

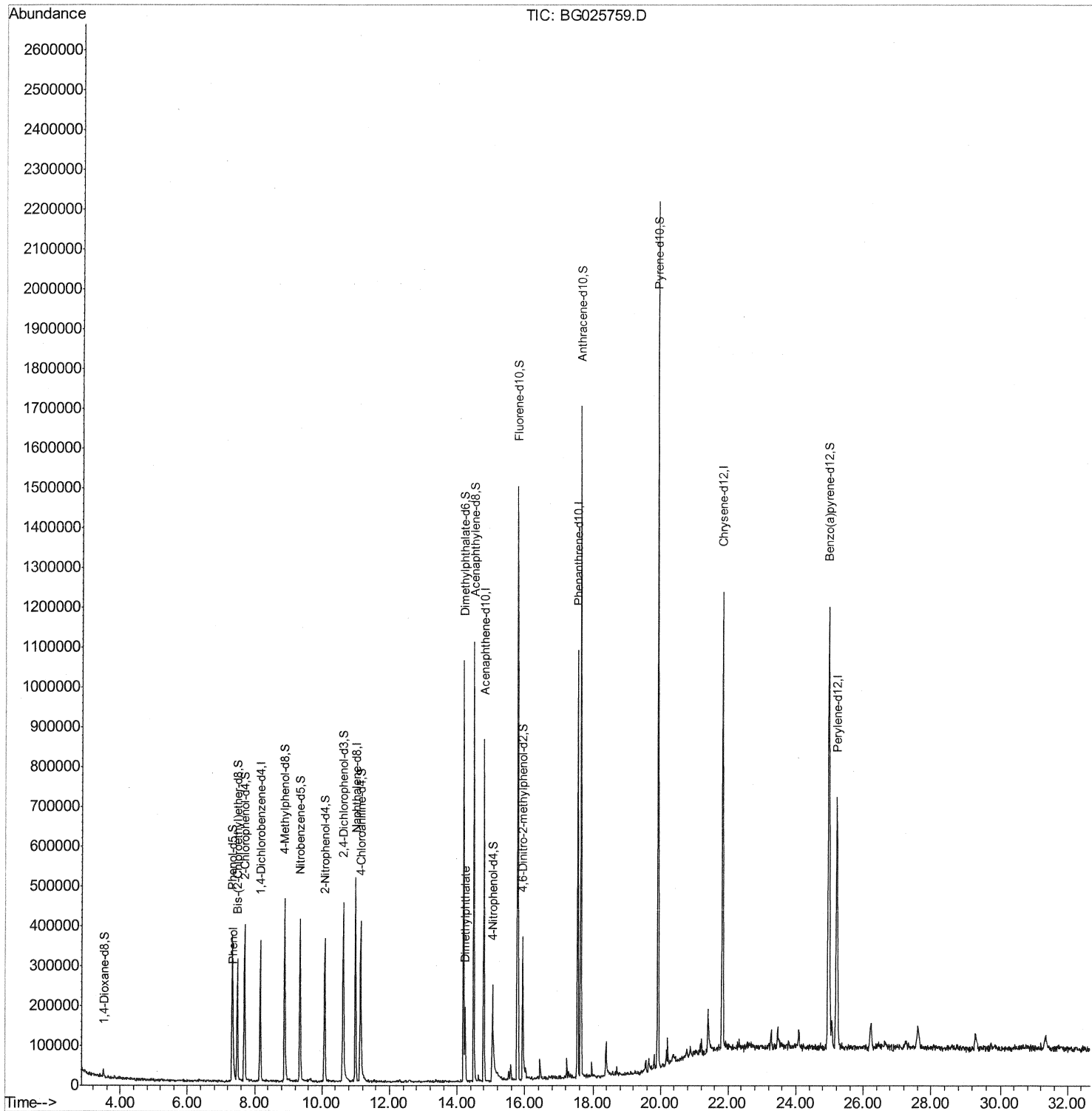
Data Path : Z:\HPCHEM1\BNA_G\Data\BG020217\
Data File : BG025759.D
Acq On : 2 Feb 2017 15:15
Operator : SJ/MA
Sample : I1495-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BDNS4

Manual Integrations
APPROVED

mohammad
2/3/2017 11:38:15 AM

Quant Time: Feb 02 16:27:15 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Feb 02 14:22:02 2017
Response via : Initial Calibration



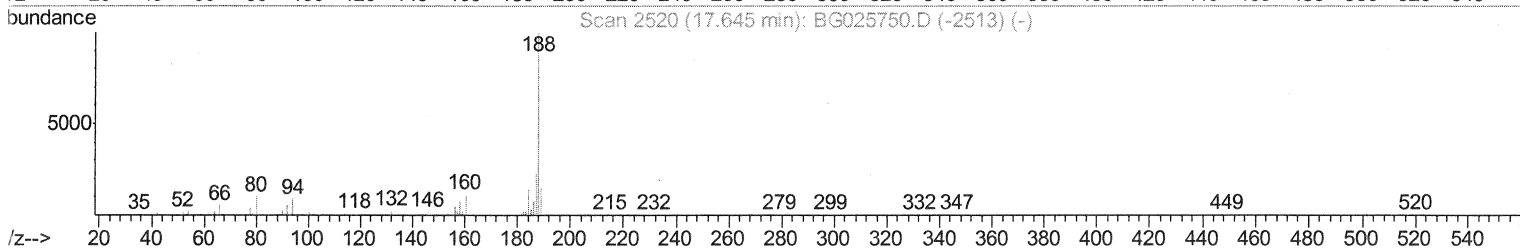
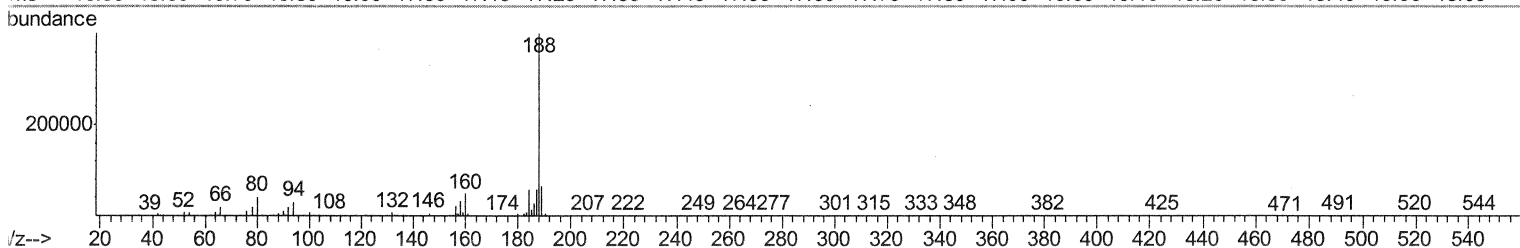
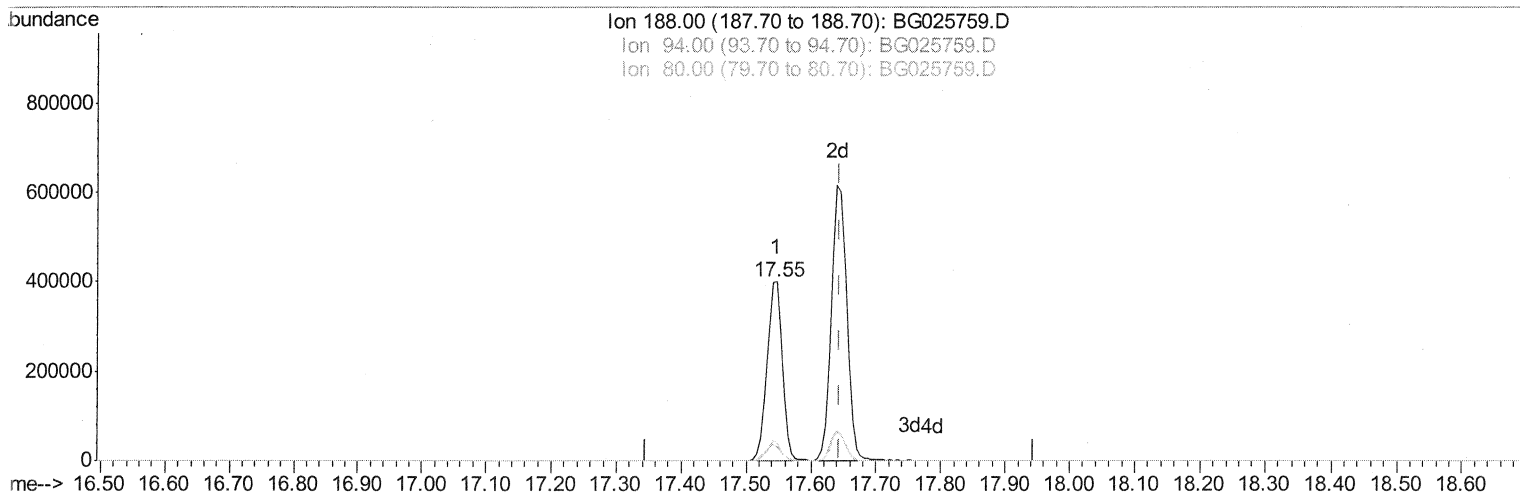
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TIC: BG025759.D

(70) Anthracene-d10 (S)

17.546min (-0.099) 23.49ng/ui

response 641238

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	7.67
80.00	9.50	10.34
0.00	0.00	0.00

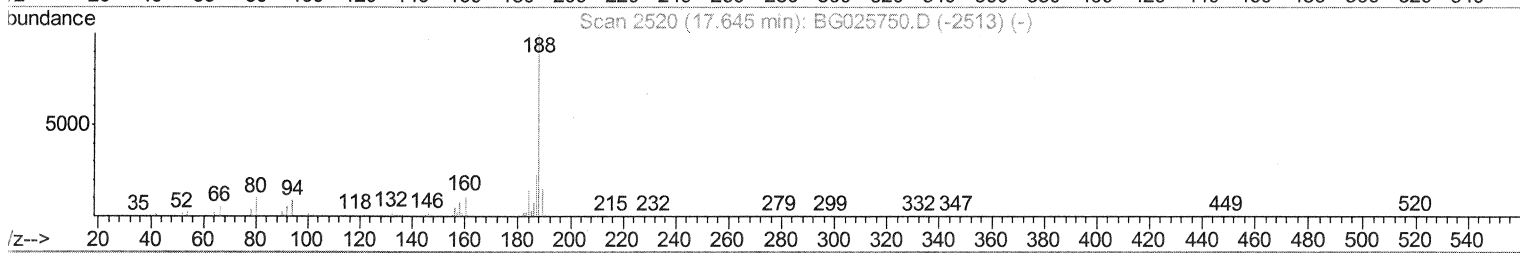
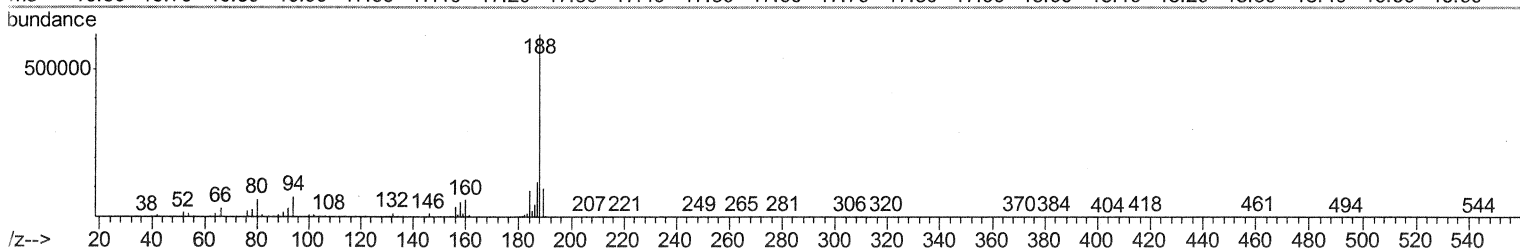
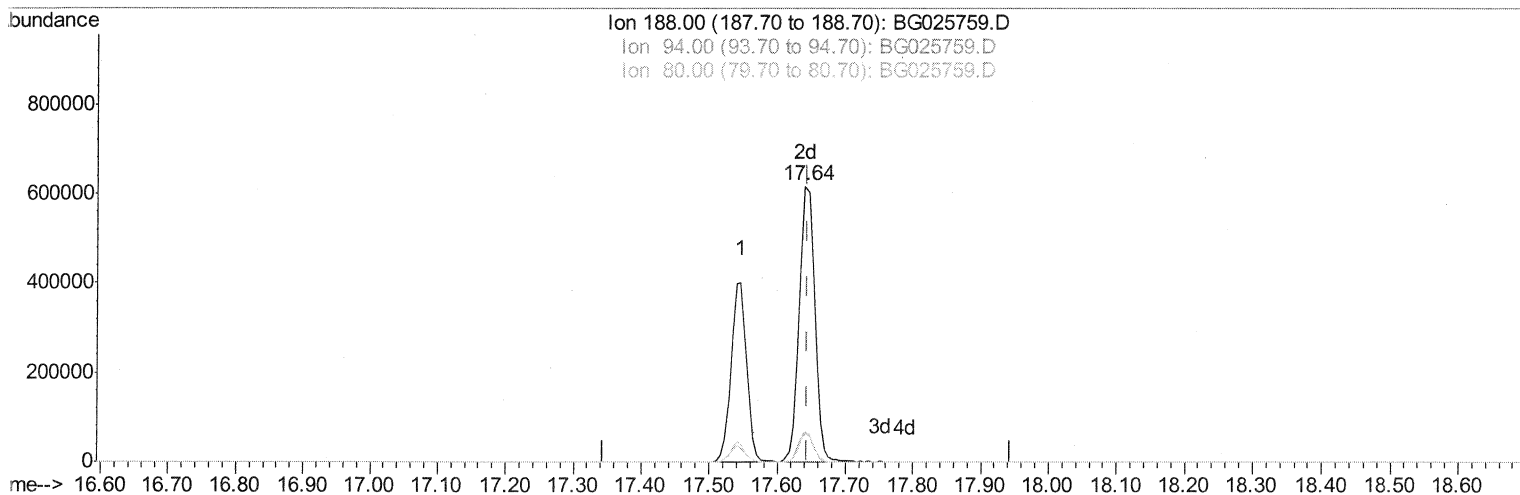
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Response via : Initial Calibration



TIC: BG025759.D

(70) Anthracene-d10 (S)

17.640min (-0.005) 35.99ng/ul m SJ 2/2/17

response 982523

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	11.14#
80.00	9.50	9.96
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\Data\BG020217\
 Data File : BG025759.D
 Acq On : 2 Feb 2017 15:15
 Operator : SJ/MA
 Sample : I1495-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BDNS4

Manual Integrations
 APPROVED

mohammad
 2/3/2017 11:38:15 AM

Quant Time: Feb 02 16:27:15 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 02 14:22:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	89479	20.00	ng/ul	0.00
18) Naphthalene-d8	10.99	136	396542	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.80	164	300784	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.55	188	641238	20.00	ng/ul	0.00
75) Chrysene-d12	21.83	240	765719	20.00	ng/ul	0.00
83) Perylene-d12	25.21	264	767052	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	10083	5.64	ng/uL	0.00
5) Phenol-d5	7.33	99	231167	30.14	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.48	67	166998	33.72	ng/ul	0.00
9) 2-Chlorophenol-d4	7.70	132	166011	30.69	ng/ul	0.00
13) 4-Methylphenol-d8	8.89	113	196942	30.21	ng/ul	0.00
19) Nitrobenzene-d5	9.36	128	84407	30.03	ng/ul	0.00
22) 2-Nitrophenol-d4	10.08	143	105290	31.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.63	165	179581	27.14	ng/ul	0.00
29) 4-Chloroaniline-d4	11.14	131	236255	31.94	ng/ul	0.00
43) Dimethylphthalate-d6	14.19	166	680283	32.03	ng/ul	0.00
46) Acenaphthylene-d8	14.50	160	791525	34.50	ng/ul	0.00
51) 4-Nitrophenol-d4	15.05	143	77334	26.57	ng/ul	0.01
57) Fluorene-d10	15.78	176	611689	30.55	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.94	200	91188	23.05	ng/ul	0.00
70) Anthracene-d10	17.64	188	982523m>	35.99	ng/ul	0.00
76) Pyrene-d10	19.92	212	1176491	36.92	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.97	264	1167188	36.20	ng/ul	0.00

SJ 2/2/17

Target Compounds

						Qvalue
6) Phenol	7.36	94	42070	5.22	ng/ul#	82
44) Dimethylphthalate	14.24	163	114219	5.47	ng/ul#	93

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