

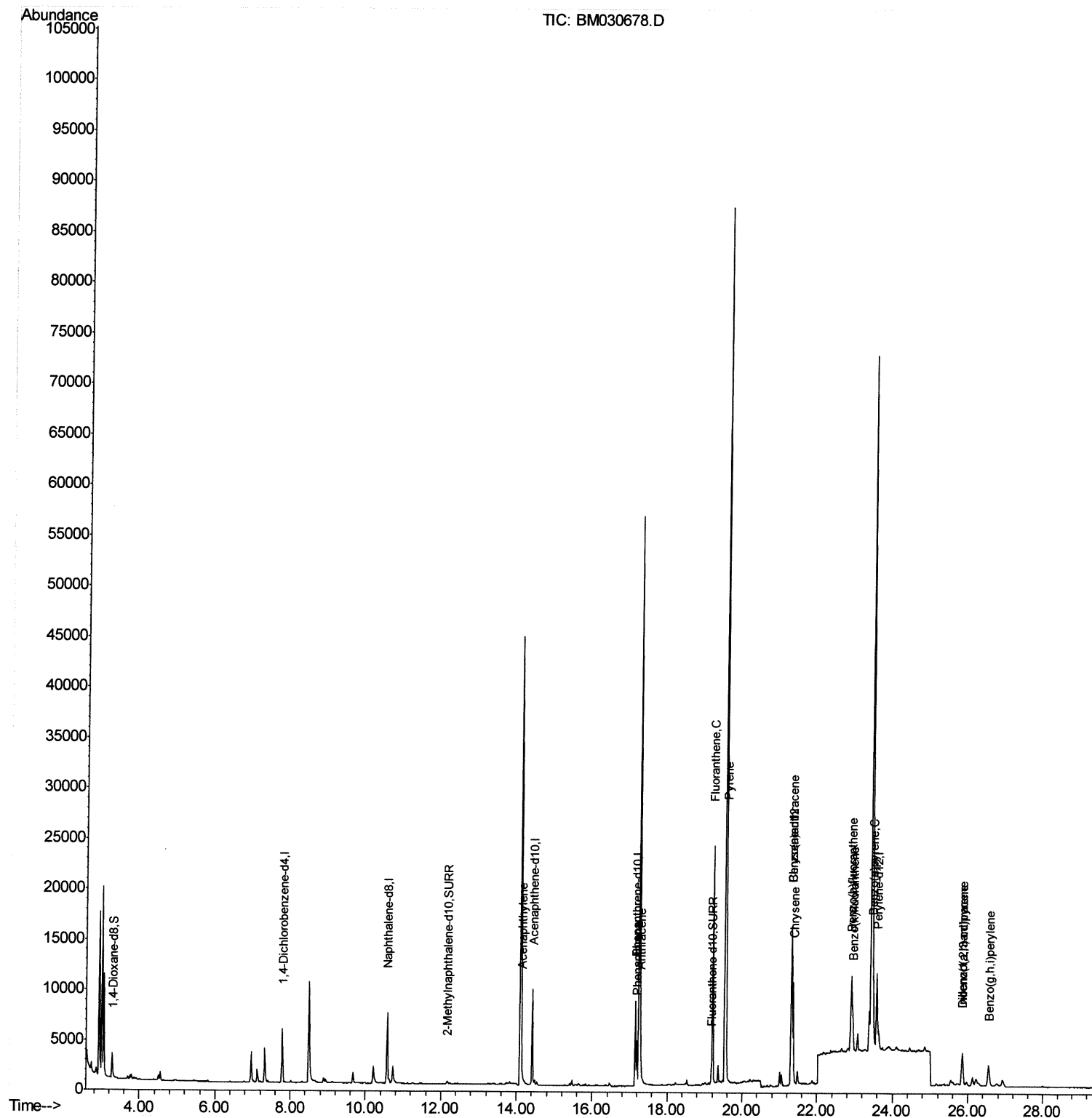
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062821\
Data File : BM030678.D
Acq On : 29 Jun 2021 06:58
Operator : CG/JU
Sample : M2704-11DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGDD4DL

Manual Integrations
APPROVED

mohammad
6/29/2021 2:32:23 PM

Quant Time: Jun 29 07:53:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM062821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 29 07:43:40 2021
Response via : Initial Calibration



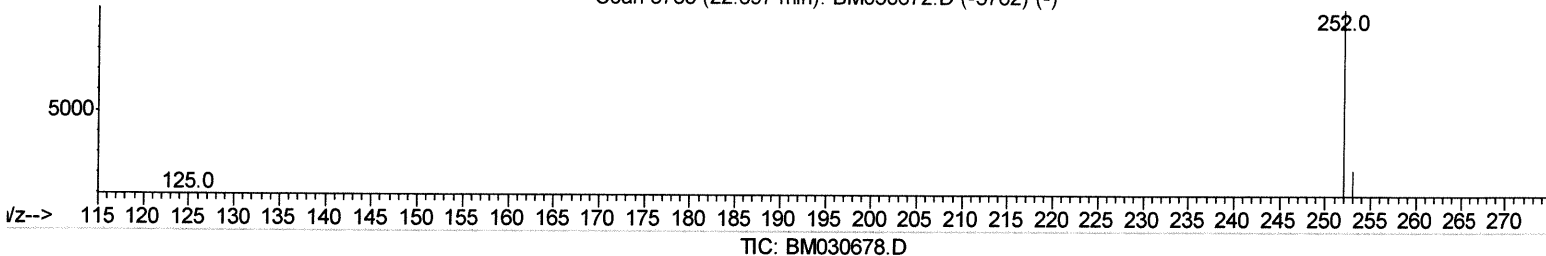
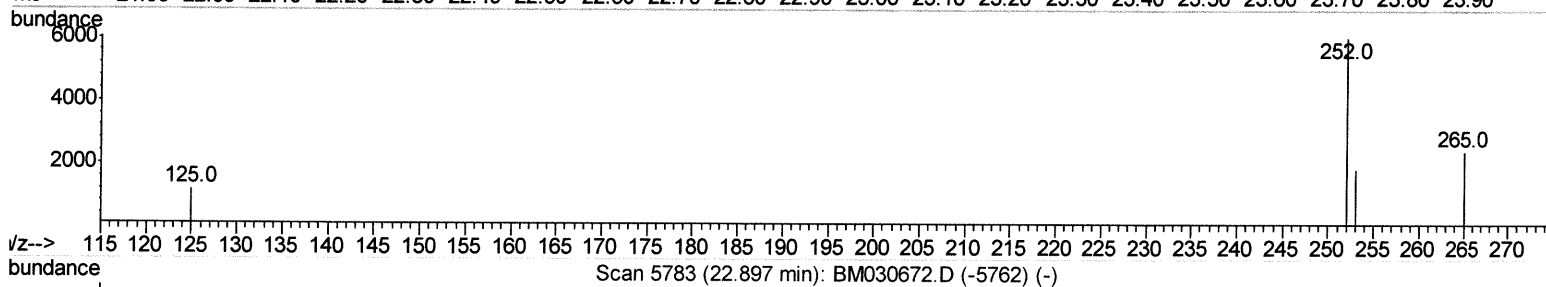
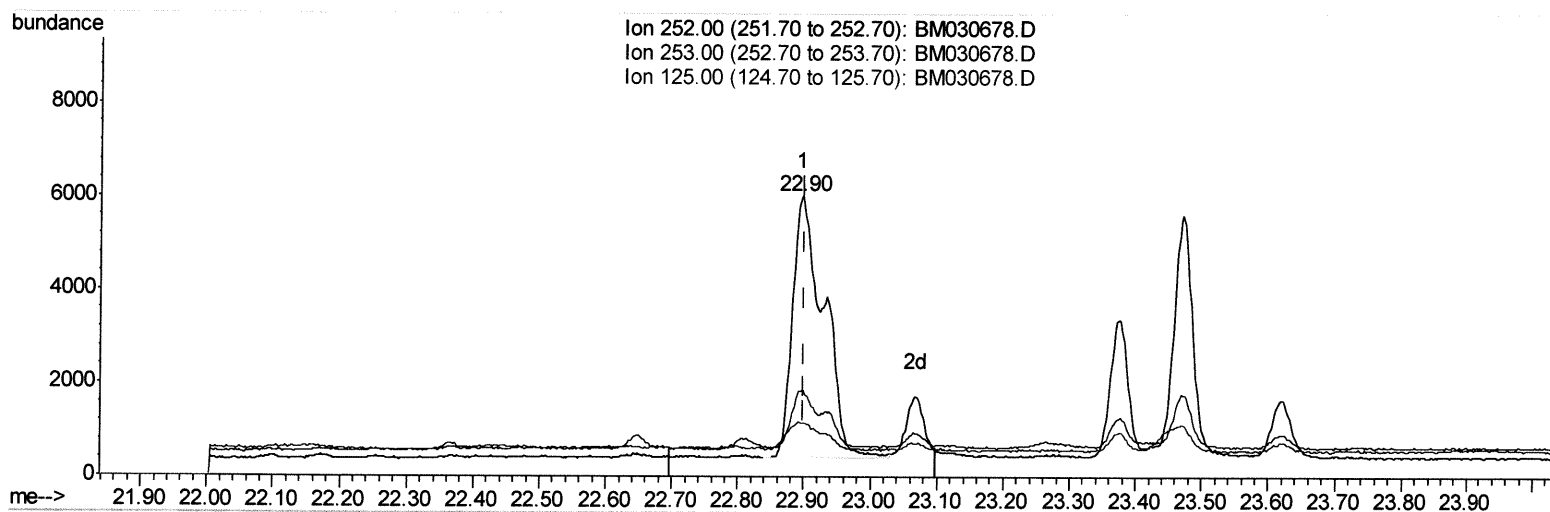
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(24) Benzo(b)fluoranthene

22.897min (-0.003) 0.40ng/ul

response 18453

Ion	Exp%	Act%
252.00	100	100
253.00	28.90	30.78
125.00	16.60	18.97
0.00	0.00	0.00

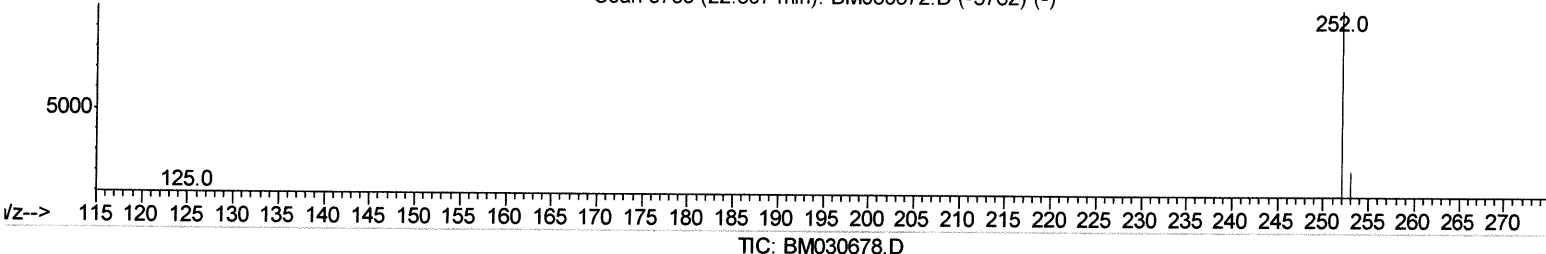
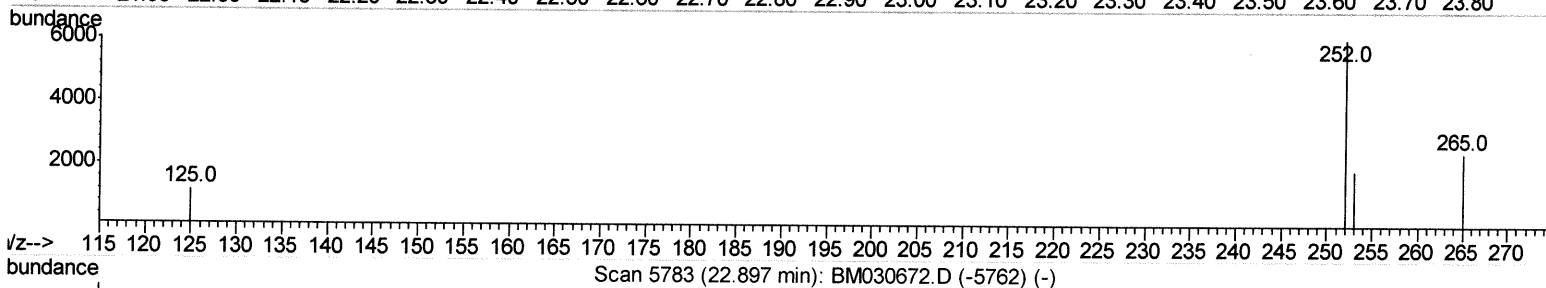
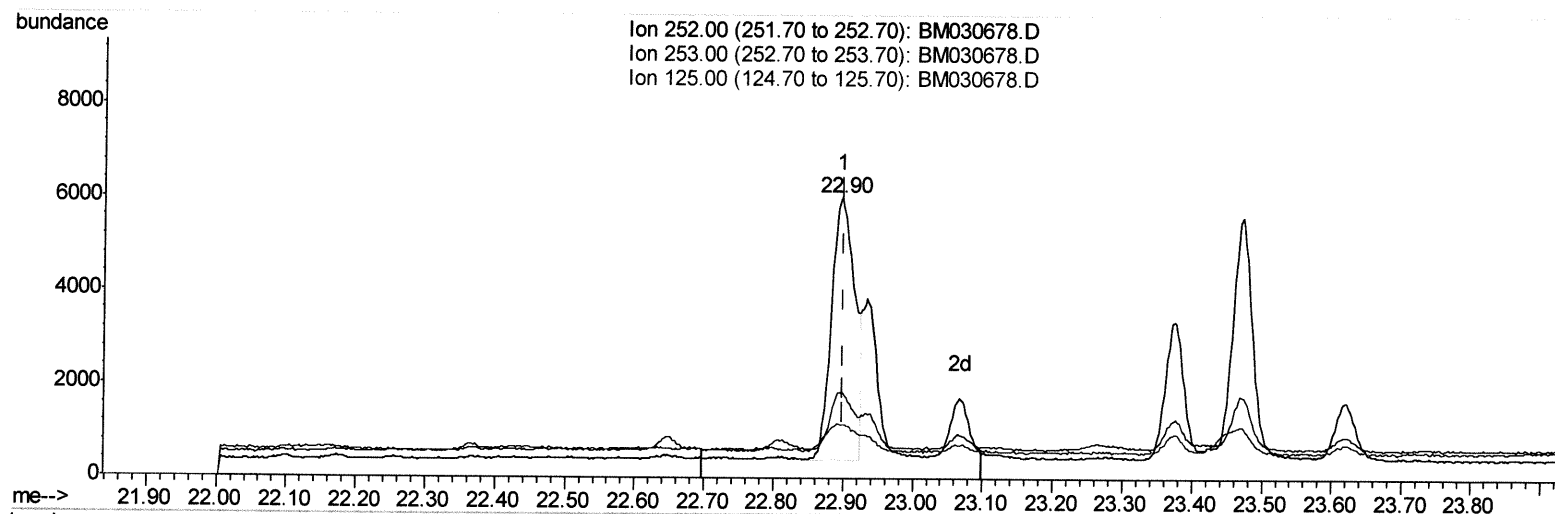
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(24) Benzo(b)fluoranthene

22.897min (-0.003) 0.28ng/ul m

JU 6/30/21

response 12926

Ion	Exp%	Act%
252.00	100	100
253.00	28.90	30.78
125.00	16.60	18.97
0.00	0.00	0.00

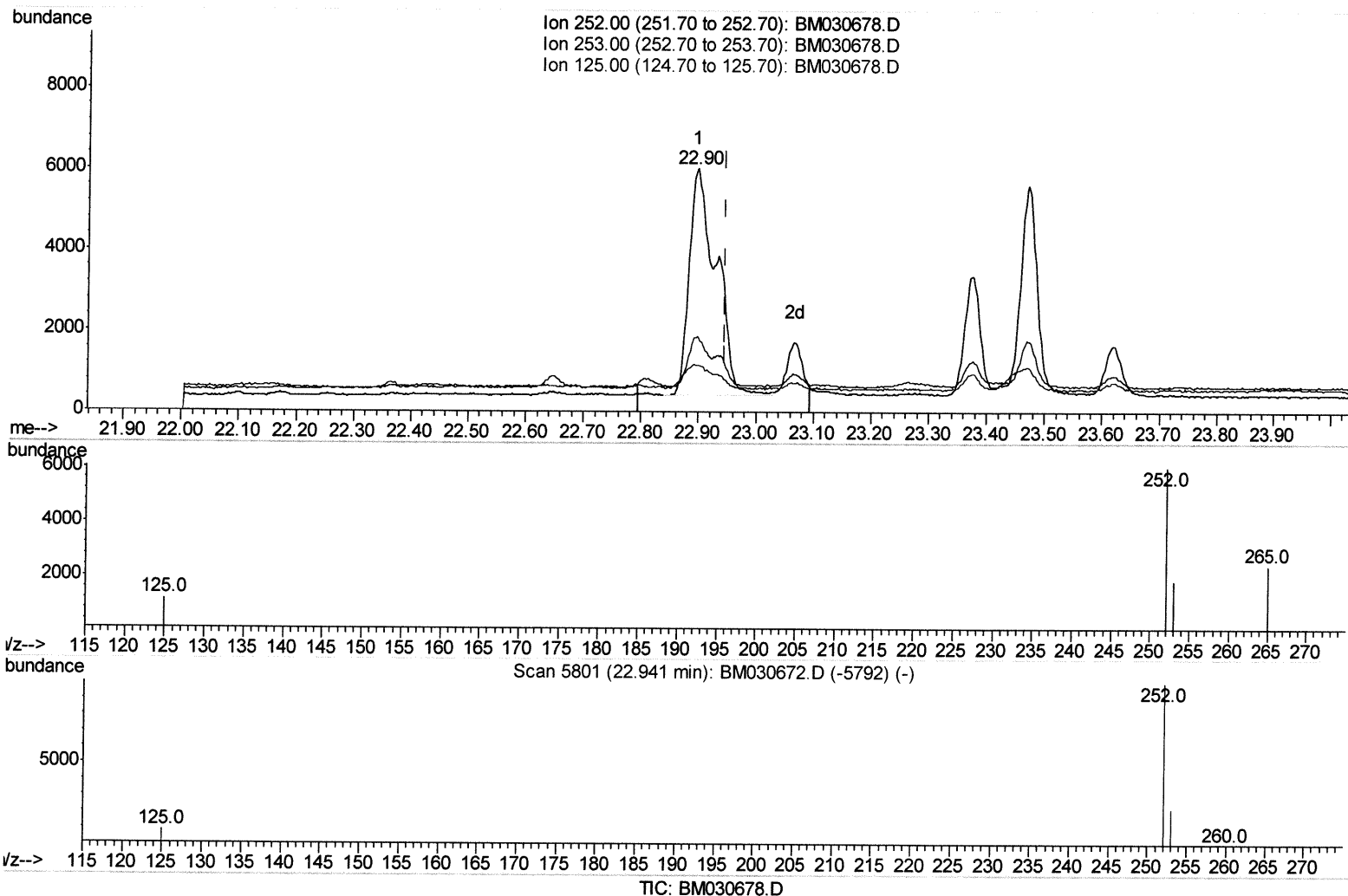
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Instrument :
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(25) Benzo(k)fluoranthene

22.897min (-0.049) 0.39ng/ul

response 18453

Ion	Exp%	Act%
252.00	100	100
253.00	29.00	30.78
125.00	15.80	18.97#
0.00	0.00	0.00

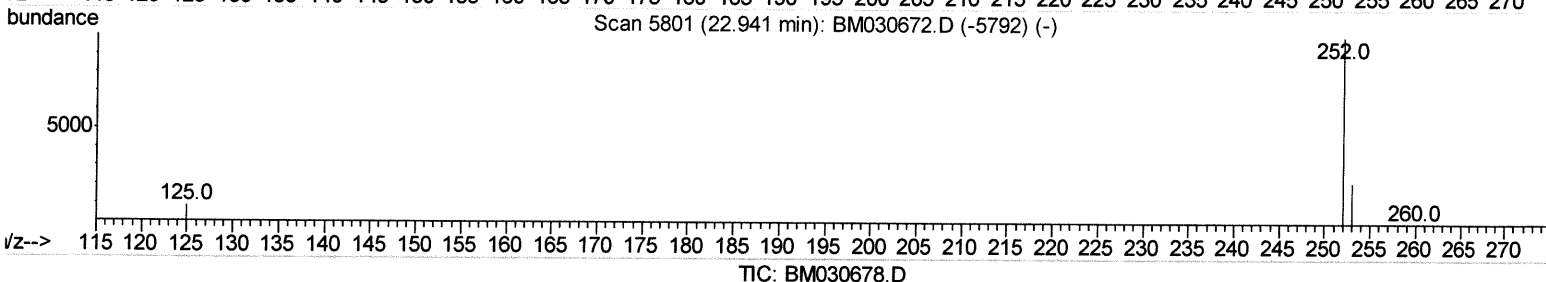
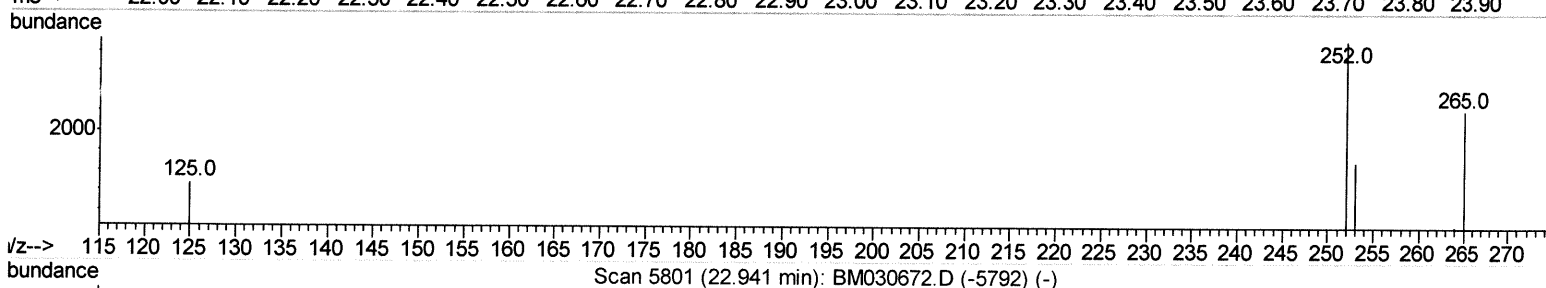
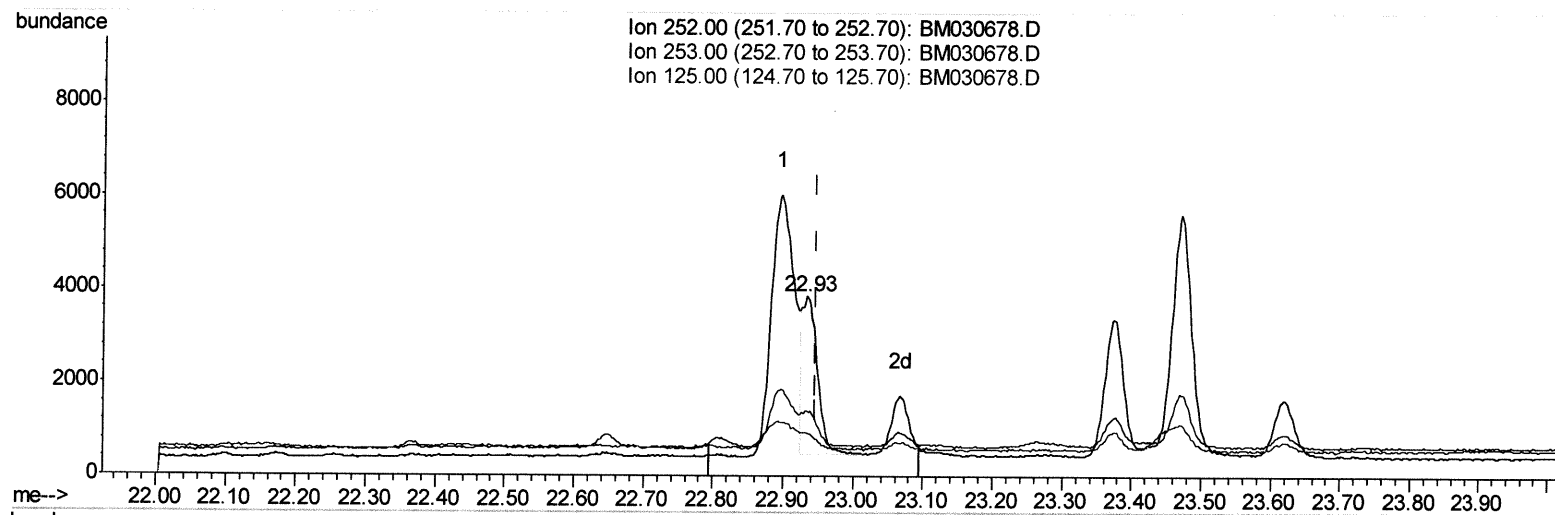
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(25) Benzo(k)fluoranthene

22.933min (-0.012) 0.11ng/ul m *JU 6/30/21*

response 5180

Ion	Exp%	Act%
252.00	100	100
253.00	29.00	36.43#
125.00	15.80	24.30#
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	2887	0.40	ng/ul	0.00
4) Naphthalene-d8	10.58	136	10392	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.42	164	5613	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.15	188	10636	0.40	ng/ul	0.00
17) Chrysene-d12	21.32	240	9683	0.40	ng/ul	0.00
23) Perylene-d12	23.57	264	10442	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.28	96	1808	0.57	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.17	152	501	0.03	ng/ul	0.00
18) Fluoranthene-d10	19.17	212	1145	0.04	ng/ul	0.00
Target Compounds						
					Ovalue	
10) Acenaphthylene	14.14	152	689	0.027	ng/ul#	77
15) Phenanthrene	17.19	178	4909	0.137	ng/ul	98
16) Anthracene	17.28	178	2532	0.081	ng/ul	93
19) Fluoranthene	19.20	202	22081	0.517	ng/ul	100
20) Pyrene	19.56	202	19366	0.451	ng/ul	98
21) Benzo(a)anthracene	21.30	228	12057	0.312	ng/ul	95
22) Chrysene	21.35	228	10202	0.236	ng/ul	97
24) Benzo(b)fluoranthene	22.90	252	12926m	0.279	ng/ul	} Jy 6/30/21
25) Benzo(k)fluoranthene	22.93	252	5180m	0.109	ng/ul	
26) Benzo(a)pyrene	23.47	252	9957	0.253	ng/ul	
27) Indeno(1,2,3-cd)pyrene	25.85	276	6028	0.118	ng/ul#	92
28) Dibenzo(a,h)anthracene	25.85	278	1492	0.036	ng/ul#	38
29) Benzo(a,h,i)perylene	26.55	276	4667	0.105	ng/ul#	92

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