

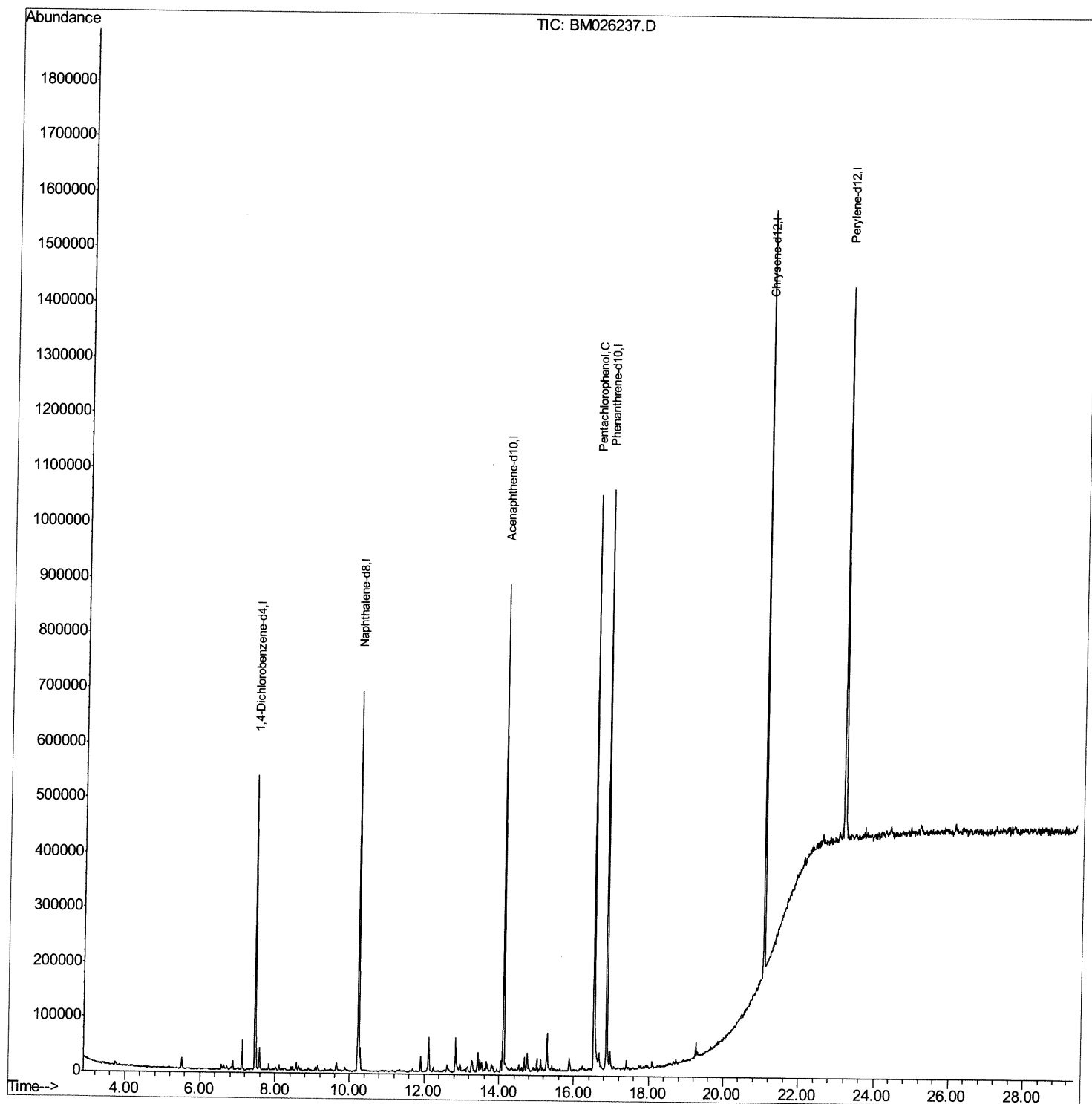
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060120\
Data File : BM026237.D
Acq On : 02 Jun 2020 20:55
Operator : JU/CG
Sample : L2829-13DL2 100X
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C5CE6DL2

Manual Integrations
APPROVED

mohammad
6/3/2020 8:43:16 AM

Quant Time: Jun 03 00:46:59 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 03 00:17:42 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

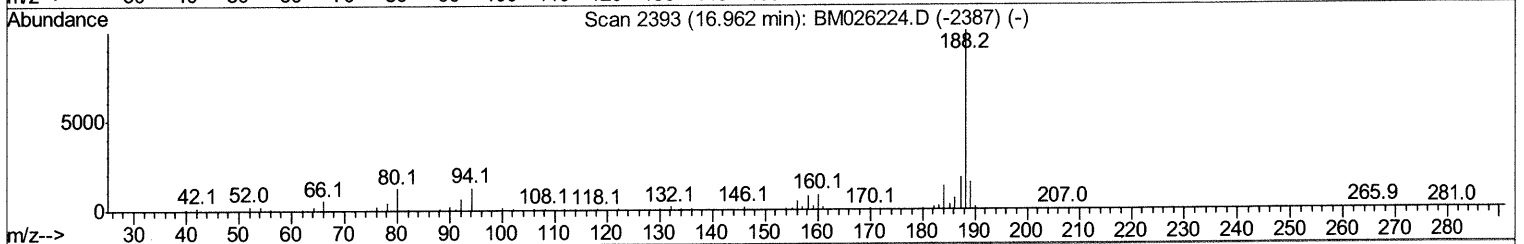
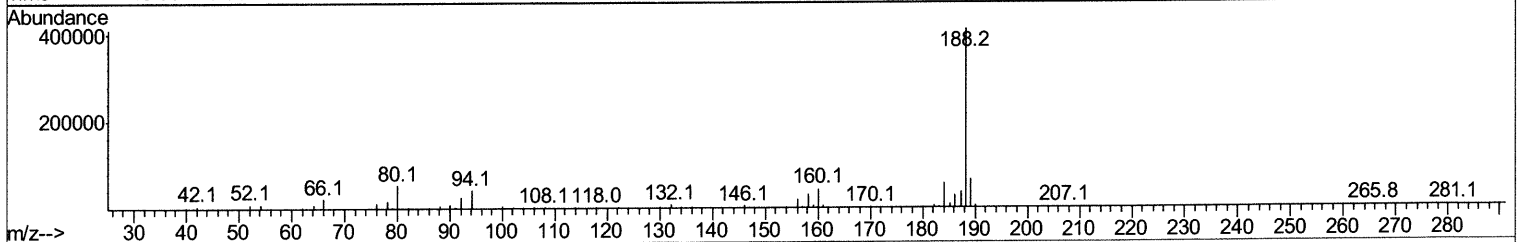
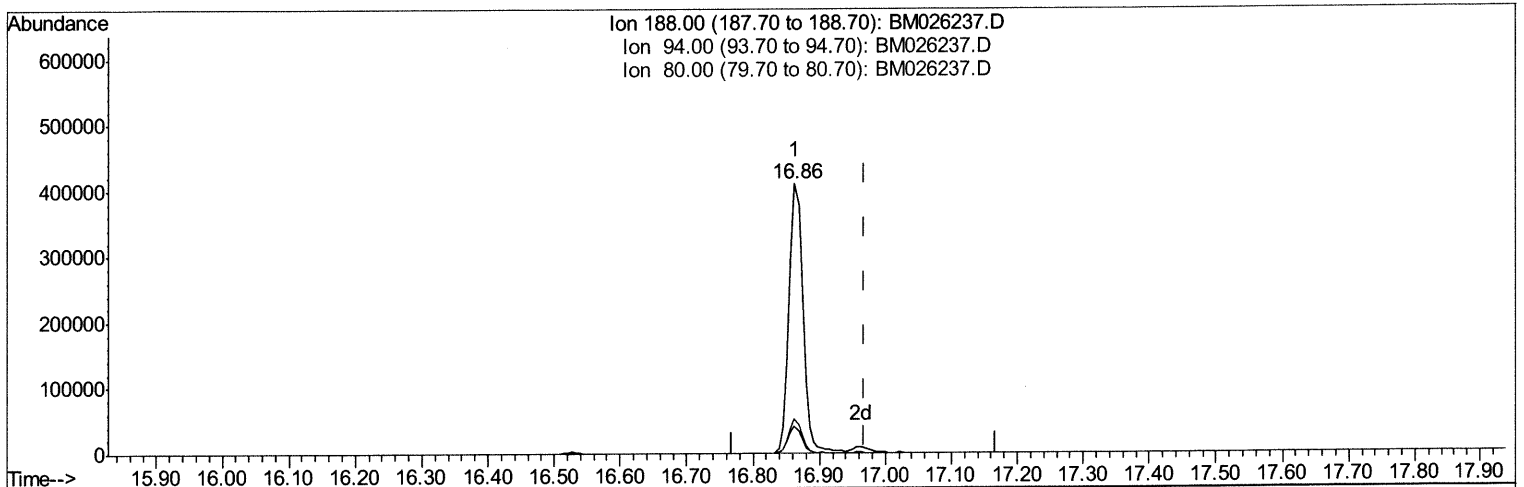
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TIC: BM026237.D

(71) Anthracene-d10 (S)

16.862min (-0.106) 20.29ng/ul

response 602238

Ion	Exp%	Act%
188.00	100	100
94.00	12.10	10.15
80.00	12.20	12.94
0.00	0.00	0.00

Quantitation Report (Qedit)

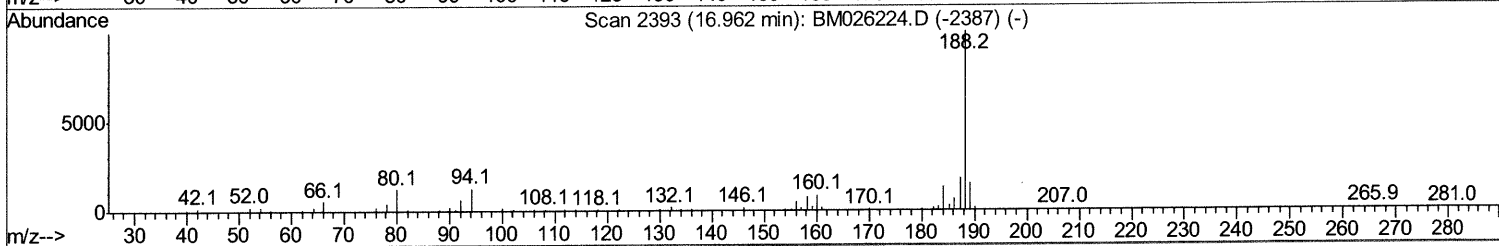
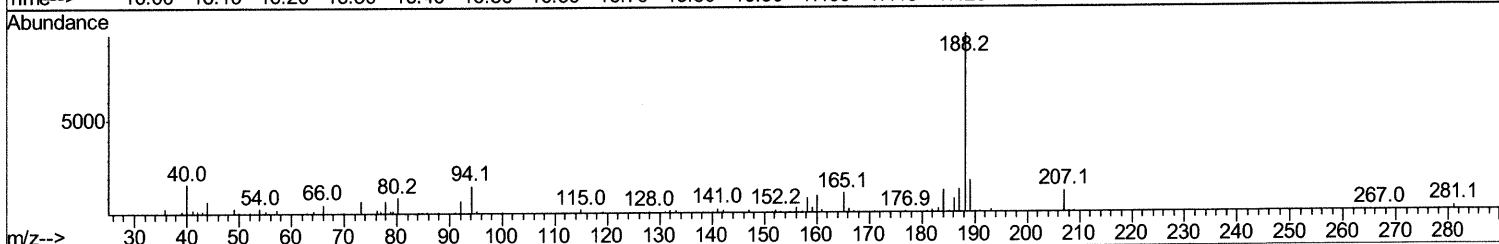
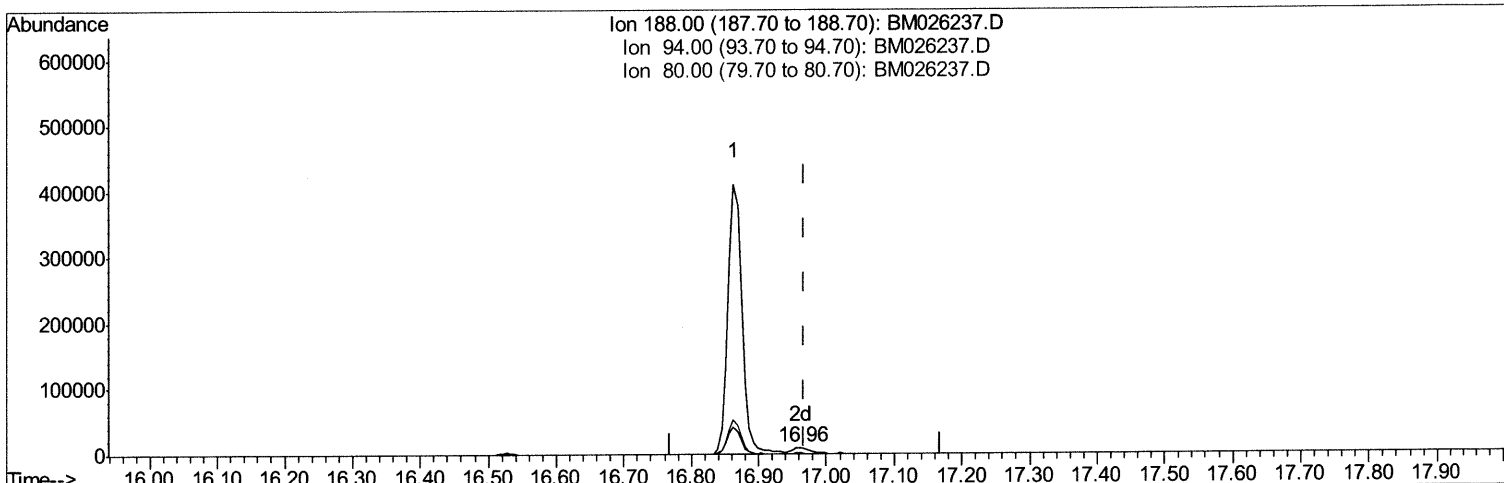
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 Response via : Initial Calibration



TIC: BM026237.D

(71) Anthracene-d10 (S)

16.962min (-0.006) 0.49ng/ul m

06/04/20 JU

response 14682

Ion	Exp%	Act%
188.00	100	100
94.00	12.10	16.53#
80.00	12.20	10.21
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060120\
 Data File : BM026237.D
 Acq On : 02 Jun 2020 20:55
 Operator : JU/CG
 Sample : L2829-13DL2 100X
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5CE6DL2

Manual Integrations
 APPROVED

mohammad
 6/3/2020 8:43:16 AM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	127963	20.00	ng/ul	0.00
18) Naphthalene-d8	10.23	136	503478	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.12	164	281840	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.86	188	602238	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	611253	20.00	ng/ul	0.00
86) Perylene-d12	23.18	264	656726	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	0.00	99	0d	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	6.83	67	2782	0.35	ng/ul	0.01
9) 2-Chlorophenol-d4	7.02	132	1918	0.21	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	1097	0.13	ng/ul	0.01
19) Nitrobenzene-d5	8.63	128	994	0.24	ng/ul	0.01
22) 2-Nitrophenol-d4	9.33	143	695	0.17	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	1628	0.20	ng/ul	0.02
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.55	166	7936	0.35	ng/ul	0.00
47) Acenaphthylene-d8	13.80	160	8690	0.32	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.12	176	7921	0.40	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
71) Anthracene-d10	16.96	188	14682m>	0.49	ng/ul >	0.00
79) Pyrene-d10	19.26	212	12367	0.39	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.05	264	10706	0.31	ng/ul	0.00

6/6/2020 JU

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
69) Pentachlorophenol	16.53	266	180380	41.535	ng/ul	97

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