

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090716\
 Data File : BM007449.D
 Acq On : 07 Sep 2016 21:18
 Operator : UM/SJ
 Sample : H4666-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

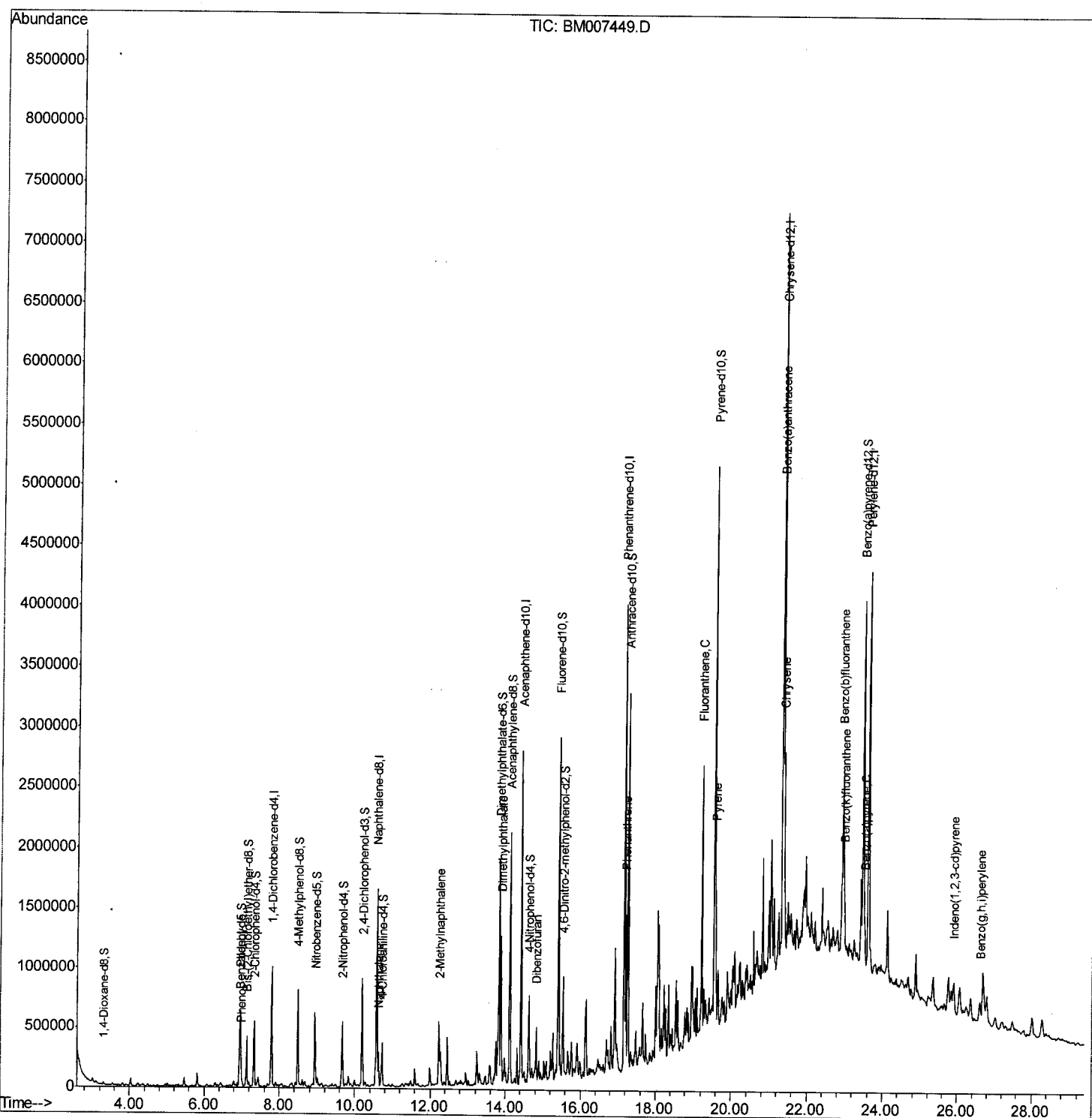
Instrument :
 BNA_M
 ClientSampleId :
 BD3A1

Manual Integrations
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Quant Time: Sep 08 04:49:57 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 08 03:10:54 2016
 Response via : Initial Calibration

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Quantitation Report (Qedit)

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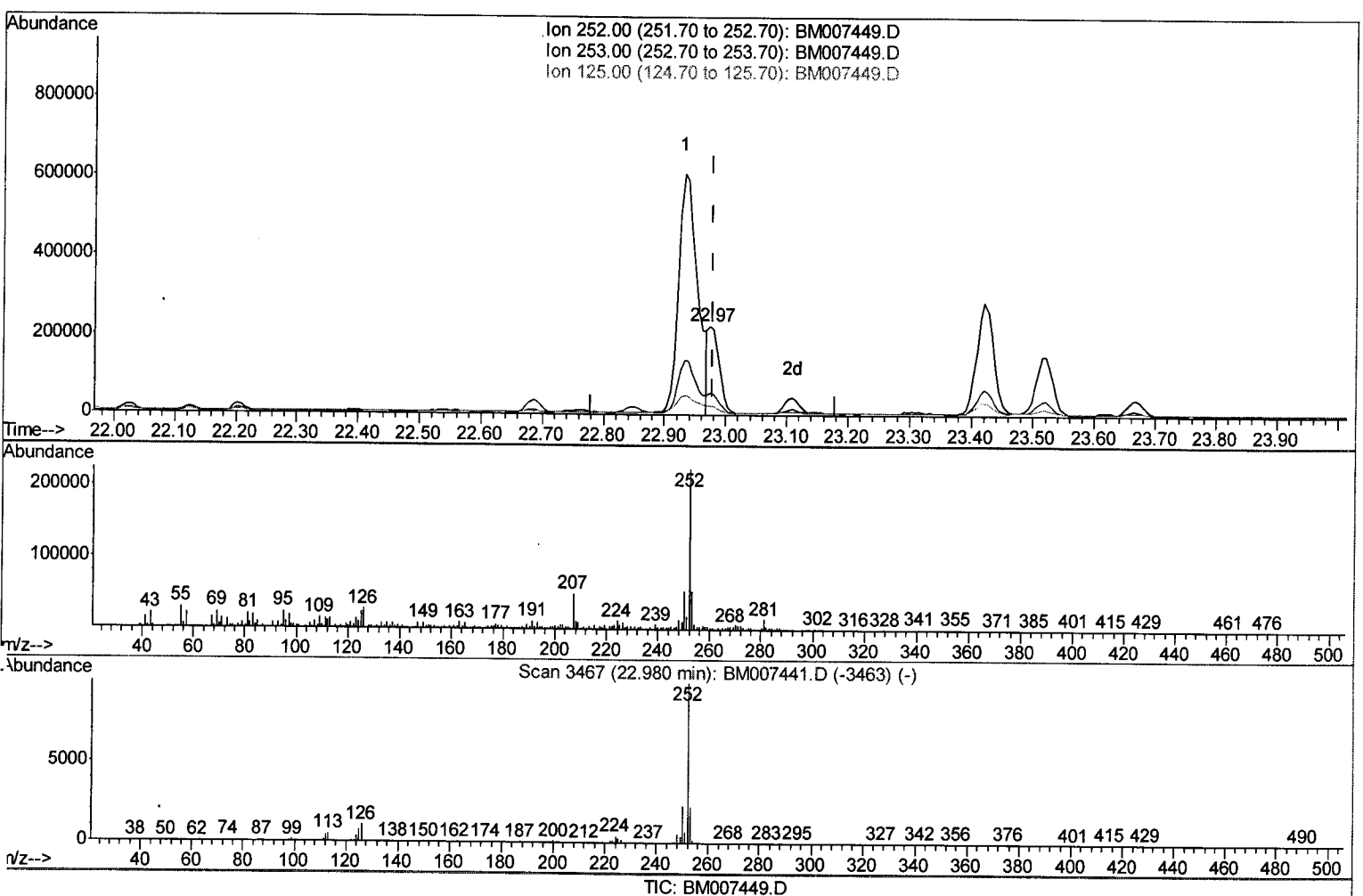
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Quant Time: Sep 08 04:43:49 2016
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(86) Benzo(k)fluoranthene

22.974min (-0.006) 1.97ng/ul m

response 275210

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	24.25
125.00	7.10	10.15#
0.00	0.00	0.00

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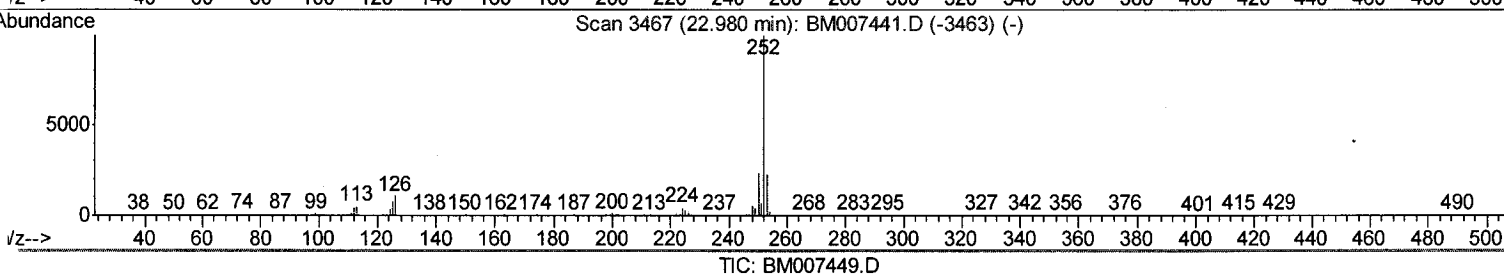
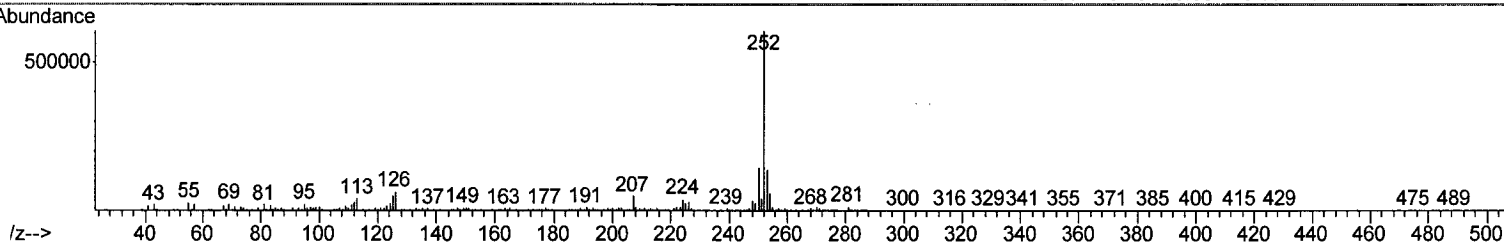
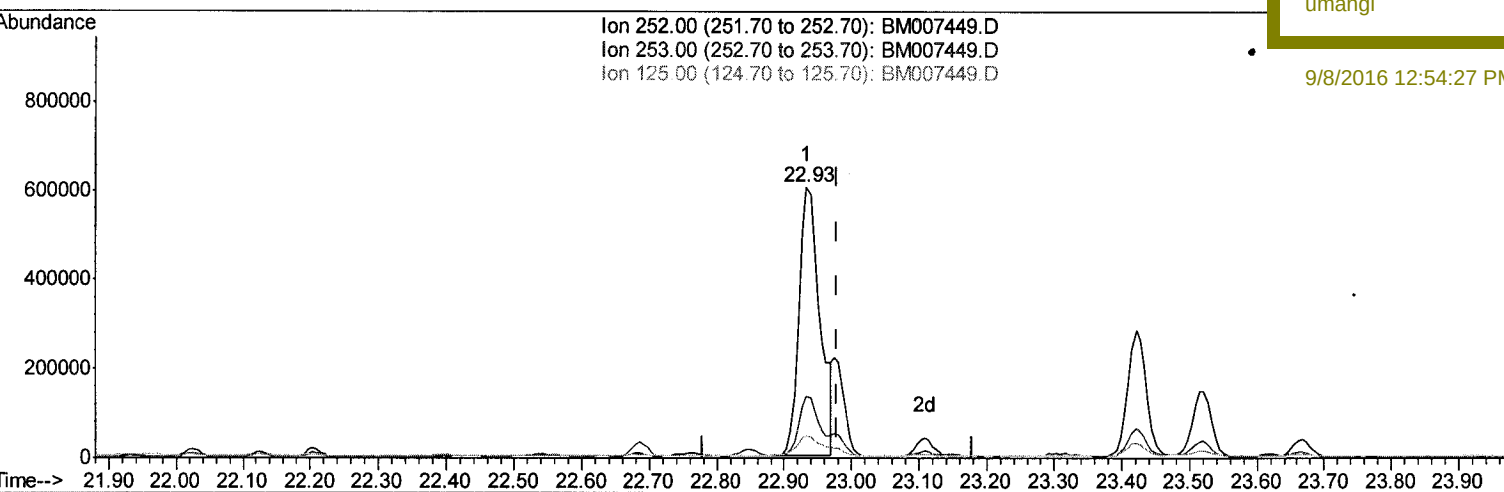
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(86) Benzo(k)fluoranthene

22.933min (-0.047) 9.29ng/ui

response 1296086

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	22.67
125.00	7.10	8.11
0.00	0.00	0.00

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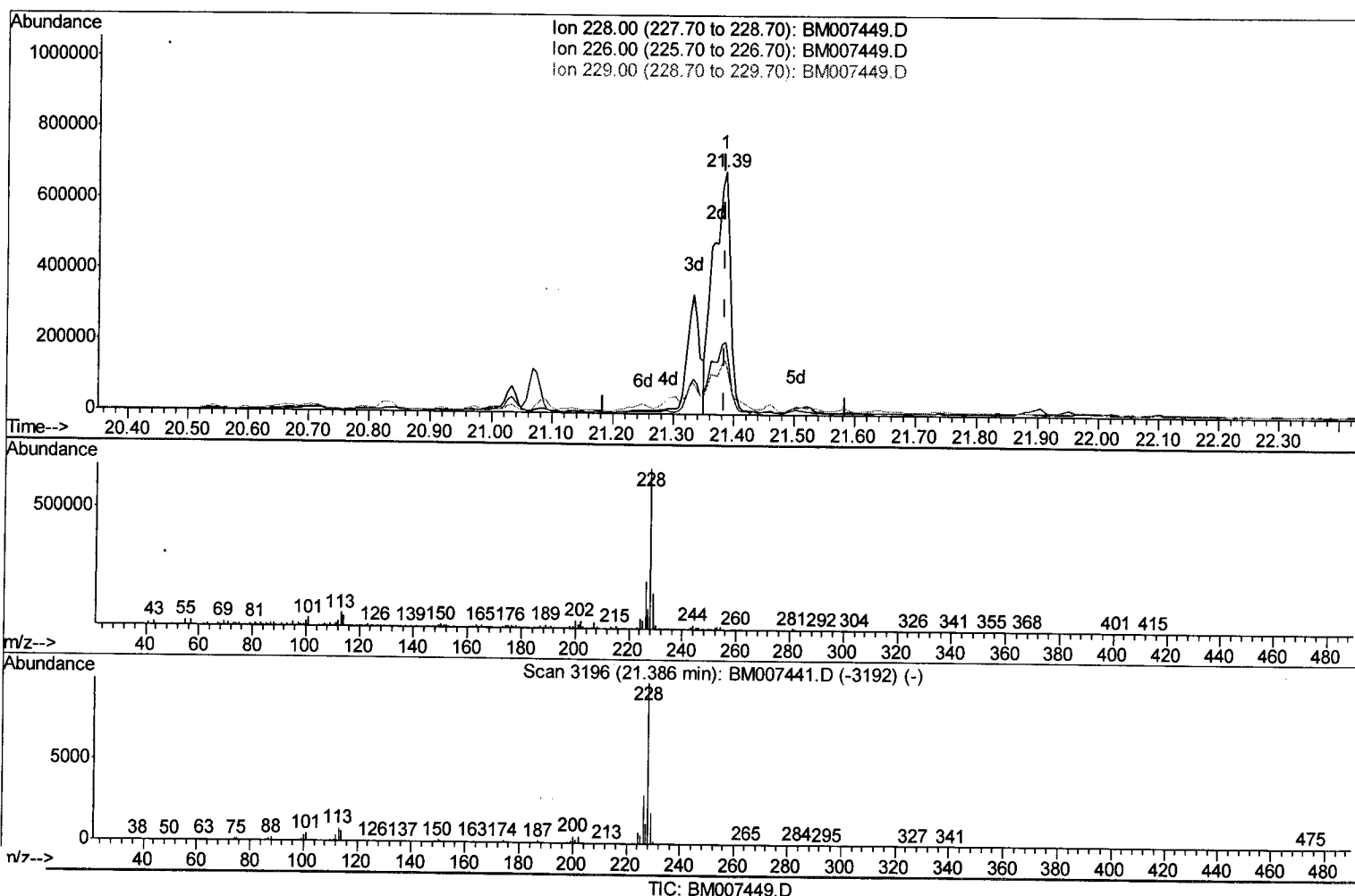
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(82) Chrysene

21.386min (+0.000) 9.09ng/ul m > 09/09/2016

response 1339622

Ion	Exp%	Act%
228.00	100	100
226.00	29.30	29.85
229.00	19.40	22.33
0.00	0.00	0.00

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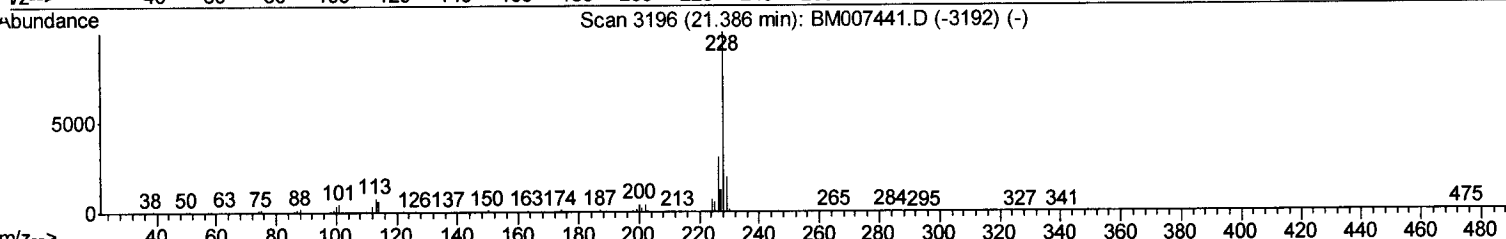
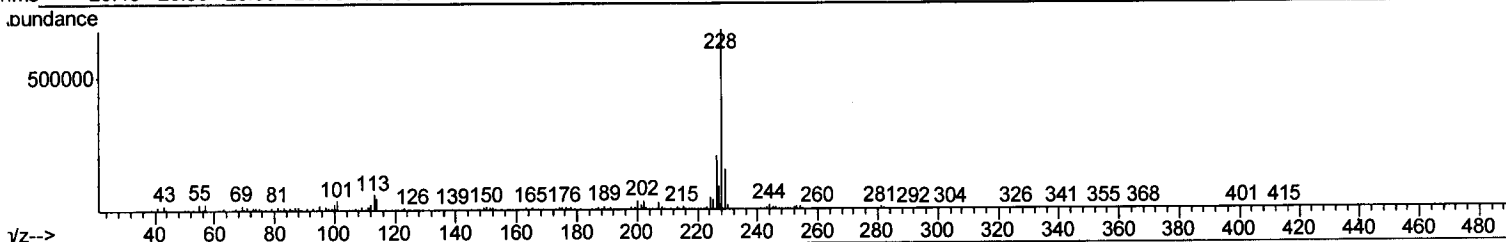
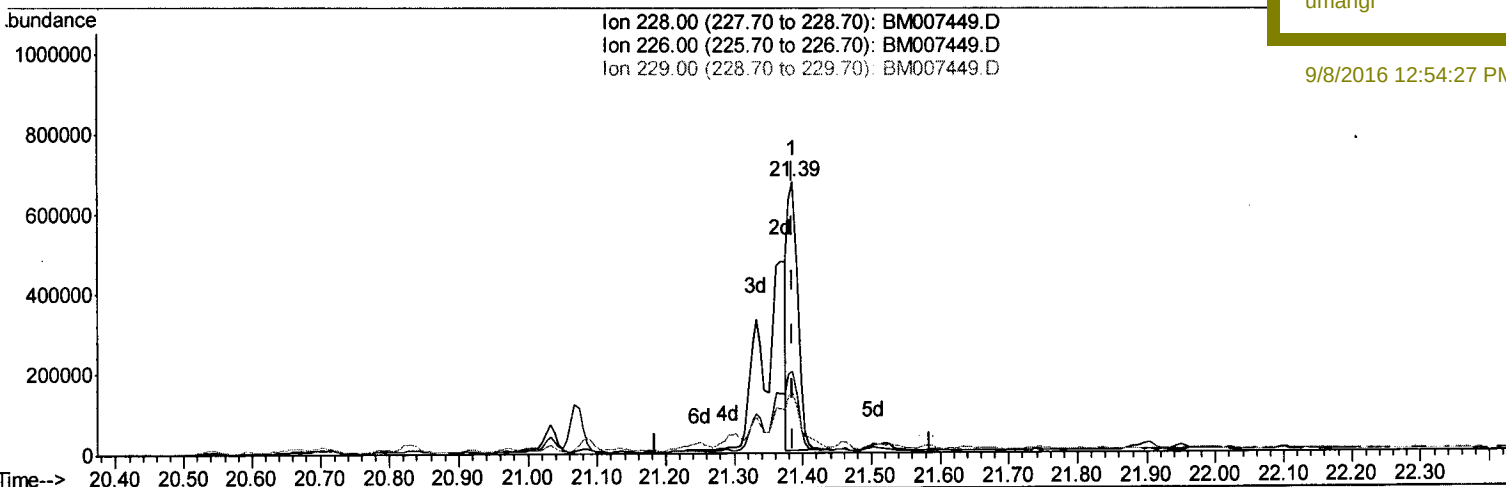
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TIC: BM007449.D

(82) Chrysene

21.386min (+0.000) 4.78ng/ul

response 704514

Ion	Exp%	Act%
228.00	100	100
226.00	29.30	29.85
229.00	19.40	22.33
0.00	0.00	0.00

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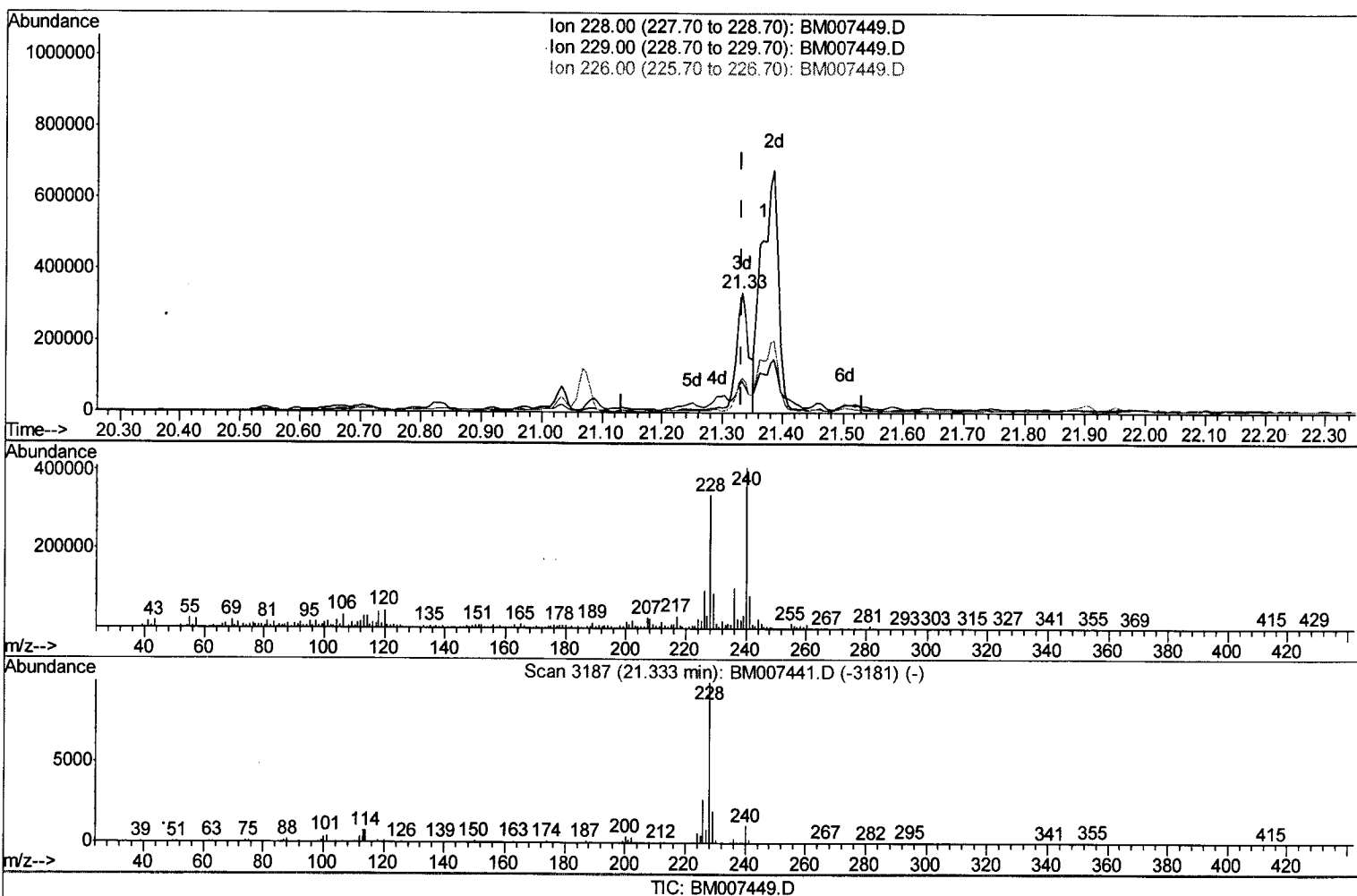
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(80) Benzo(a)anthracene

21.333min (+0.000) 3.23ng/ul m

response 525952

Ion	Exp%	Act%
228.00	100	100
229.00	19.50	26.54#
226.00	27.30	29.41
0.00	0.00	0.00

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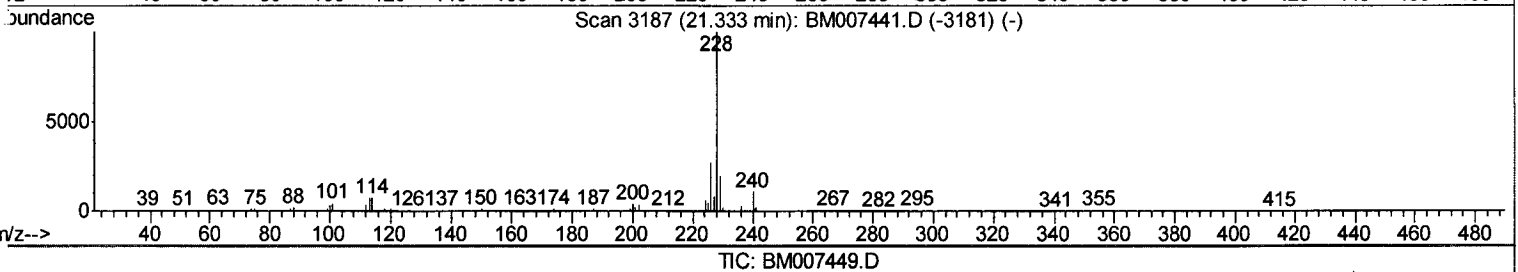
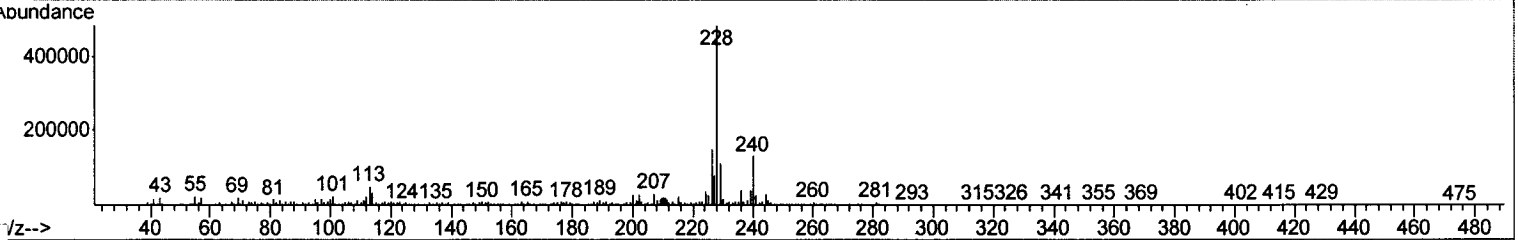
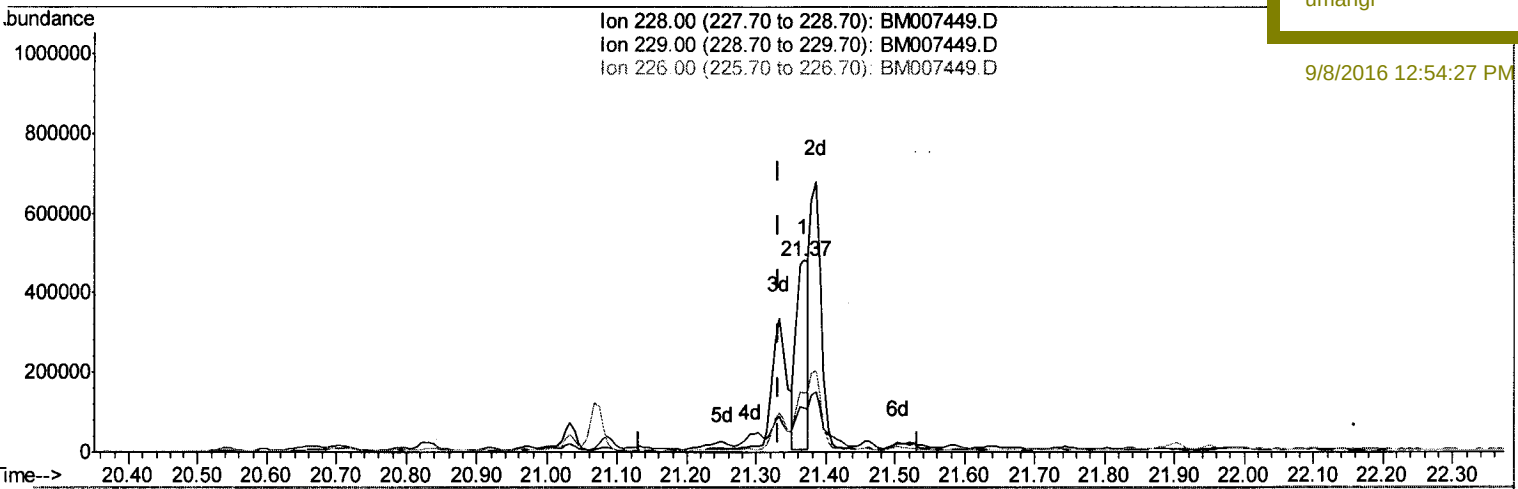
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Instrument :
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ClientSampleId :
BD3A1

Manual Integrations
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(80) Benzo(a)anthracene

21.368min (+0.036) 3.68ng/ul

response 599820

Ion	Exp%	Act%
228.00	100	100
229.00	19.50	23.31
226.00	27.30	30.83
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	265505	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	1318101	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.42	164	901796	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	2229581	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	2761415	20.00	ng/ul	0.00
83) Perylene-d12	23.62	264	2551931	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.30	96	9686	1.86	ng/uL	0.00
5) Phenol-d5	6.96	99	324477	12.94	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	176462	13.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	241734	13.04	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	298793	13.66	ng/ul	0.00
19) Nitrobenzene-d5	8.95	128	139576	14.42	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	166430	14.73	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	328351	14.60	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	186540	6.65	ng/ul	0.00
43) Dimethylphthalate-d6	13.83	166	1208677	15.54	ng/ul	0.00
46) Acenaphthylene-d8	14.12	160	1452312	16.10	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	168215	11.45	ng/ul	0.00
57) Fluorene-d10	15.42	176	1088351	16.17	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	193789	13.82	ng/ul	0.00
70) Anthracene-d10	17.26	188	1769784	16.99	ng/ul	0.00
76) Pyrene-d10	19.56	212	2140015	18.53	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.47	264	1940388	16.51	ng/ul	0.00

Target Compounds

					Qvalue
4) Benzaldehyde	6.94	77	95653	6.69 ng/ul	97
6) Phenol	6.99	94	28479	1.09 ng/ul	91
28) Naphthalene	10.63	128	223251	3.41 ng/ul	98
34) 2-Methylnaphthalene	12.24	142	223318	4.16 ng/ul	100
44) Dimethylphthalate	13.88	163	757668	9.78 ng/ul	100
53) Dibenzofuran	14.82	168	164474	1.82 ng/ul	96
69) Phenanthrene	17.21	178	786242	6.57 ng/ul	99
74) Fluoranthene	19.22	202	1111275	7.40 ng/ul	99
77) Pyrene	19.59	202	747489	5.05 ng/ul	100
80) Benzo(a)anthracene	21.33	228	525952m	3.23 ng/ul	
82) Chrysene	21.39	228	1339622m	9.09 ng/ul	
85) Benzo(b)fluoranthene	22.93	252	1296086	8.59 ng/ul	97
86) Benzo(k)fluoranthene	22.97	252	275210m	1.97 ng/ul	
88) Benzo(a)pyrene	23.52	252	284957	1.97 ng/ul#	93
89) Indeno(1,2,3-cd)pyrene	25.90	276	243247	1.37 ng/ul	98
91) Benzo(g,h,i)perylene	26.61	276	199450	1.33 ng/ul	98

} U.M.
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(#) = qualifier out of range (m) = manual integration (+) = signals summed