

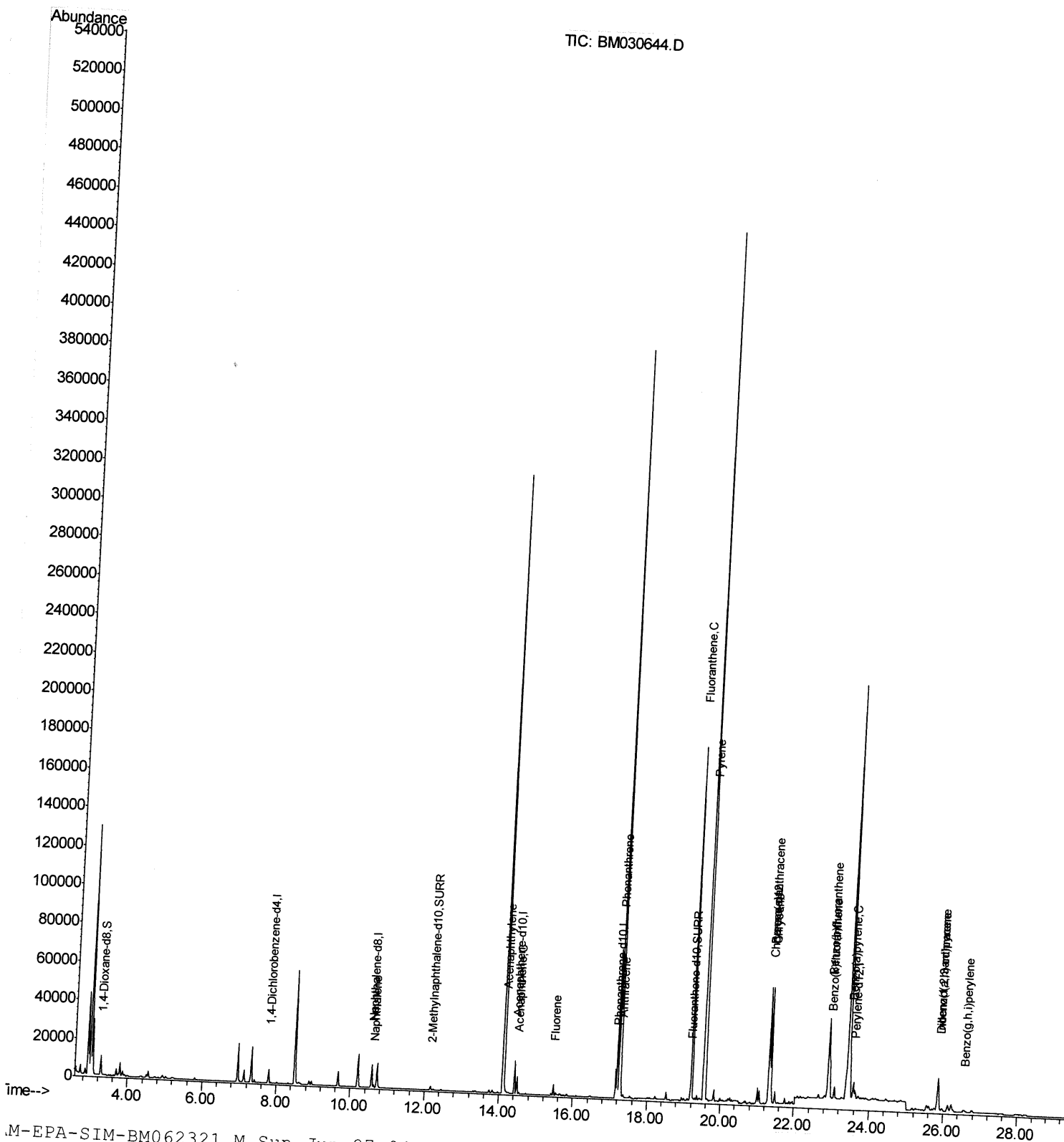
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062521\
Data File : BM030644.D
Acq On : 26 Jun 2021 18:57
Operator : CG/JU
Sample : M2682-17
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGDC5

Manual Integrations
APPROVED

mohammad
6/28/2021 1:23:18 PM

Quant Time: Jun 27 04:10:23 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM062321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 25 10:09:59 2021
Response via : Initial Calibration



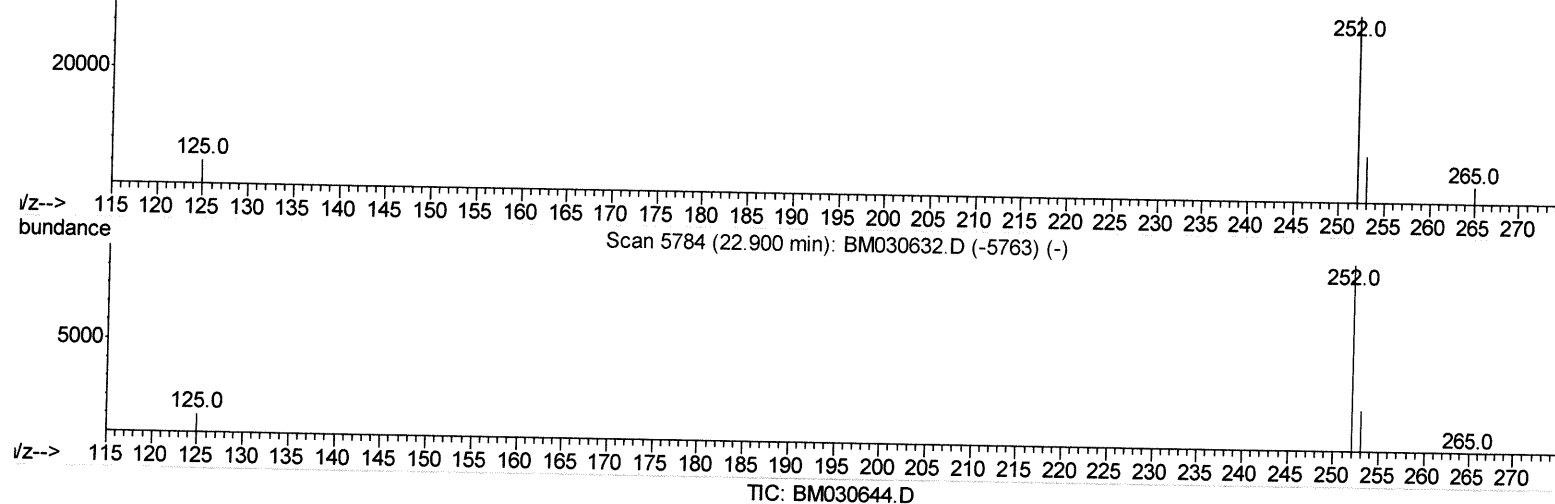
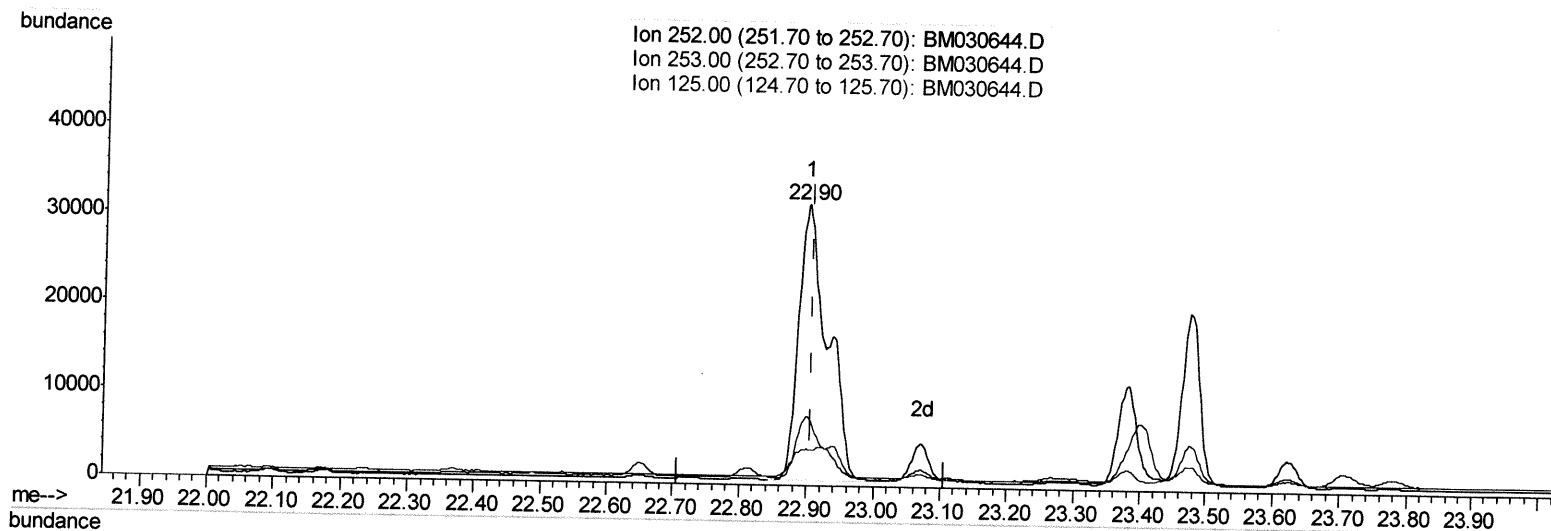
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Quant Time: Jun 27 02:42:16 2021
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(24) Benzo(b)fluoranthene

22.900min (-0.007) 1.85ng/ul

response 93603

Ion	Exp%	Act%
252.00	100	100
253.00	25.10	24.78
125.00	13.90	12.99
0.00	0.00	0.00

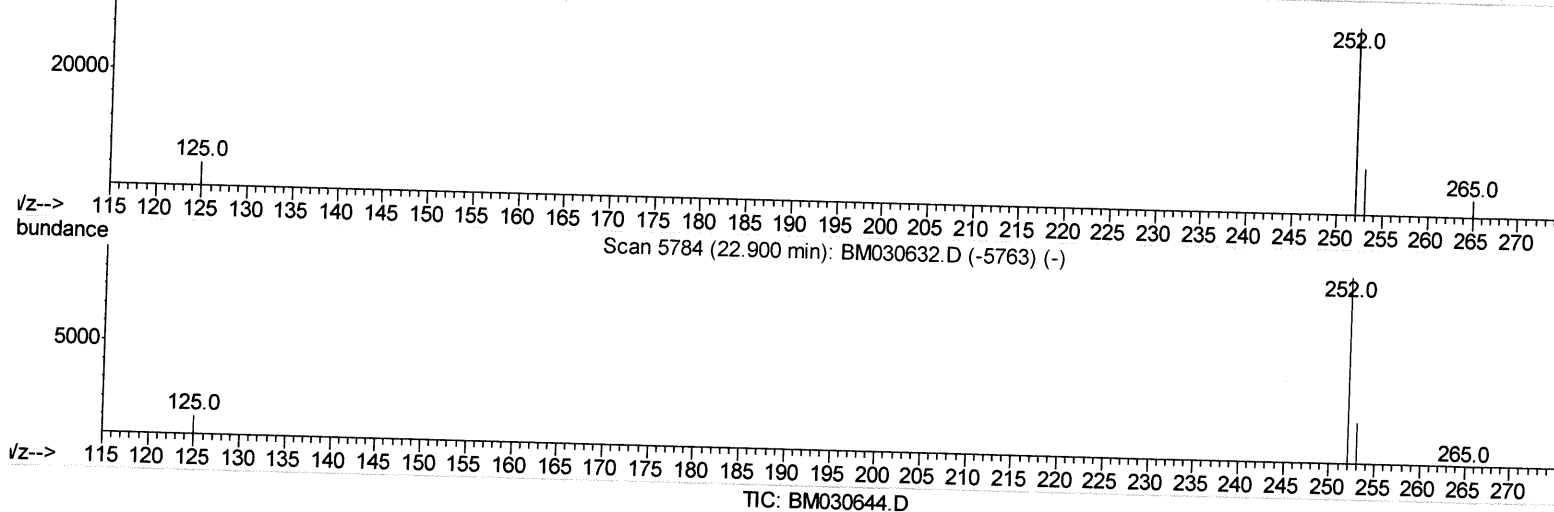
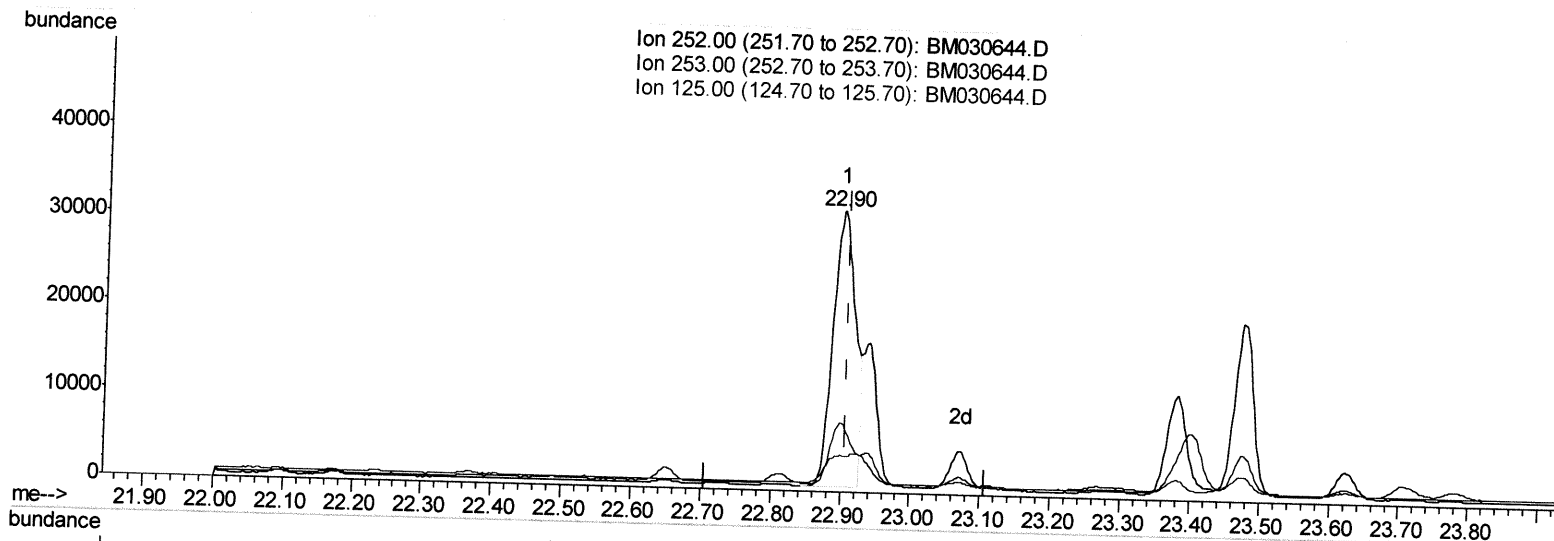
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(24) Benzo(b)fluoranthene

22.900min (-0.007) 1.35ng/ul m

response 68363

Ion	Exp%	Act%
252.00	100	100
253.00	25.10	24.78
125.00	13.90	12.99
0.00	0.00	0.00

JU 6/30/21

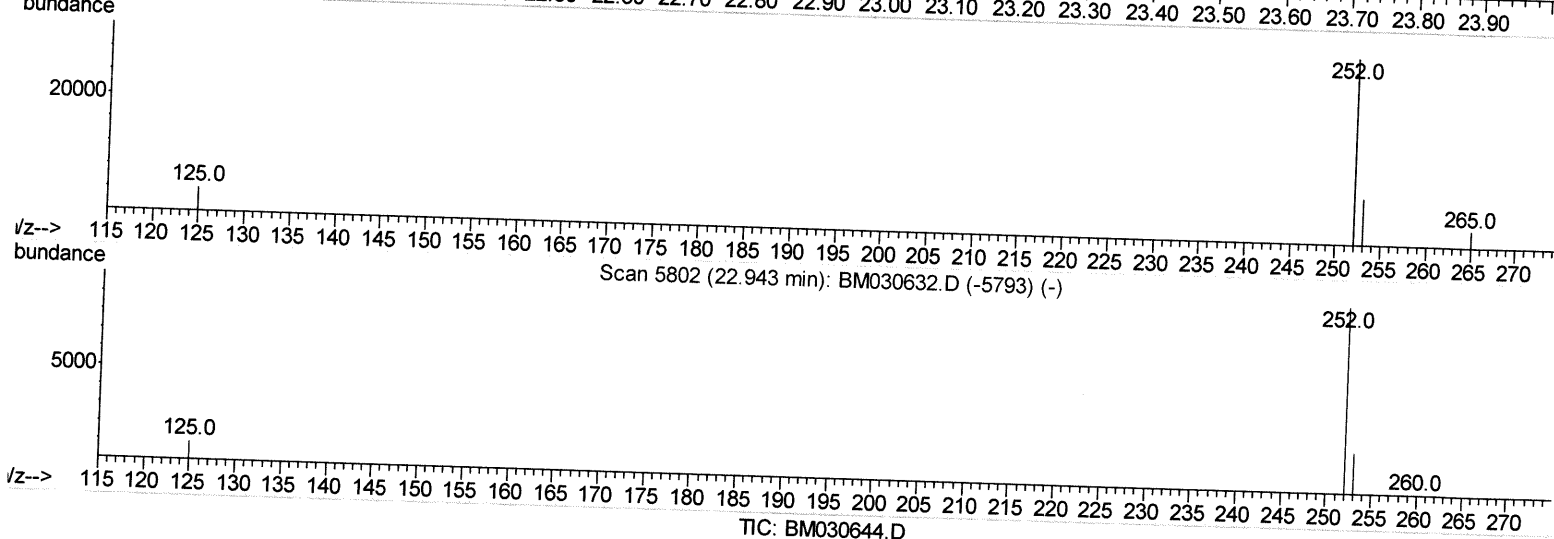
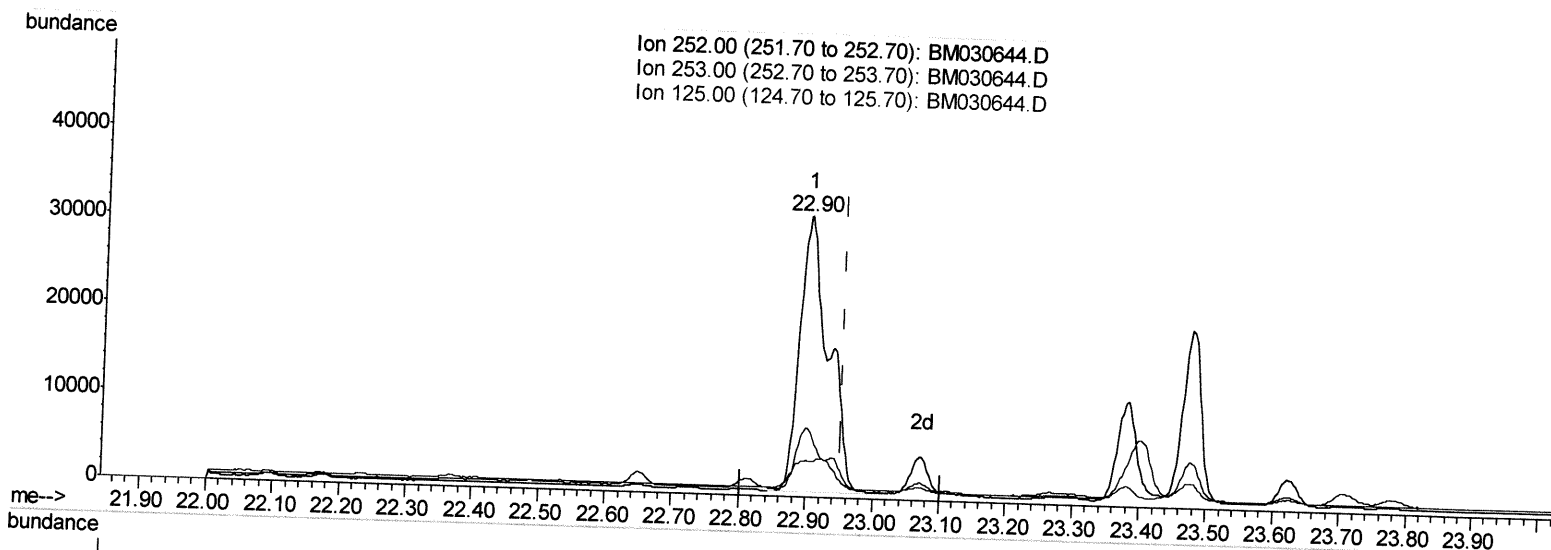
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Instrument :
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Quant Time: Jun 27 04:09:11 2021
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(25) Benzo(k)fluoranthene

22.900min (-0.053) 1.76ng/ul

response 93603

Ion	Exp%	Act%
252.00	100	100
253.00	25.80	24.78
125.00	13.70	12.99
0.00	0.00	0.00

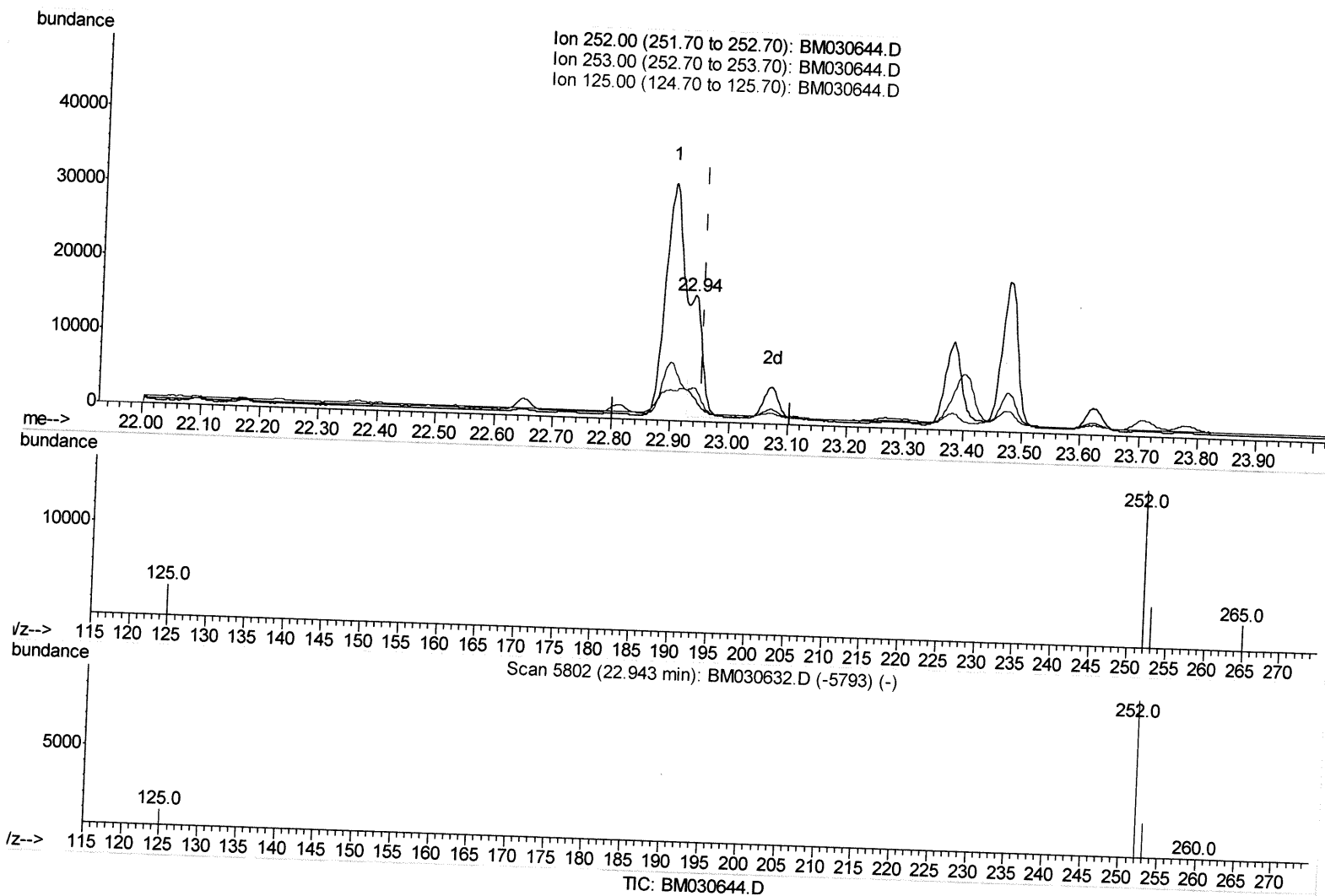
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(25) Benzo(k)fluoranthene

22.939min (-0.014) 0.47ng/ul m

response 24949

Ion	Exp%	Act%
252.00	100	100
253.00	25.80	26.87
125.00	13.70	20.01#
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	3751	0.40	ng/ul	0.00
4) Naphthalene-d8	10.58	136	15692	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.42	164	9559	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.16	188	17890	0.40	ng/ul	0.00
17) Chrysene-d12	21.32	240	11637	0.40	ng/ul	0.00
23) Perylene-d12	23.58	264	11628	0.40	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.28	96	6947	1.73	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.17	152	2841	0.11	ng/ul	0.00
18) Fluoranthene-d10	19.18	212	6402	0.18	ng/ul	0.00
Target Compounds						
5) Naphthalene	10.63	128	1846	0.040	ng/ul#	88
10) Acenaphthylene	14.14	152	2221	0.050	ng/ul#	88
11) Acenaphthene	14.48	153	5407	0.148	ng/ul	97
12) Fluorene	15.46	166	3649	0.090	ng/ul	98
15) Phenanthrene	17.20	178	78729	1.236	ng/ul	99
16) Anthracene	17.29	178	9398	0.168	ng/ul	97
19) Fluoranthene	19.21	202	159648	2.984	ng/ul	98
20) Pyrene	19.57	202	119058	2.172	ng/ul	97
21) Benzo(a)anthracene	21.31	228	51388	1.112	ng/ul	97
22) Chrysene	21.36	228	58050	1.131	ng/ul	97
24) Benzo(b)fluoranthene	22.90	252	68363m	1.352	ng/ul	97
25) Benzo(k)fluoranthene	22.94	252	24949m	0.470	ng/ul	96
26) Benzo(a)pyrene	23.47	252	37149	0.828	ng/ul	92
27) Indeno(1,2,3-cd)pyrene	25.85	276	27221	0.477	ng/ul#	79
28) Dibenzo(a,h)anthracene	25.86	278	7745	0.170	ng/ul#	69
29) Benzo(a,h,i)perylene	26.56	276	2906	0.059	ng/ul#	

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