

Quantitation Report (Qedit)

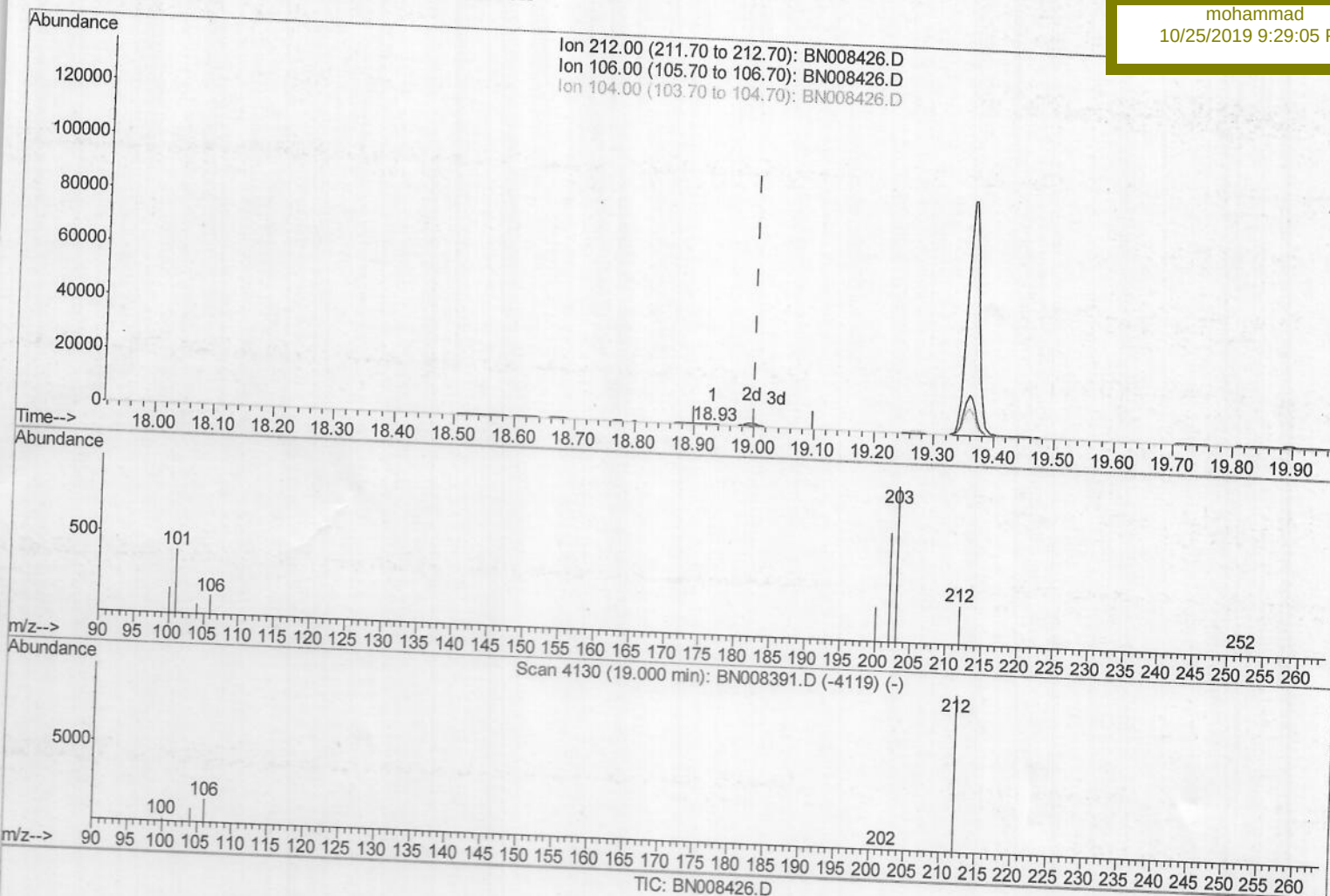
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102419\
 Data File : BN008426.D
 Acq On : 25 Oct 2019 14:39
 Operator : JU
 Sample : K5236-17DL 5X
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Quant Time: Oct 25 15:18:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 25 13:14:06 2019
 Response via : Initial Calibration

Instrument :
 BNA_N
 ClientSampleId :
 BEPN0DL

Manual Integrations
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(14) Fluoranthene-d10 (SURR)

18.930min (-0.070) 0.00ng/ul

response 121

Ion	Exp%	Act%
212.00	100	100
106.00	12.70	12.40
104.00	7.20	6.61
0.00	0.00	0.00

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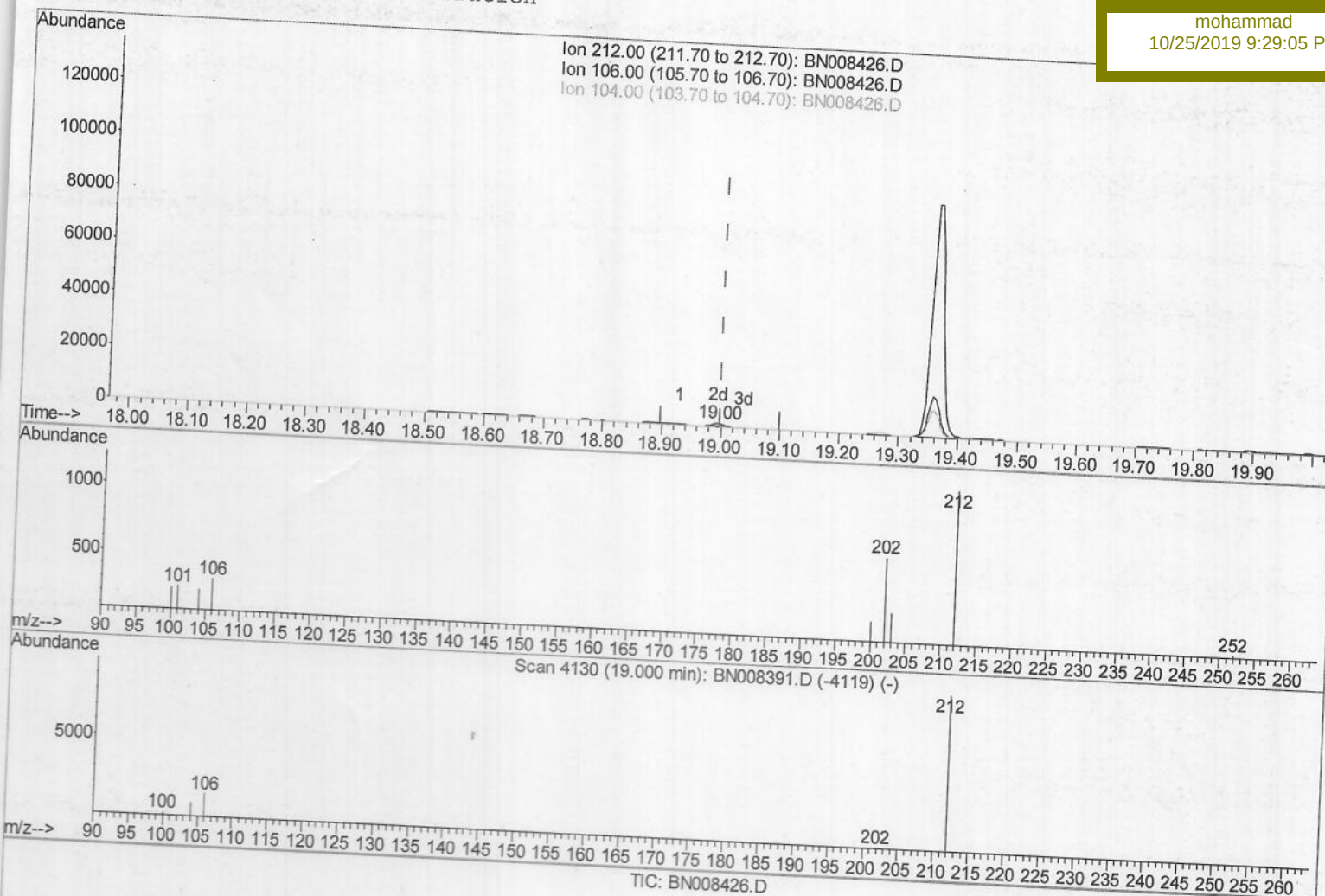
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(14) Fluoranthene-d10 (SURR)

18.995min (-0.005) 0.04ng/ul m →

response 1877

Ion	Exp%	Act%
212.00	100	100
106.00	12.70	0.80#
104.00	7.20	0.43#
0.00	0.00	0.00

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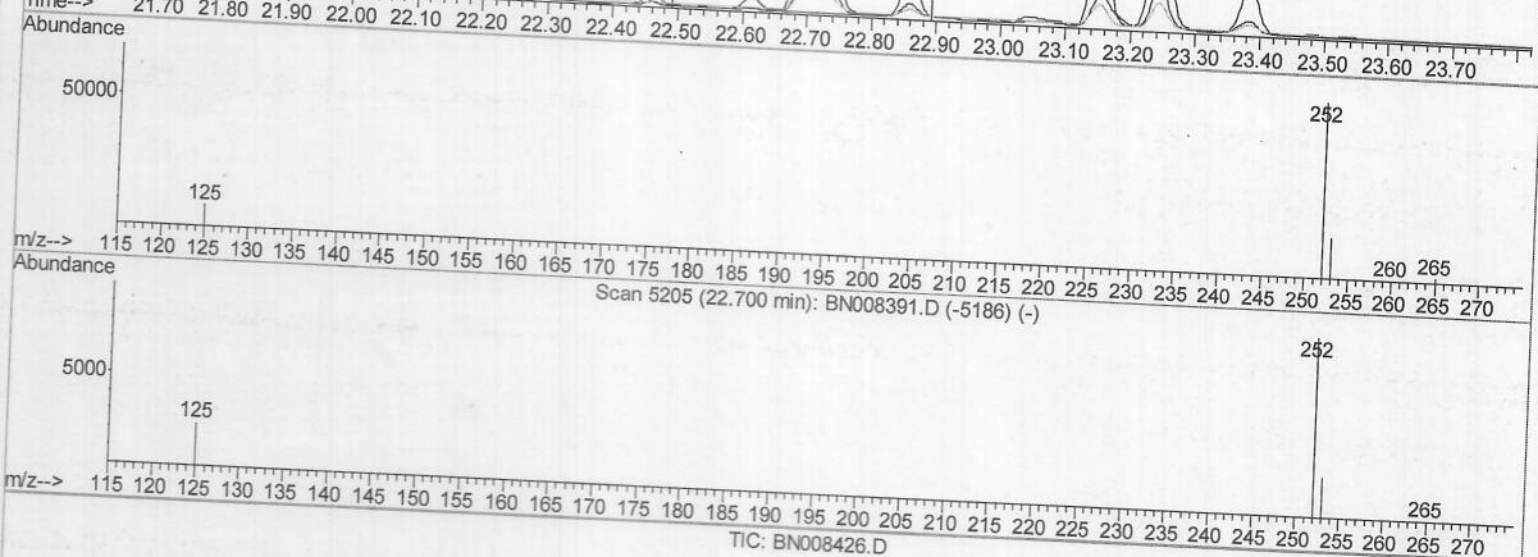
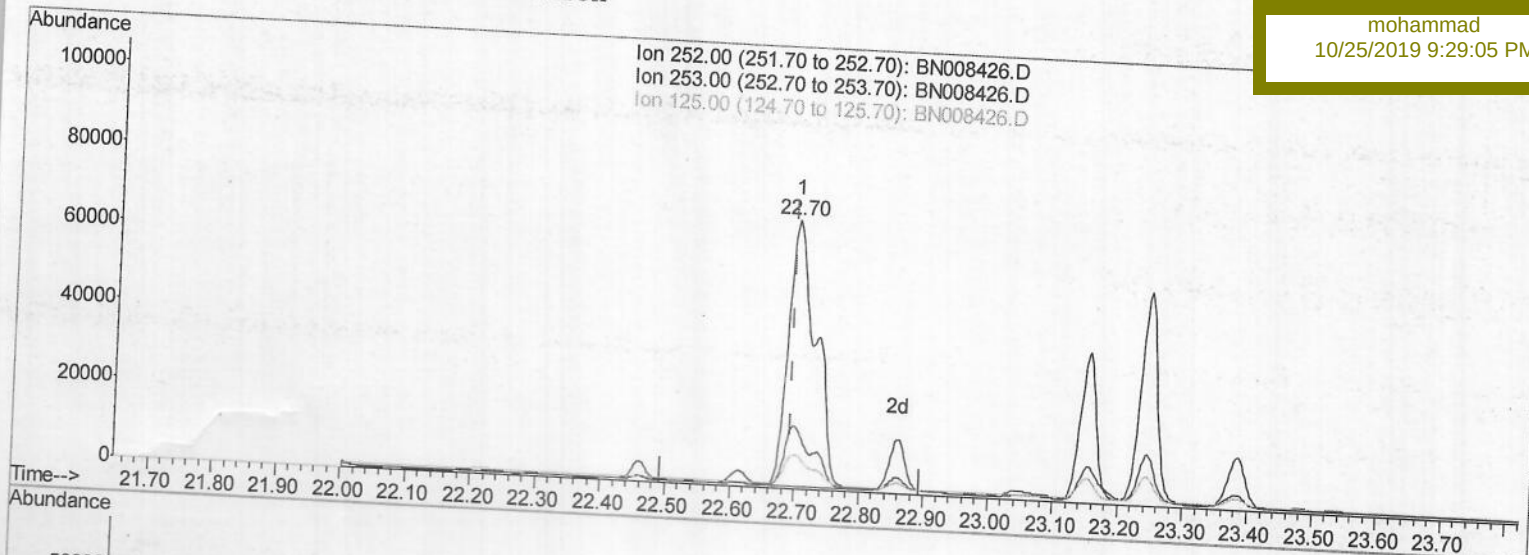
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(21) Benzo(b)fluoranthene
 22.697min (+0.003) 6.22ng/ul
 response 198018

Ion	Exp%	Act%
252.00	100	100
253.00	26.70	23.28
125.00	16.20	12.52
0.00	0.00	0.00

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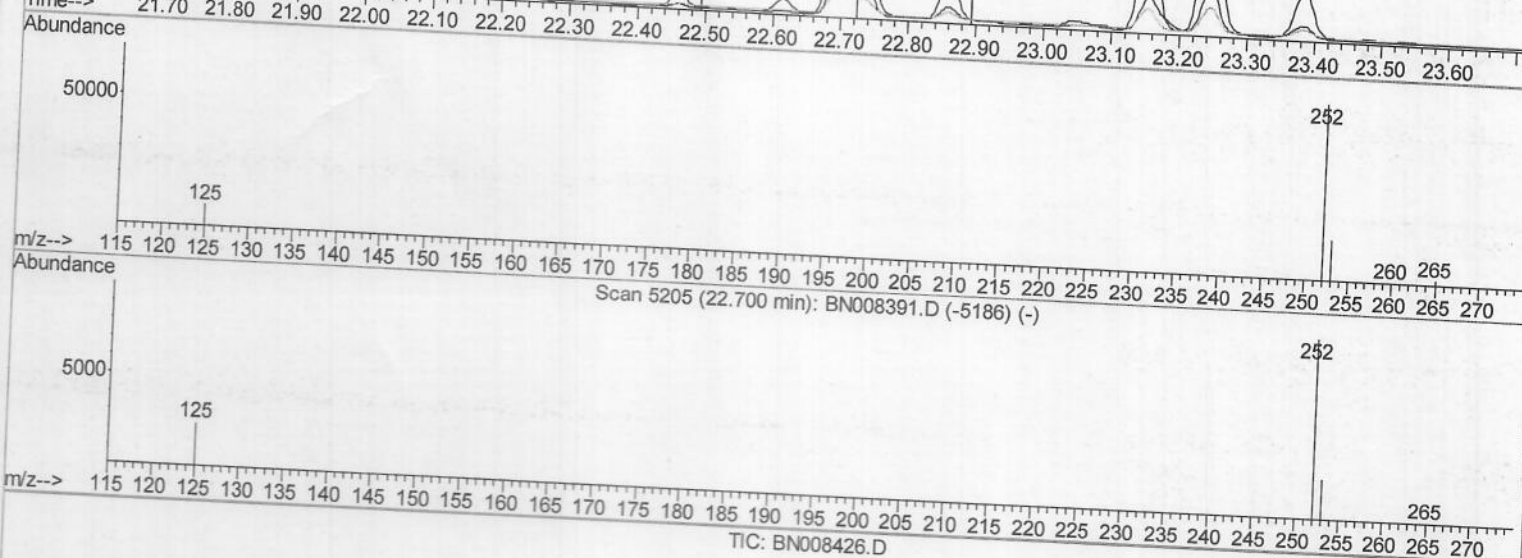
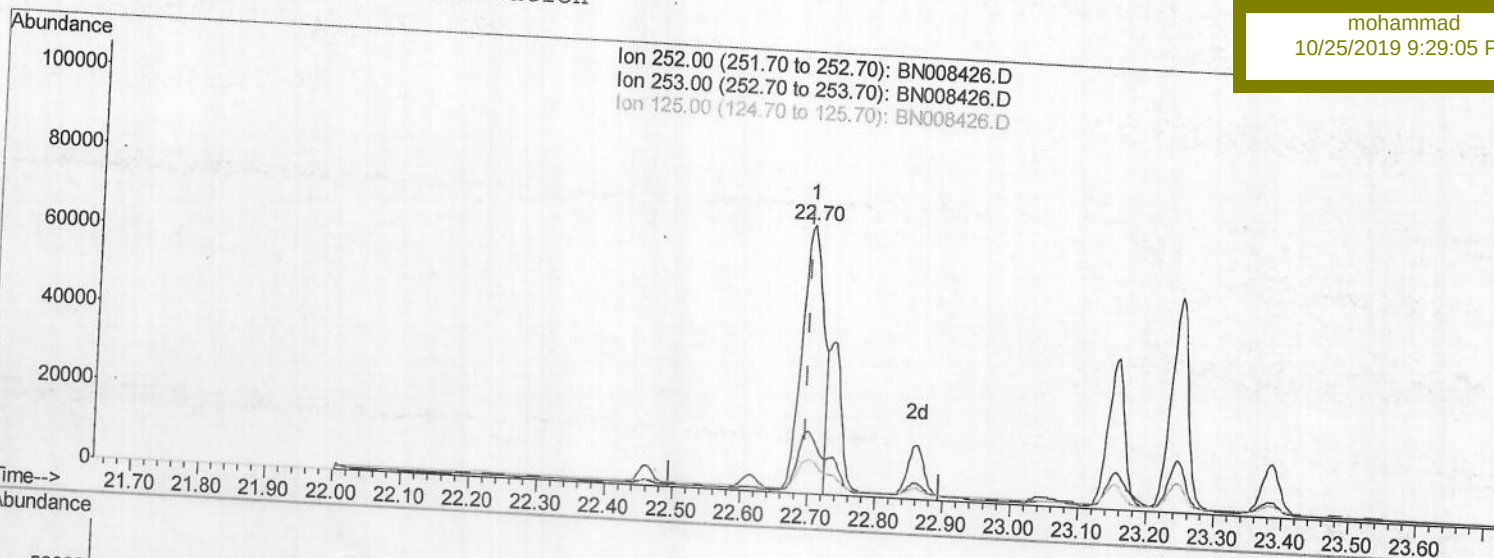
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(21) Benzo(b)fluoranthene

22.697min (+0.003) 4.60ng/ul m →

response 146394

Ion	Exp%	Act%
252.00	100	100
253.00	26.70	23.28
125.00	16.20	12.52
0.00	0.00	0.00

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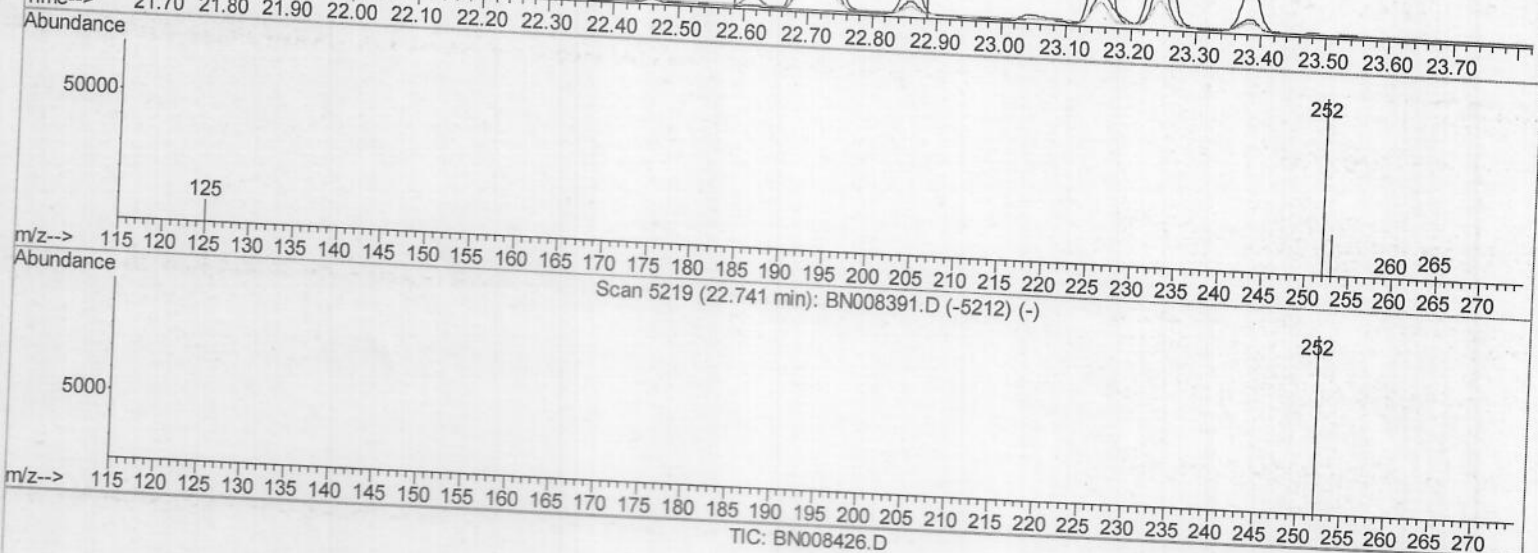
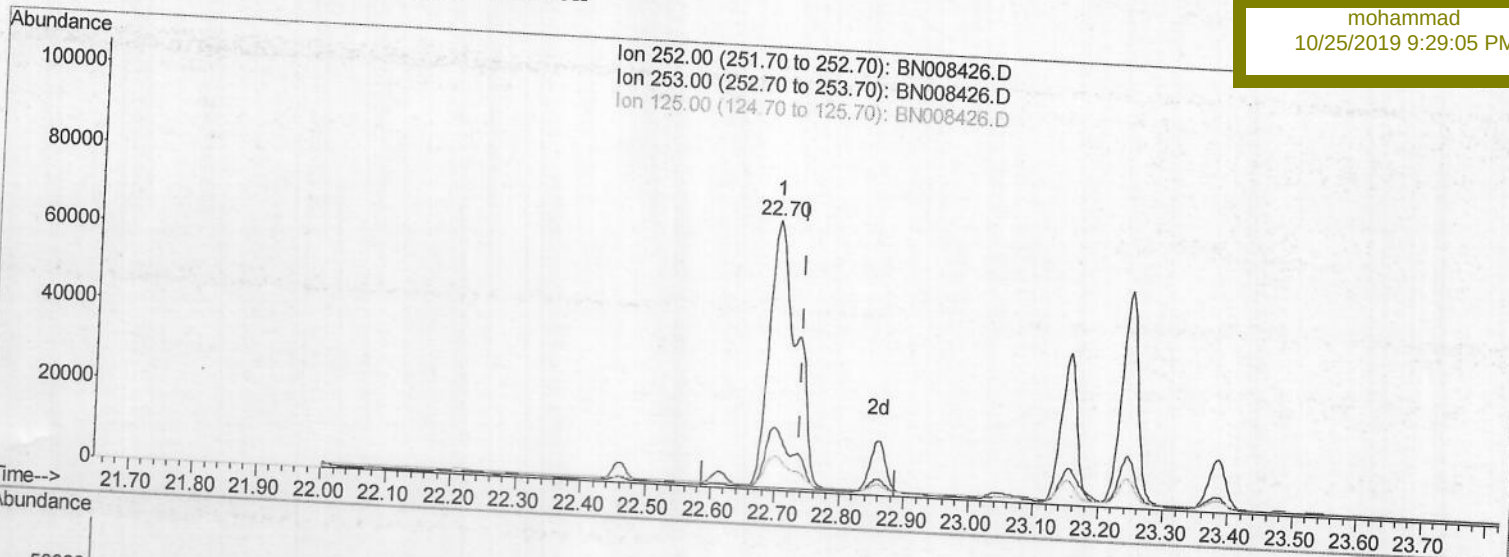
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(22) Benzo(k)fluoranthene

22.697min (-0.041) 5.49ng/ul

response 198018

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	23.28
125.00	15.40	12.52
0.00	0.00	0.00

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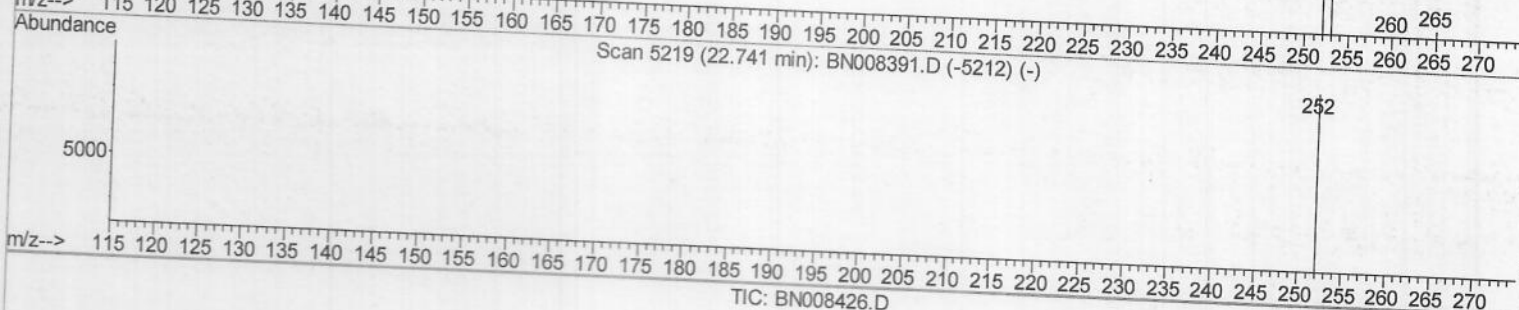
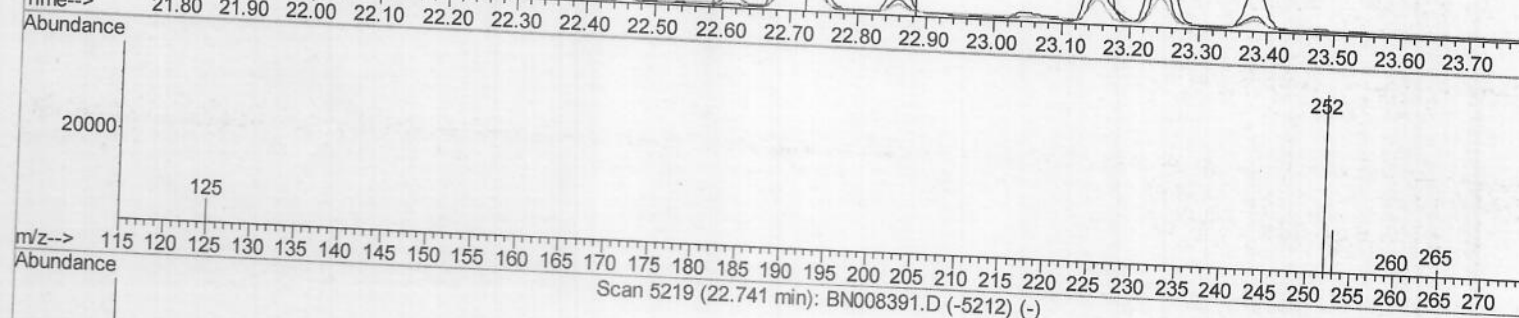
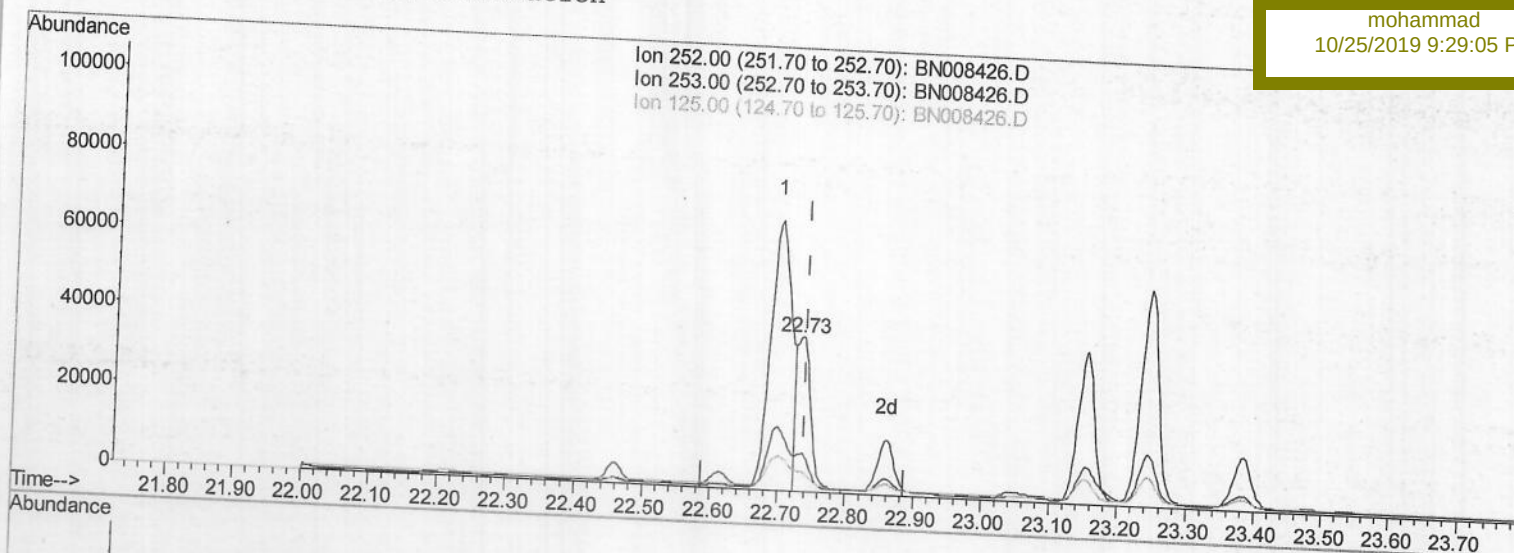
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(22) Benzo(k)fluoranthene

22.732min (-0.006) 1.54ng/ul m →

response 55692

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	24.39
125.00	15.40	13.18
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	7.57	152	1771	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	9065	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	6126	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	13405	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	11384	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	9778	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	461	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1877m	0.04	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.38	128	1801	0.070	ng/ul#	89
5) 2-Methylnaphthalene	12.02	142	1476	0.083	ng/ul	99
7) Acenaphthylene	13.93	152	13049	0.447	ng/ul	99
8) Acenaphthene	14.27	153	4231	0.186	ng/ul	98
9) Fluorene	15.26	166	9707	0.377	ng/ul	98
12) Phenanthrene	17.00	178	176890	4.507	ng/ul	99
13) Anthracene	17.09	178	31842	0.922	ng/ul	99
15) Fluoranthene	19.03	202	338888	6.538	ng/ul	100
17) Pyrene	19.39	202	264987	5.846	ng/ul	99
18) Benzo(a)anthracene	21.15	228	137154	3.314	ng/ul	98
19) Chrysene	21.20	228	132073	2.600	ng/ul	98
21) Benzo(b)fluoranthene	22.70	252	146394m	4.599	ng/ul	98
22) Benzo(k)fluoranthene	22.73	252	55692m	1.545	ng/ul	86
23) Benzo(a)pyrene	23.24	252	95602	2.854	ng/ul#	100
24) Indeno(1,2,3-cd)pyrene	25.48	276	56929	1.443	ng/ul	93
25) Dibenzo(a,h)anthracene	25.49	278	14396	0.462	ng/ul	98
26) Benzo(g,h,i)perylene	26.14	276	51629	1.427	ng/ul	98

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

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