

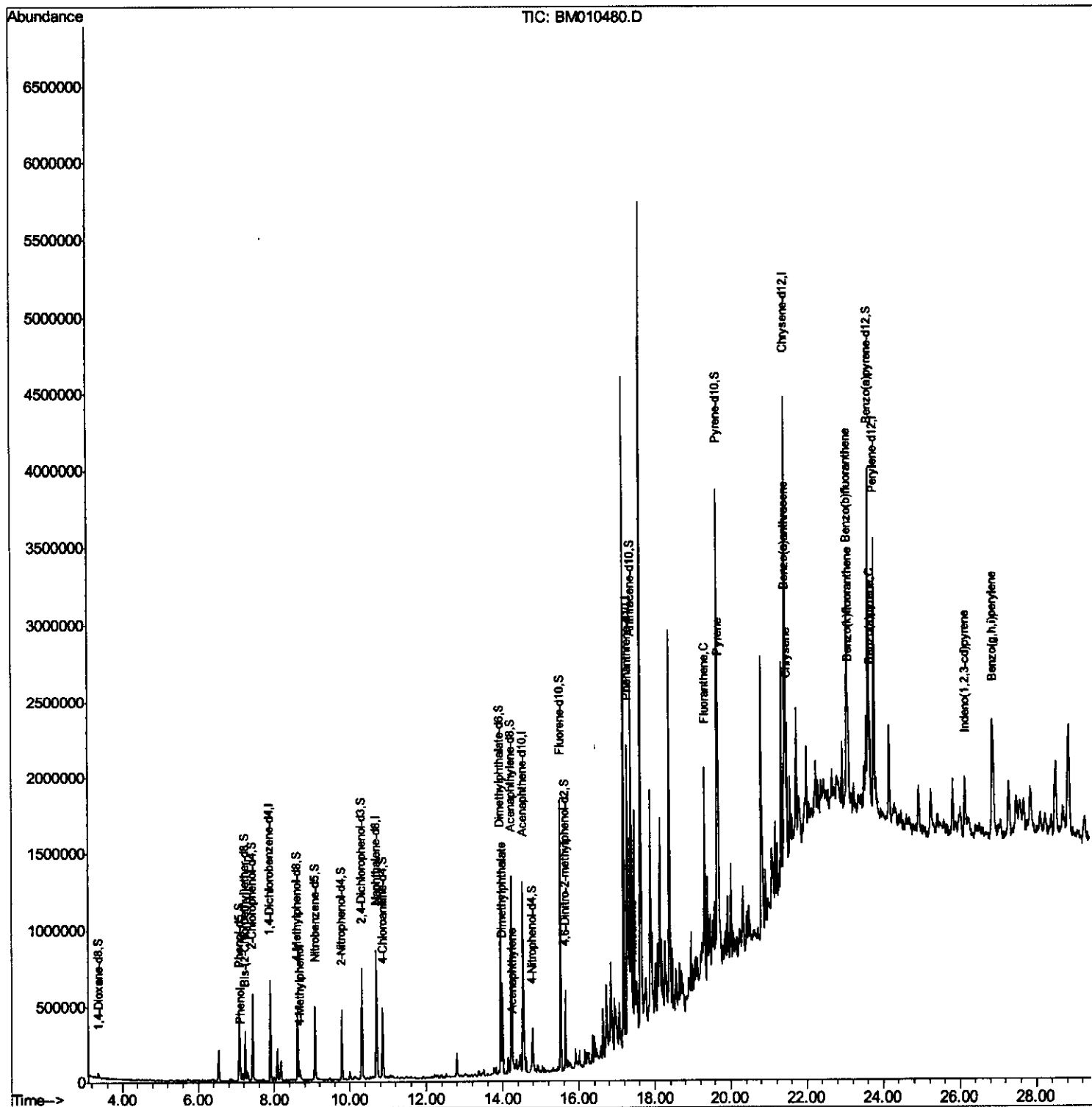
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061017\
Data File : BM010480.D
Acq On : 10 Jun 2017 11:58
Operator : SJ/MA
Sample : I3521-16
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
Client Sample ID :
F5C36

Manual Integrations
APPROVED

mohammad
6/12/2017 3:40:30 PM

Quant Time: Jun 12 05:36:03 2017
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 12 04:15:09 2017
Response via : Initial Calibration



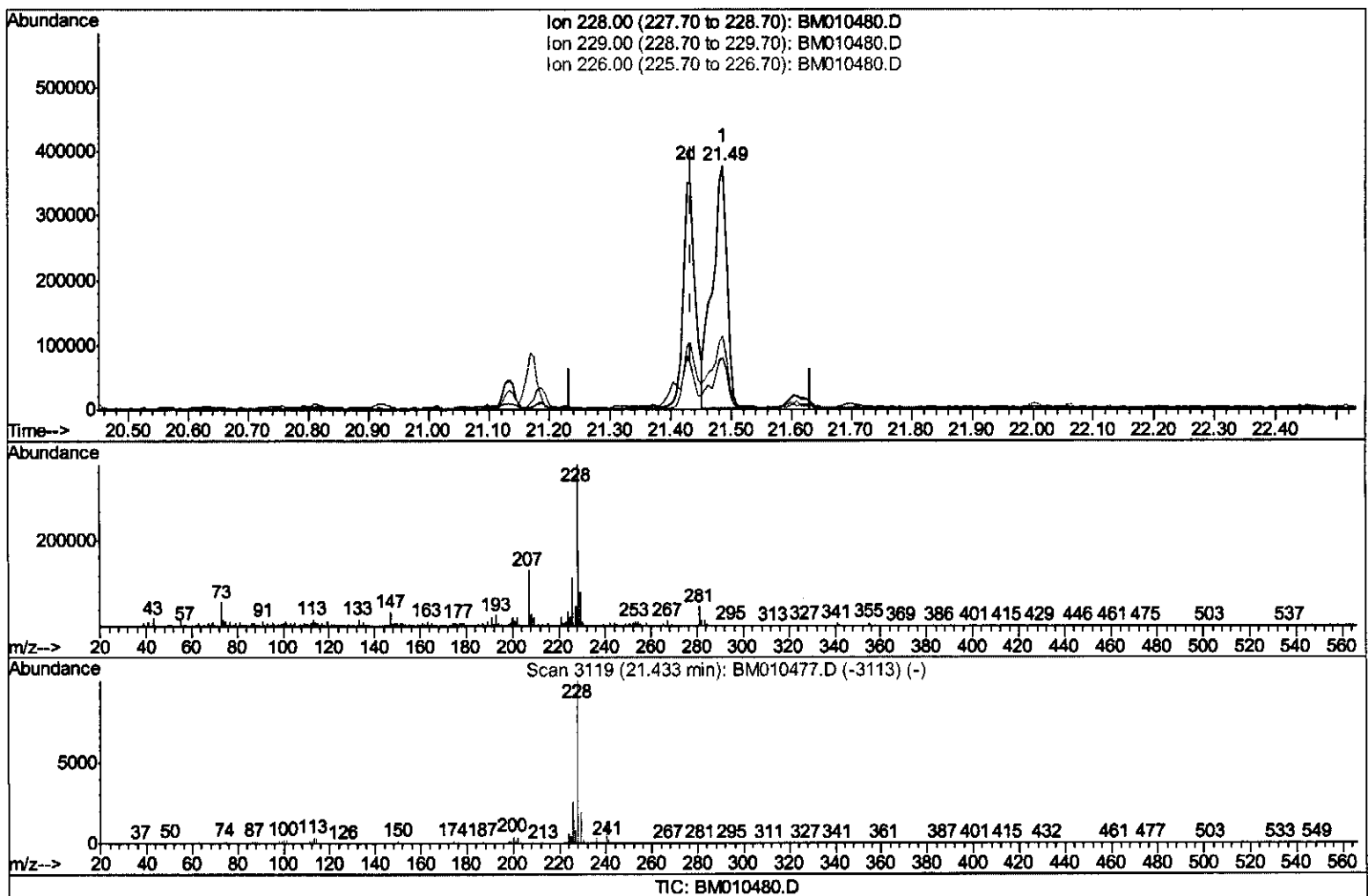
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(80) Benzo(a)anthracene

21.486min (+0.053) 8.34ng/ul

response 619096

Ion	Exp%	Act%
228.00	100	100
229.00	19.20	21.45
226.00	26.70	30.34
0.00	0.00	0.00

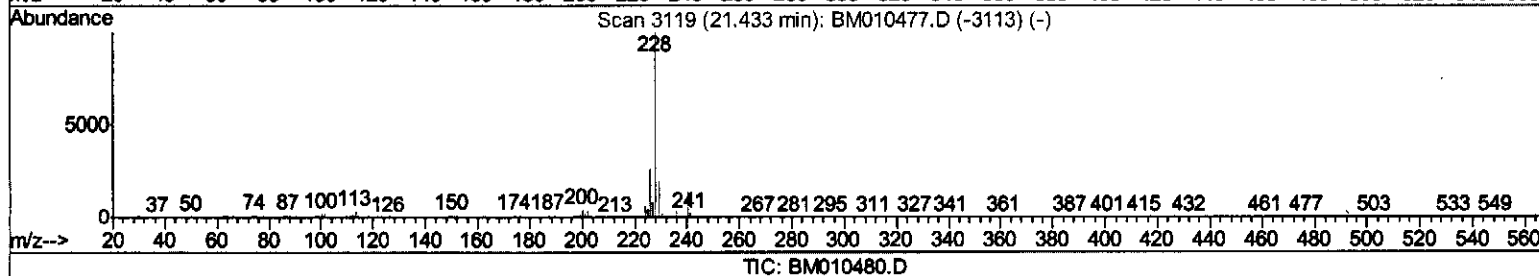
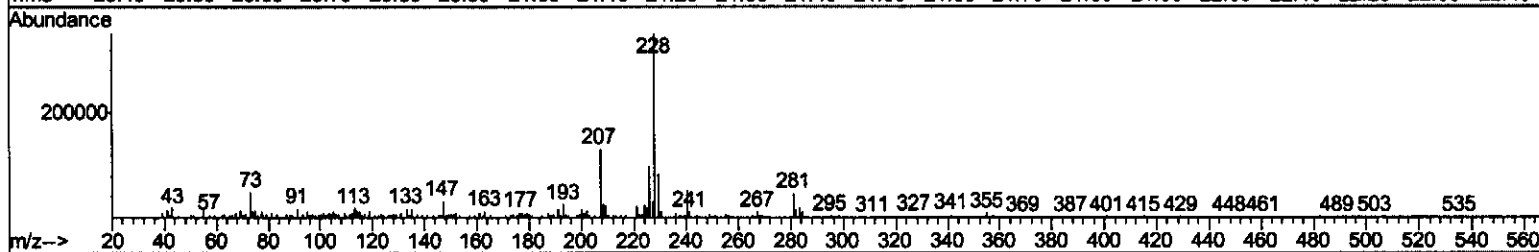
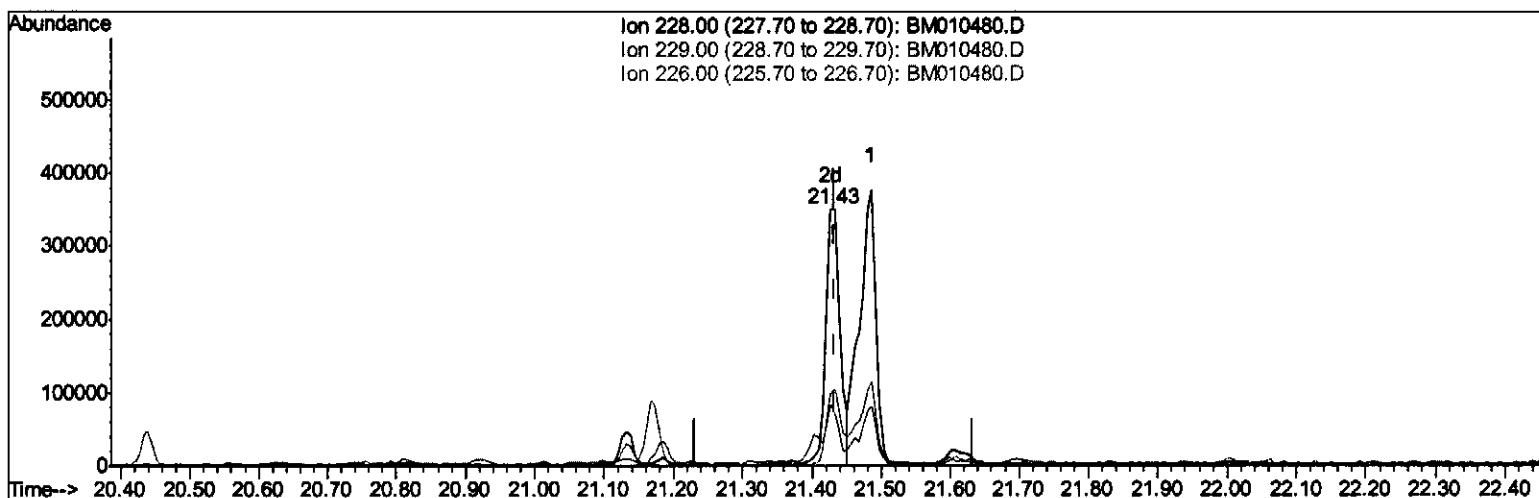
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(80) Benzo(a)anthracene

21.427min (-0.006) 6.80ng/ul m

response 504419

Ion	Exp%	Act%
228.00	100	100
229.00	19.20	23.83#
226.00	26.70	28.28
0.00	0.00	0.00

SJ 06/19/17

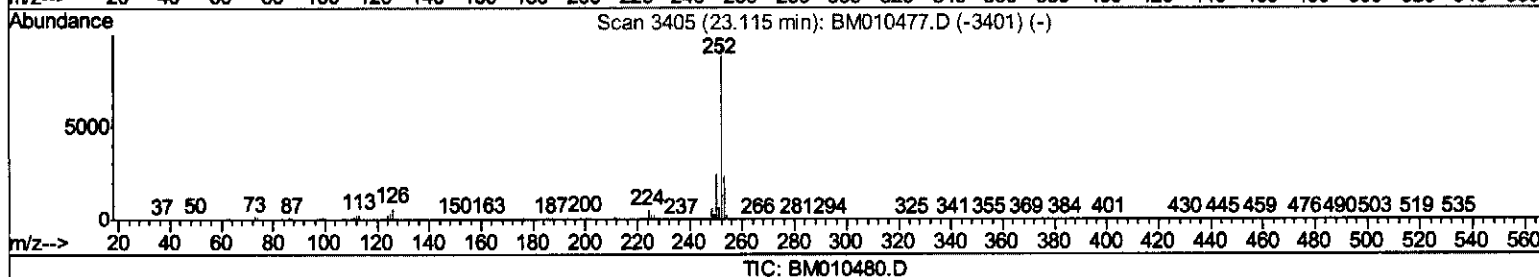
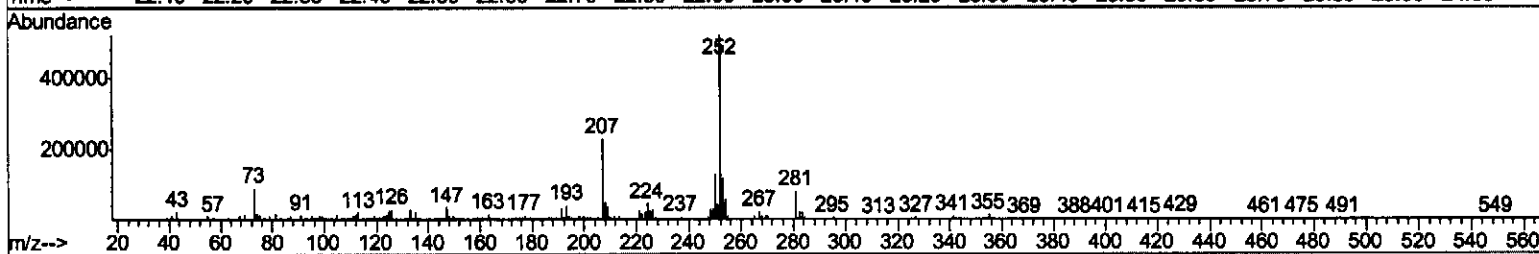
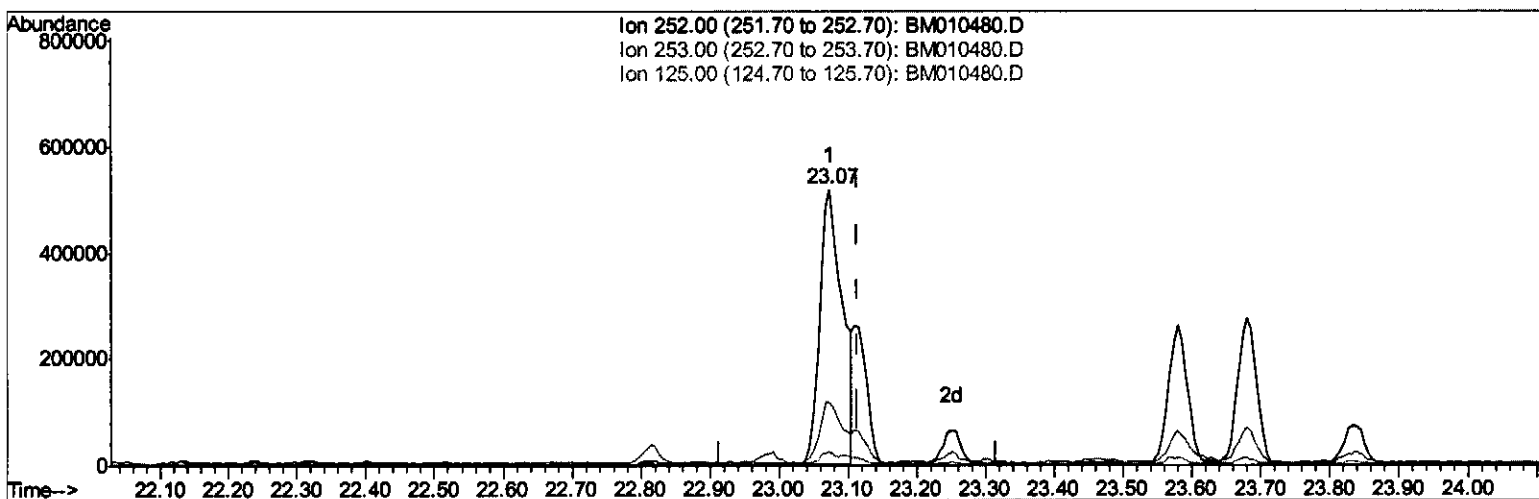
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Manual Integrations
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(86) Benzo(k)fluoranthene

23.074min (-0.041) 15.75ng/ul

response 1165682

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	22.85
125.00	4.30	4.95
0.00	0.00	0.00

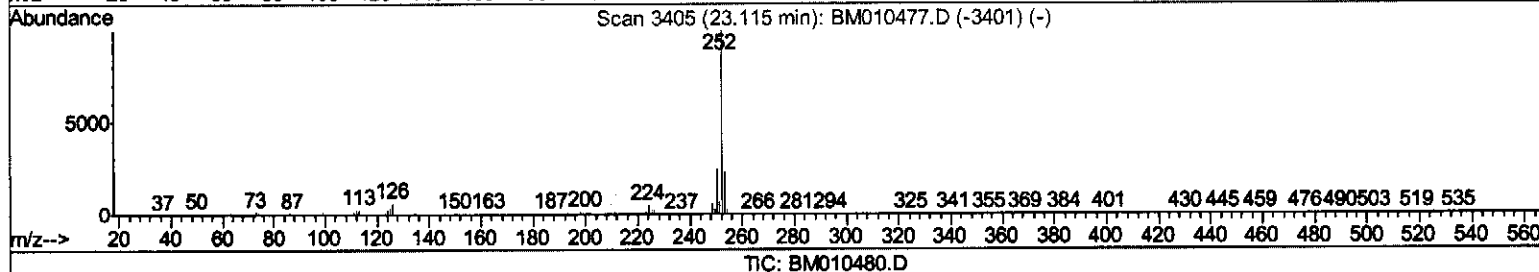
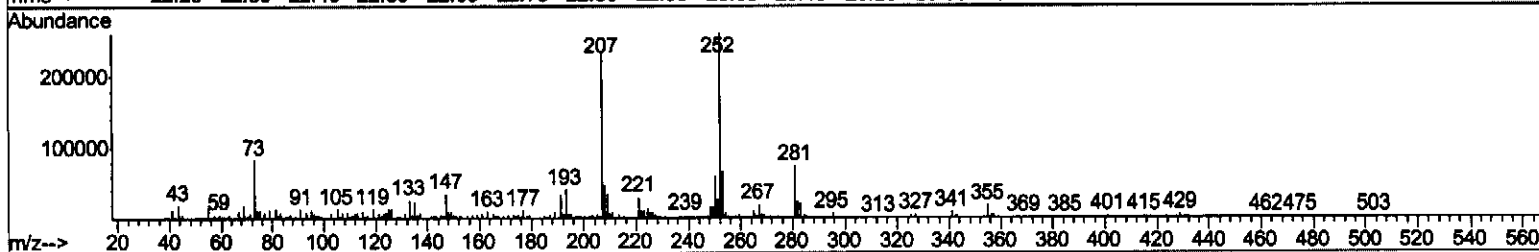
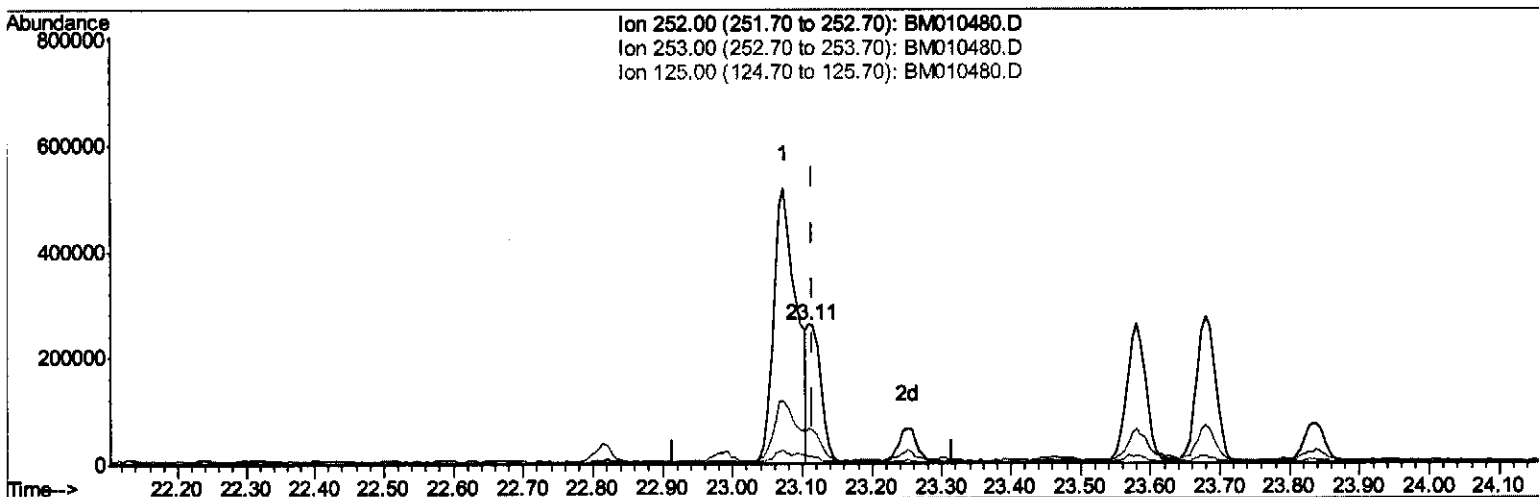
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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Manual Integrations
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Response via : Initial Calibration



(86) Benzo(k)fluoranthene

23.109min (-0.006) 4.82ng/ul m

response 357117

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	25.21
125.00	4.30	5.33#
0.00	0.00	0.00

SS 06/19/17

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	149112	20.00	ng/ul	0.00
18) Naphthalene-d8	10.70	136	545970	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.53	164	379391	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.28	188	926547	20.00	ng/ul	0.00
75) Chrysene-d12	21.44	240	1315656	20.00	ng/ul	0.00
83) Perylene-d12	23.79	264	1331483	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.36	96	14991	4.12	ng/uL	0.00
5) Phenol-d5	7.08	99	197104	17.44	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.24	67	124884	19.67	ng/ul	0.00
9) 2-Chlorophenol-d4	7.43	132	184472	20.48	ng/ul	0.00
13) 4-Methylphenol-d8	8.62	113	148257	16.26	ng/ul	0.00
19) Nitrobenzene-d5	9.09	128	86366	21.60	ng/ul	0.00
22) 2-Nitrophenol-d4	9.79	143	108889	23.50	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.33	165	219668	23.08	ng/ul	0.00
29) 4-Chloroaniline-d4	10.87	131	134423	14.64	ng/ul	0.00
43) Dimethylphthalate-d6	13.95	166	742010	23.54	ng/ul	0.00
46) Acenaphthylene-d8	14.23	160	819547	22.29	ng/ul	0.00
51) 4-Nitrophenol-d4	14.78	143	63451	13.45	ng/ul	0.00
57) Fluorene-d10	15.52	176	648287	23.40	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.65	200	114651	17.47	ng/ul	0.00
70) Anthracene-d10	17.38	188	1060147	23.51	ng/ul	0.00
76) Pyrene-d10	19.67	212	1415634	26.02	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.63	264	1496262	24.14	ng/ul	0.00

Target Compounds

					Qvalue
6) Phenol	7.11	94	22090	1.907 ng/ul#	93
16) 4-Methylphenol	8.67	108	20504	2.211 ng/ul	87
44) Dimethylphthalate	13.99	163	320360	10.334 ng/ul	98
47) Acenaphthylene	14.26	152	38345	1.067 ng/ul#	89
69) Phenanthrene	17.32	178	164729	3.330 ng/ul	95
71) Anthracene	17.42	178	54181	1.049 ng/ul#	95
74) Fluoranthene	19.33	202	778121	12.792 ng/ul#	95
77) Pyrene	19.70	202	811245	11.966 ng/ul#	94
80) Benzo(a)anthracene	21.43	228	504419m	6.799 ng/ul	94
82) Chrysene	21.49	228	619096	9.231 ng/ul	97
85) Benzo(b)fluoranthene	23.07	252	1165682	15.046 ng/ul	99
86) Benzo(k)fluoranthene	23.11	252	357117m	4.825 ng/ul	99
88) Benzo(a)pyrene	23.68	252	523675	6.819 ng/ul	95
89) Indeno(1,2,3-cd)pyrene	26.16	276	395443	4.076 ng/ul	97
91) Benzo(g,h,i)perylene	26.90	276	308021	3.853 ng/ul	99

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