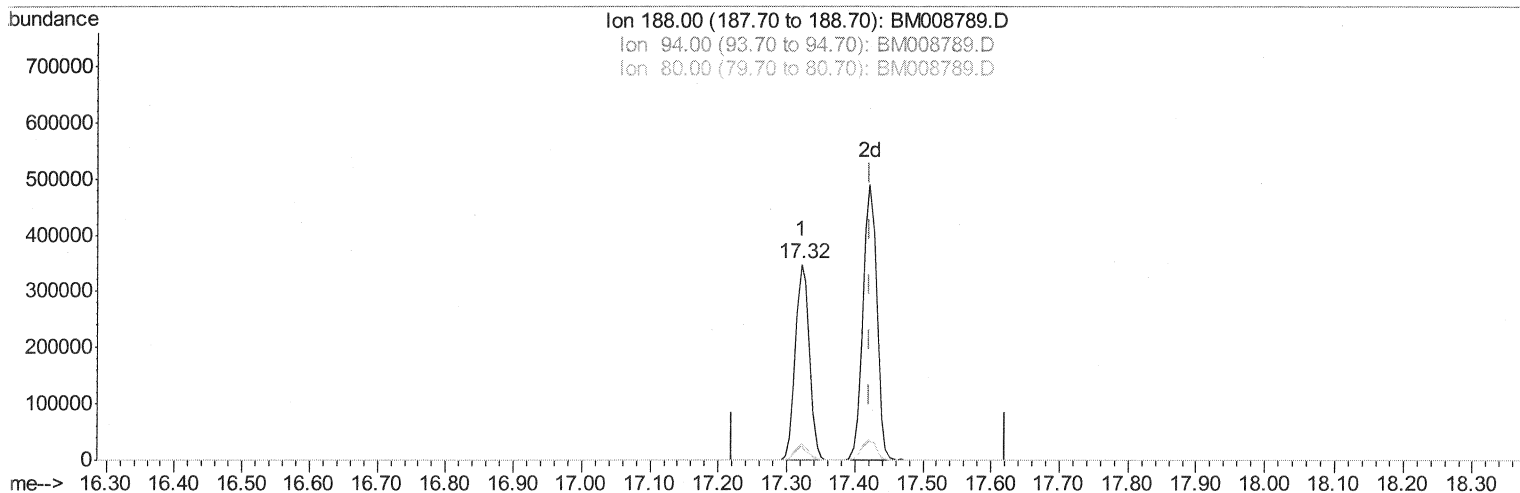
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Quant Time: Jan 23 05:03:24 2017
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TIC: BM008789.D

(70) Anthracene-d10 (S)

17.321min (-0.100) 23.21ng/ul

response 497231

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	7.24
80.00	9.20	8.21
0.00	0.00	0.00

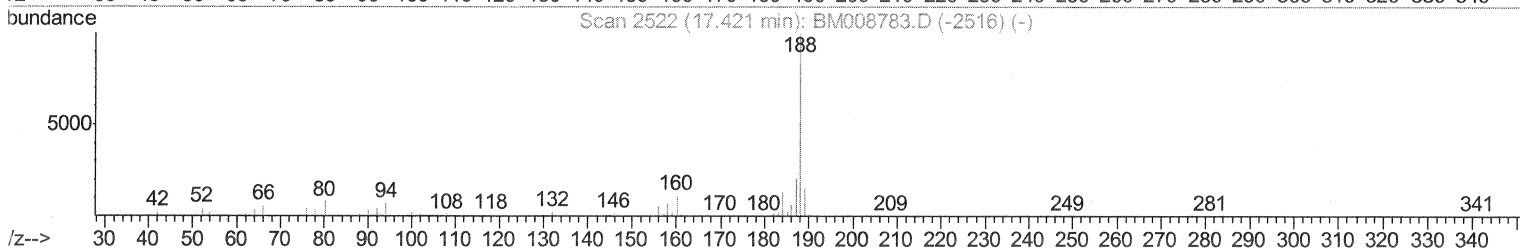
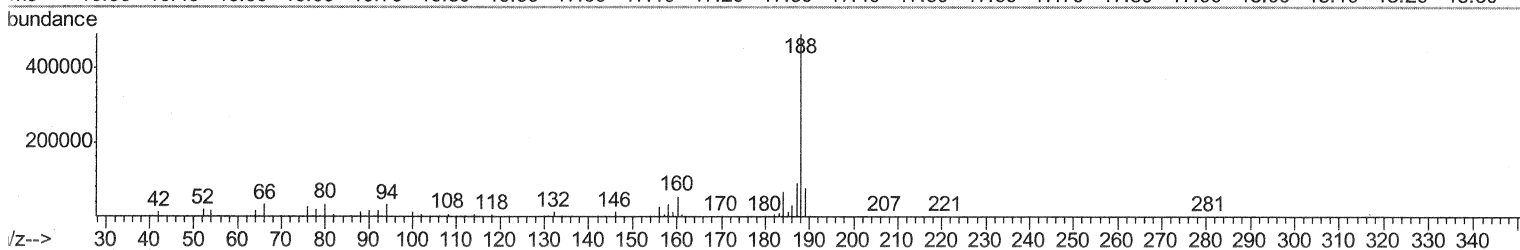
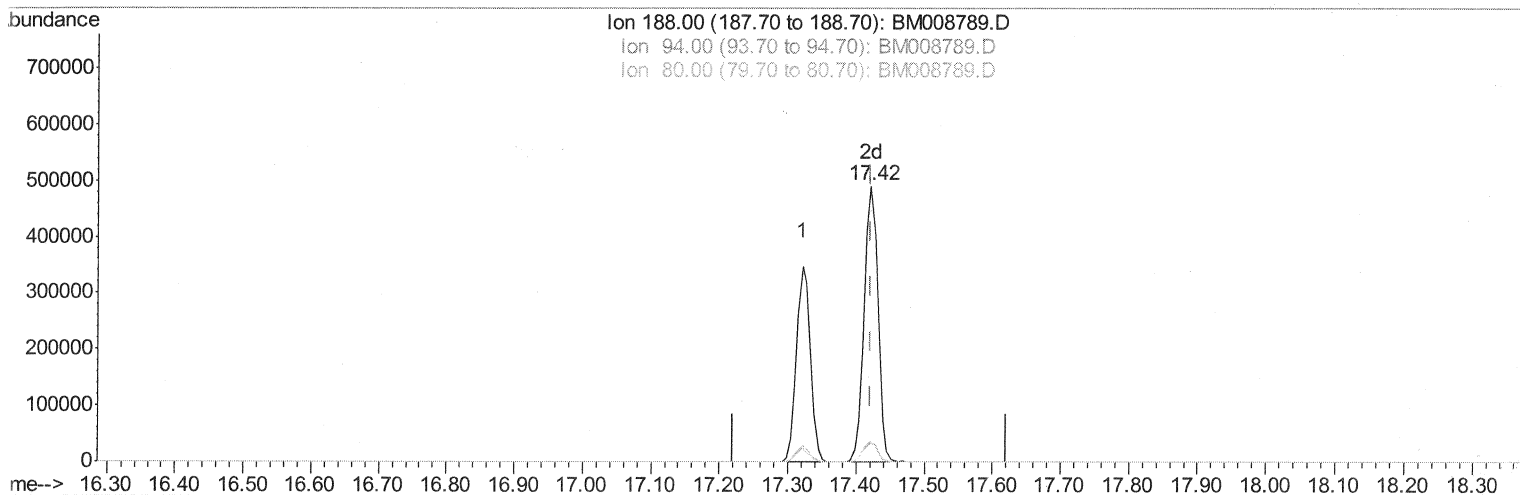
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Manual Integrations
 APPROVED

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TIC: BM008789.D

(70) Anthracene-d10 (S)

17.421min (+0.000) 31.97ng/ul m

SJ 1/22/17

response 684878

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	6.74#
80.00	9.20	7.16#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012117\
 Data File : BM008789.D
 Acq On : 22 Jan 2017 02:16
 Operator : UM/SJ
 Sample : I1316-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BDP13

Manual Integrations
 APPROVED

mohammad
 1/23/2017 2:49:11 PM

Quant Time: Jan 23 05:53:31 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010617.M
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	73411	20.00	ng/ul	0.00
18) Naphthalene-d8	10.75	136	283162	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.58	164	217334	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.32	188	497231	20.00	ng/ul	0.00
75) Chrysene-d12	21.49	240	729367	20.00	ng/ul	0.00
83) Perylene-d12	23.85	264	670292	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.40	96	5987	3.77	ng/uL	0.00
5) Phenol-d5	7.10	99	157896	26.18	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.28	67	135826	33.19	ng/ul	0.00
9) 2-Chlorophenol-d4	7.48	132	113223	26.96	ng/ul	0.00
13) 4-Methylphenol-d8	8.65	113	116262	24.85	ng/ul	0.00
19) Nitrobenzene-d5	9.12	128	54262	31.05	ng/ul	0.00
22) 2-Nitrophenol-d4	9.83	143	62766	33.05	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.36	165	126065	28.83	ng/ul	0.00
29) 4-Chloroaniline-d4	10.89	131	105819	21.84	ng/ul	0.00
43) Dimethylphthalate-d6	13.99	166	400918	27.96	ng/ul	0.00
46) Acenaphthylene-d8	14.27	160	492298	29.12	ng/ul	0.00
51) 4-Nitrophenol-d4	14.76	143	51879	22.12	ng/ul	0.00
57) Fluorene-d10	15.57	176	395251	29.89	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.68	200	62131	24.24	ng/ul	0.00
70) Anthracene-d10	17.42	188	684878m)	31.97	ng/ul	0.00
76) Pyrene-d10	19.71	212	867587	27.00	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.70	264	859204	31.43	ng/ul	0.00

SJ 1/23/17

Target Compounds

					Qvalue
6) Phenol	7.13	94	27194	4.20 ng/ul#	54
44) Dimethylphthalate	14.03	163	88342	6.16 ng/ul	98

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