

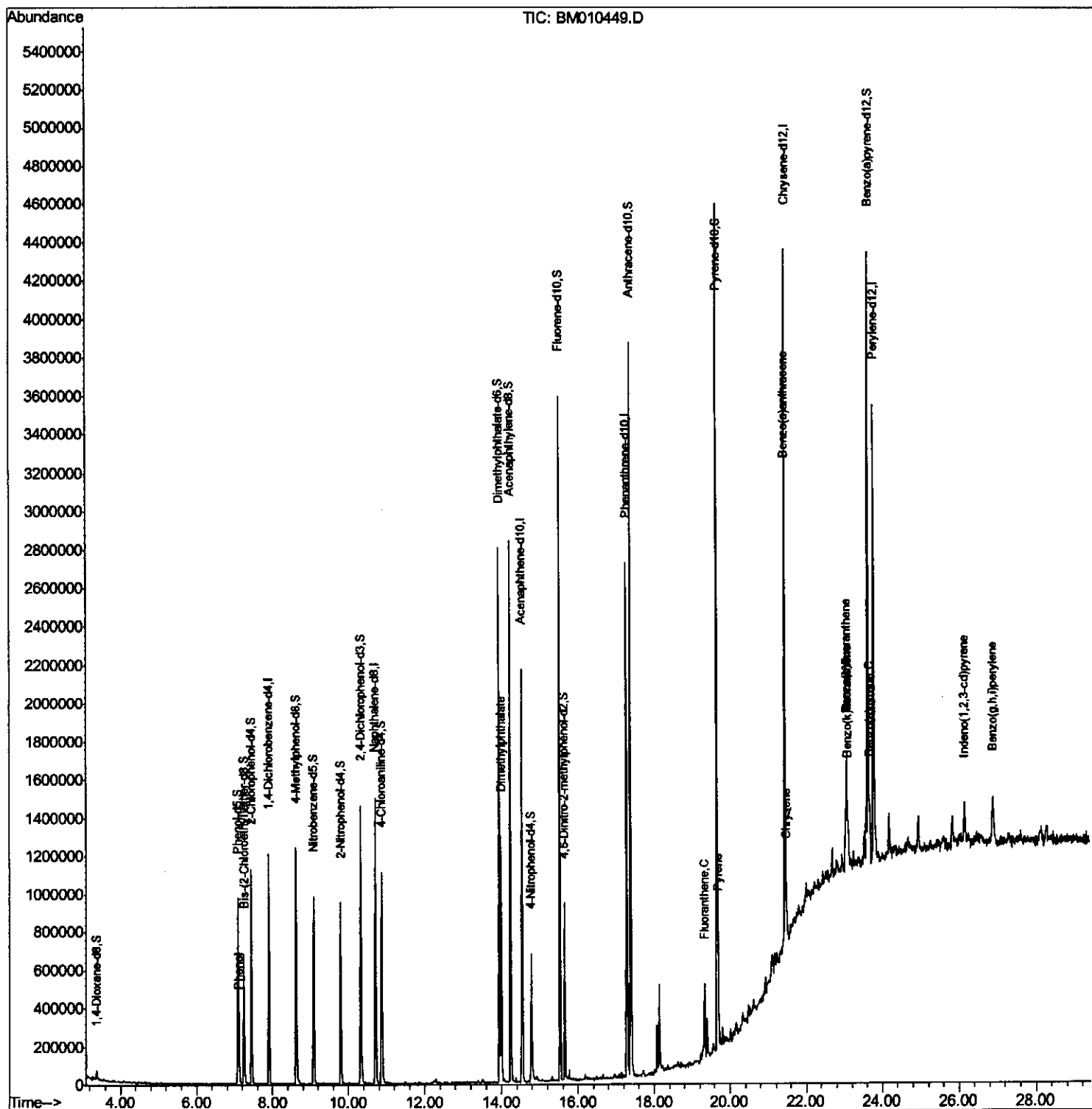
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
Data File : BM010449.D
Acq On : 09 Jun 2017 12:28
Operator : SJ/MA
Sample : I3521-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
F5C60

Manual Integrations
APPROVED

mohammad
6/12/2017 11:17:52 AM

Quant Time: Jun 10 02:00:42 2017
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jun 10 01:45:58 2017
Response via : Initial Calibration



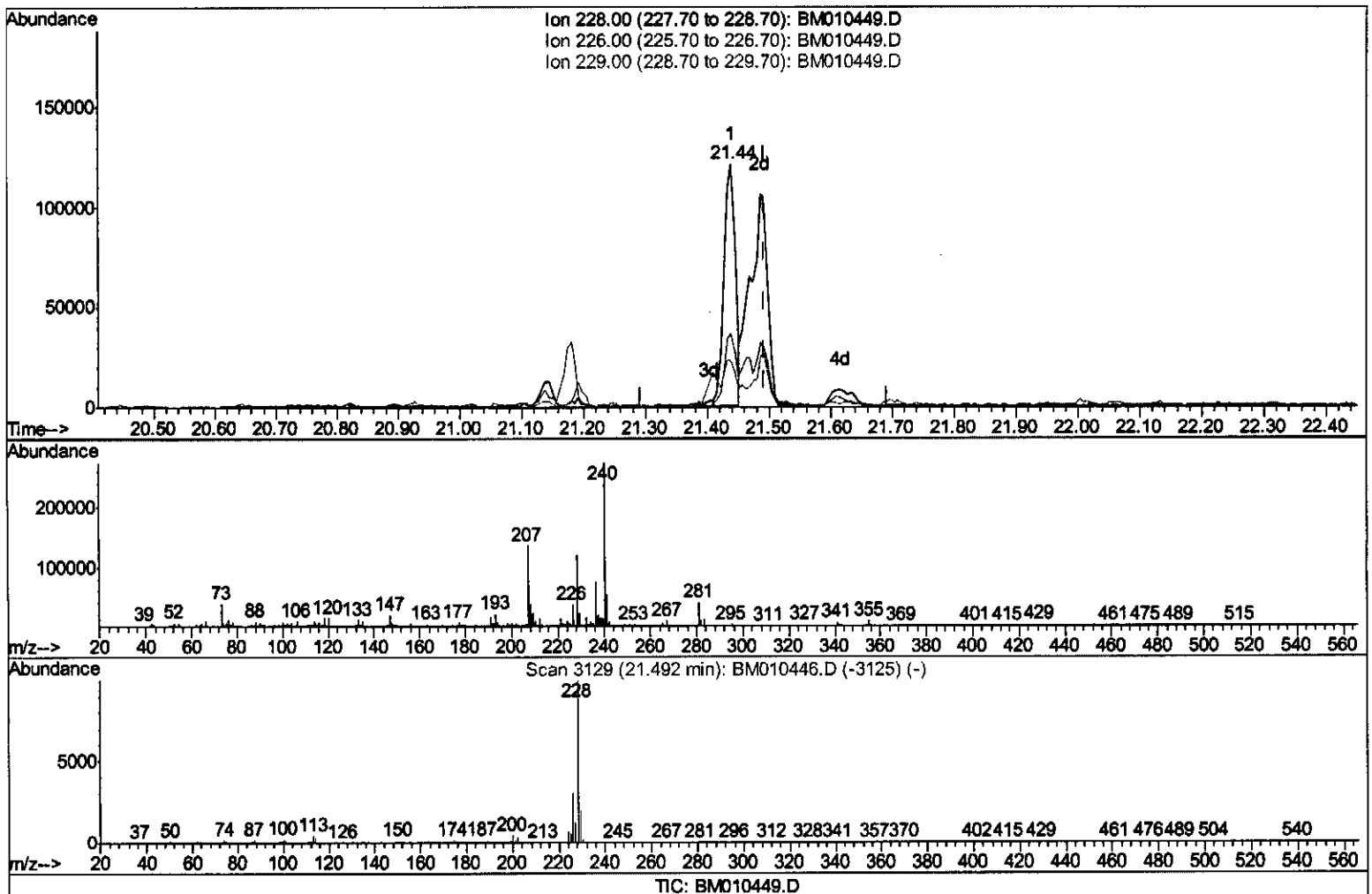
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Quant Time: Jun 10 01:47:16 2017
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(82) Chrysene

21.439min (-0.053) 1.85ng/ul

response 156587

Ion	Exp%	Act%
228.00	100	100
226.00	29.30	30.02
229.00	19.40	19.18
0.00	0.00	0.00

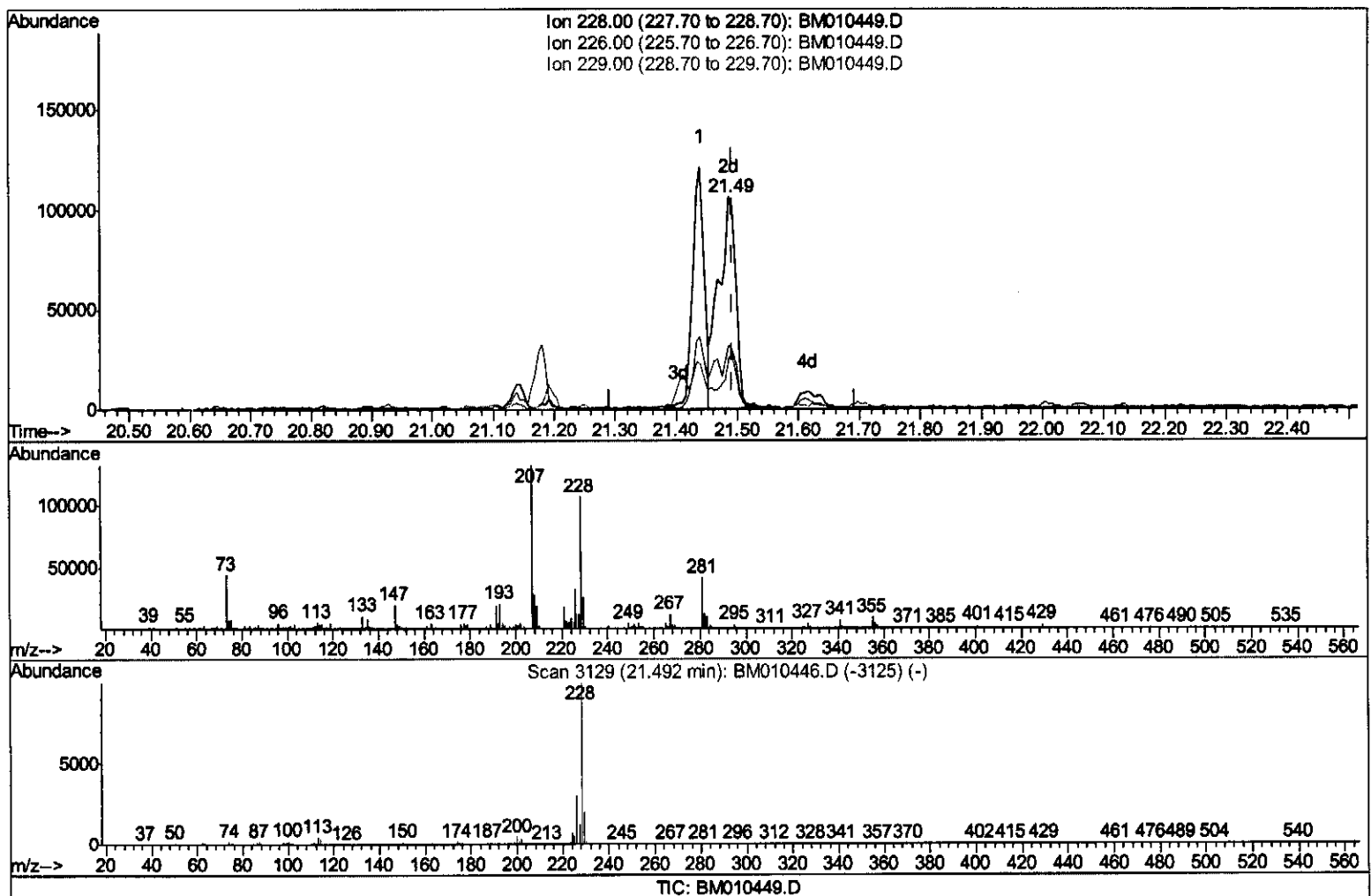
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(82) Chrysene

21.486min (-0.006) 2.45ng/ul m

response 208005

Ion	Exp%	Act%
228.00	100	100
226.00	29.30	30.22
229.00	19.40	23.71#
0.00	0.00	0.00

SJ
06/19/17

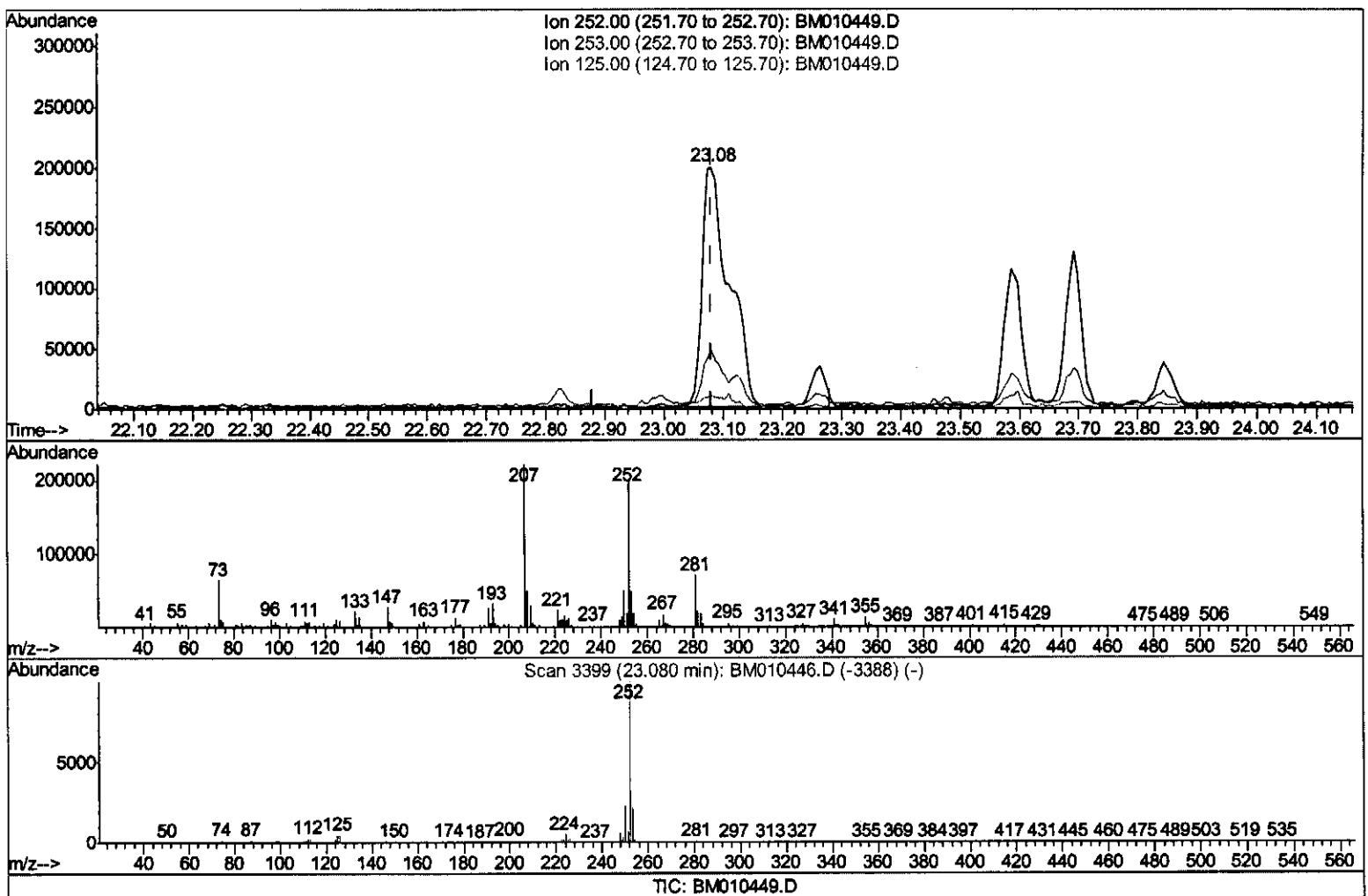
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(85) Benzo(b)fluoranthene

23.080min (-0.000) 6.25ng/ul

response 620974

Ion	Exp%	Act%
252.00	100	100
253.00	22.40	24.65
125.00	4.50	5.29
0.00	0.00	0.00

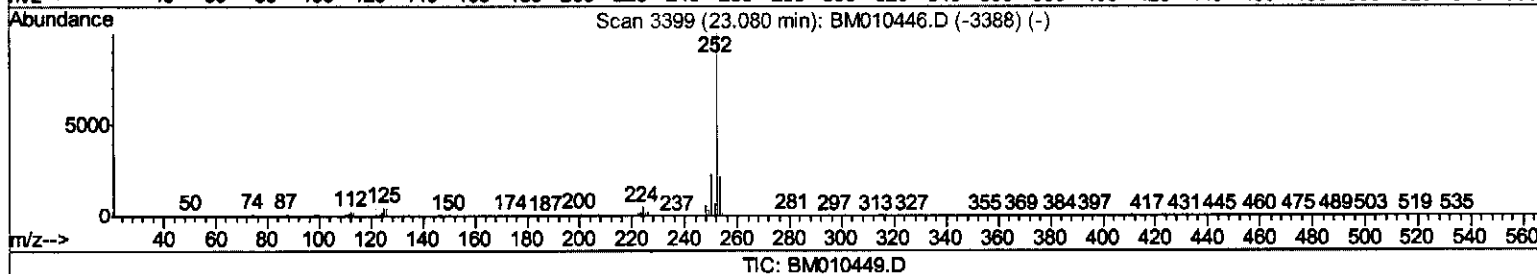
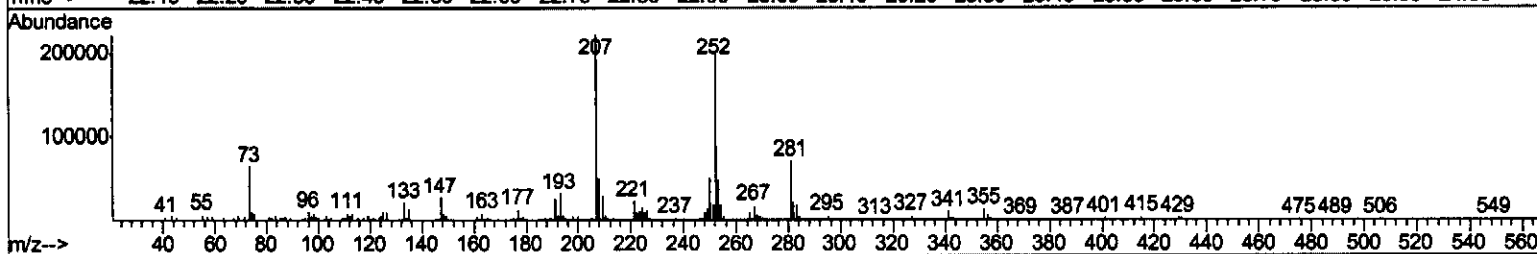
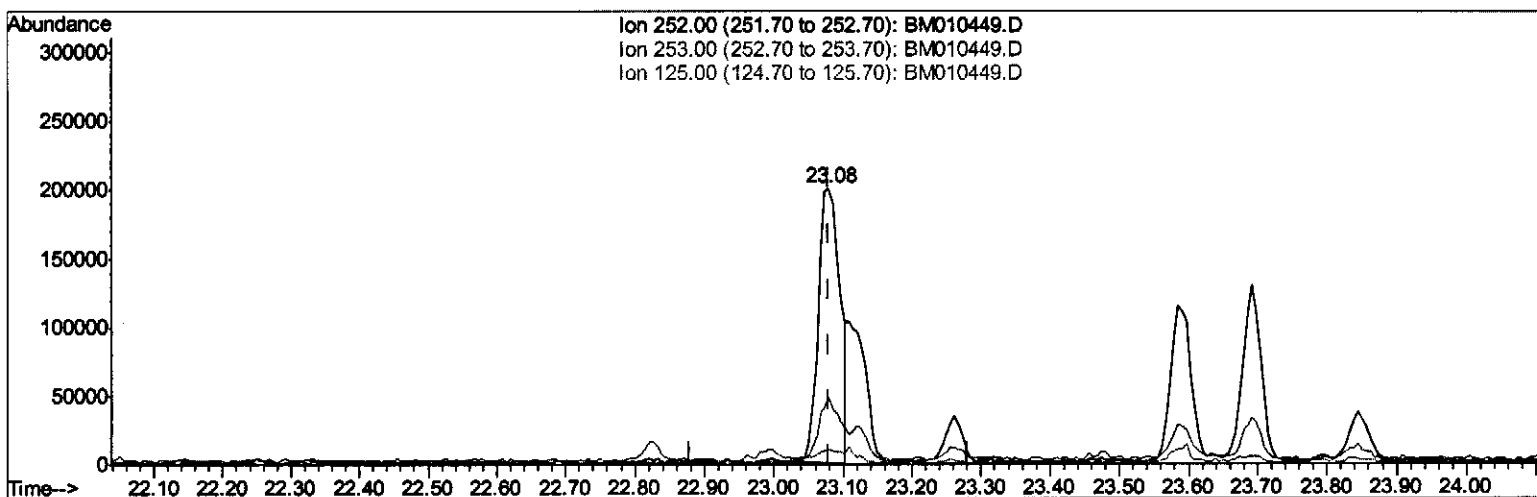
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(85) Benzo(b)fluoranthene

23.080min (-0.000) 4.42ng/ul m

SJ
06/19/17

response 438934

Ion	Exp%	Act%
252.00	100	100
253.00	22.40	24.65
125.00	4.50	5.29
0.00	0.00	0.00

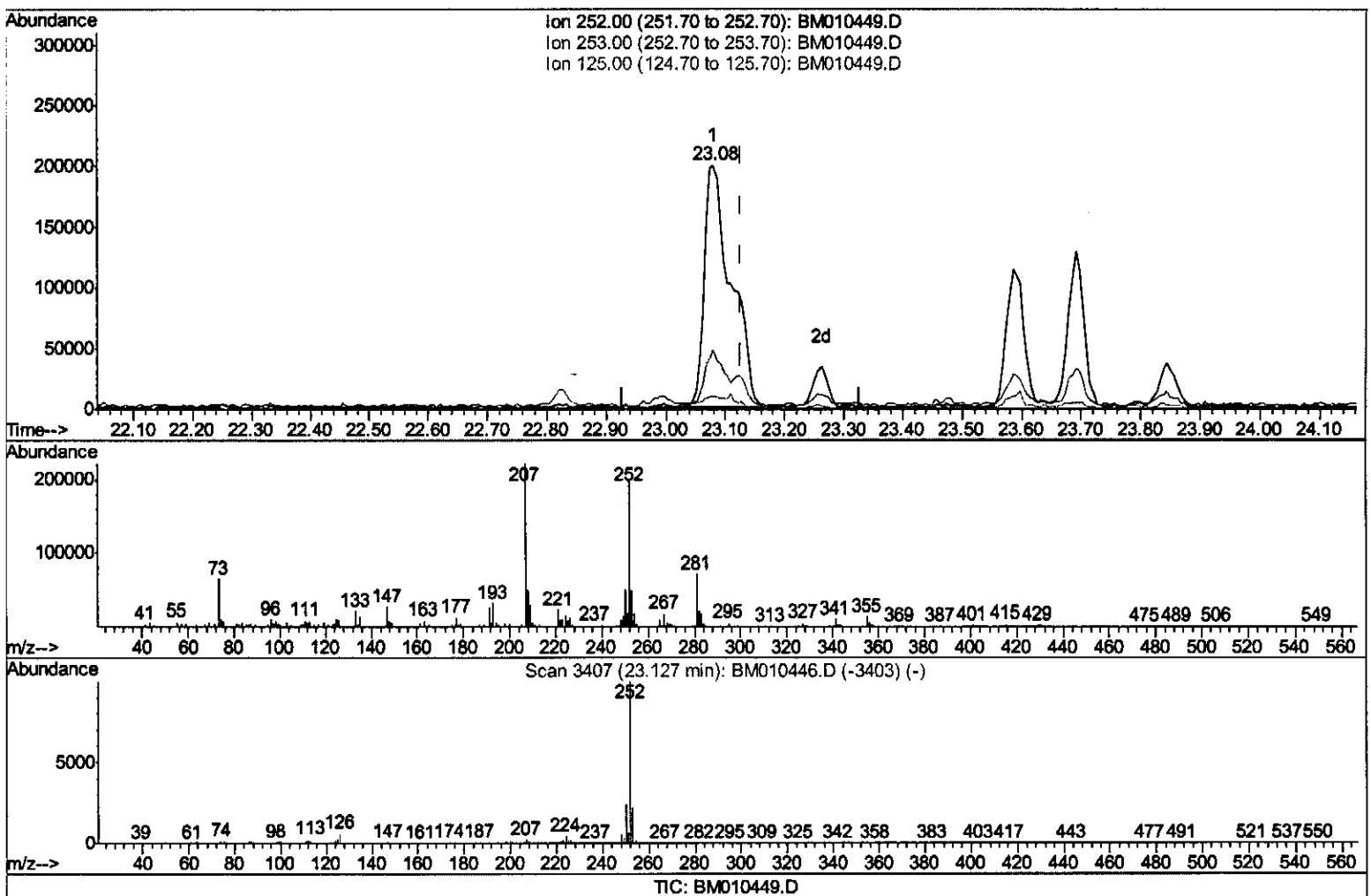
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Sample : I3521-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Manual Integrations
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mohammad
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Quant Time: Jun 10 01:59:41 2017
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jun 10 01:45:58 2017
Response via : Initial Calibration



(86) Benzo(k)fluoranthene

23.080min (-0.047) 6.54ng/ul

response 620974

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	24.65
125.00	4.30	5.29#
0.00	0.00	0.00

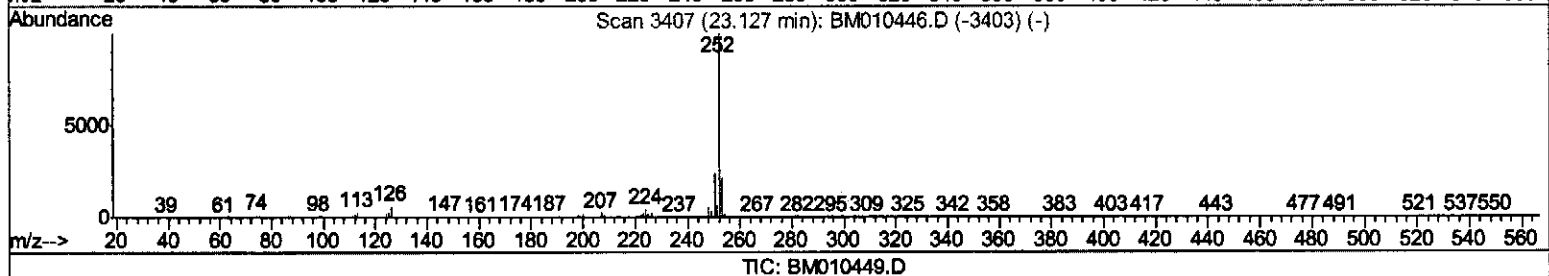
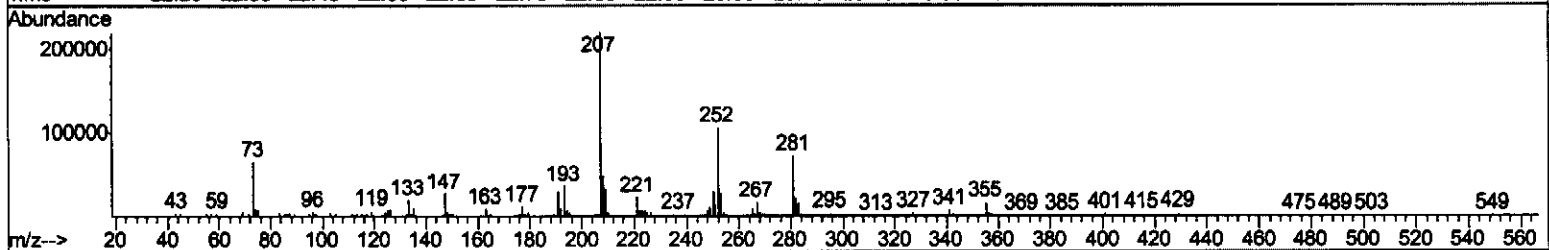
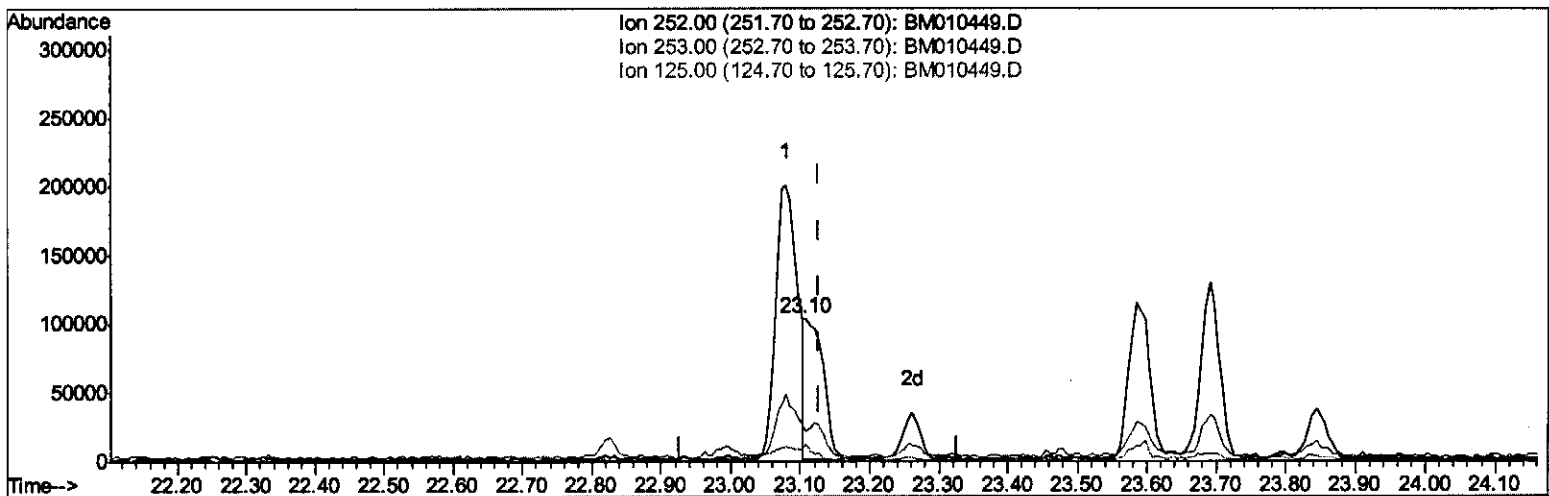
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Data File : BM010449.D
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Operator : SJ/MA
Sample : I3521-17
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
F5C60

Manual Integrations
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Quant Time: Jun 10 01:59:41 2017
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



(86) Benzo(k)fluoranthene

23.103min (-0.024) 1.95ng/ul m

response 185634

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	25.86#
125.00	4.30	7.98#
0.00	0.00	0.00

SJ
06/19/17

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
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 Operator : SJ/MA
 Sample : I3521-17
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	272494	20.00	ng/ul	0.00
18) Naphthalene-d8	10.71	136	992801	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.54	164	659974	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.29	188	1518261	20.00	ng/ul	0.00
75) Chrysene-d12	21.45	240	1664720	20.00	ng/ul	0.00
83) Perylene-d12	23.80	264	1708514	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.37	96	29100	4.38	ng/uL	0.00
5) Phenol-d5	7.09	99	467153	22.62	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.25	67	262541	22.62	ng/ul	0.00
9) 2-Chlorophenol-d4	7.45	132	396129	24.07	ng/ul	0.00
13) 4-Methylphenol-d8	8.63	113	380370	22.83	ng/ul	0.00
19) Nitrobenzene-d5	9.09	128	182708	25.12	ng/ul	0.00
22) 2-Nitrophenol-d4	9.80	143	230438	27.35	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.33	165	456810	26.40	ng/ul	0.00
29) 4-Chloroaniline-d4	10.87	131	459028	27.49	ng/ul	0.00
43) Dimethylphthalate-d6	13.96	166	1508018	27.50	ng/ul	0.00
46) Acenaphthylene-d8	14.24	160	1666724	26.06	ng/ul	0.00
51) 4-Nitrophenol-d4	14.78	143	134595	16.41	ng/ul	0.00
57) Fluorene-d10	15.53	176	1280013	26.56	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.66	200	215814	20.07	ng/ul	0.00
70) Anthracene-d10	17.39	188	1945482	26.33	ng/ul	0.00
76) Pyrene-d10	19.67	212	2180302	31.67	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.64	264	2111283	26.54	ng/ul	0.00

Target Compounds

					Qvalue	
6) Phenol	7.12	94	43279	2.044	ng/ul	93
44) Dimethylphthalate	14.00	163	725214	13.448	ng/ul	99
74) Fluoranthene	19.33	202	228603	2.294	ng/ul#	96
77) Pyrene	19.70	202	301438	3.514	ng/ul#	95
80) Benzo(a)anthracene	21.44	228	156587	1.668	ng/ul	96
82) Chrysene	21.49	228	208005m	2.451	ng/ul	} SJ 06/19/17
85) Benzo(b)fluoranthene	23.08	252	438934m	4.415	ng/ul	
86) Benzo(k)fluoranthene	23.10	252	185634m	1.955	ng/ul	
88) Benzo(a)pyrene	23.69	252	238800	2.423	ng/ul	94
89) Indeno(1,2,3-cd)pyrene	26.17	276	211712	1.700	ng/ul#	95
91) Benzo(g,h,i)perylene	26.92	276	166988	1.628	ng/ul	99

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