

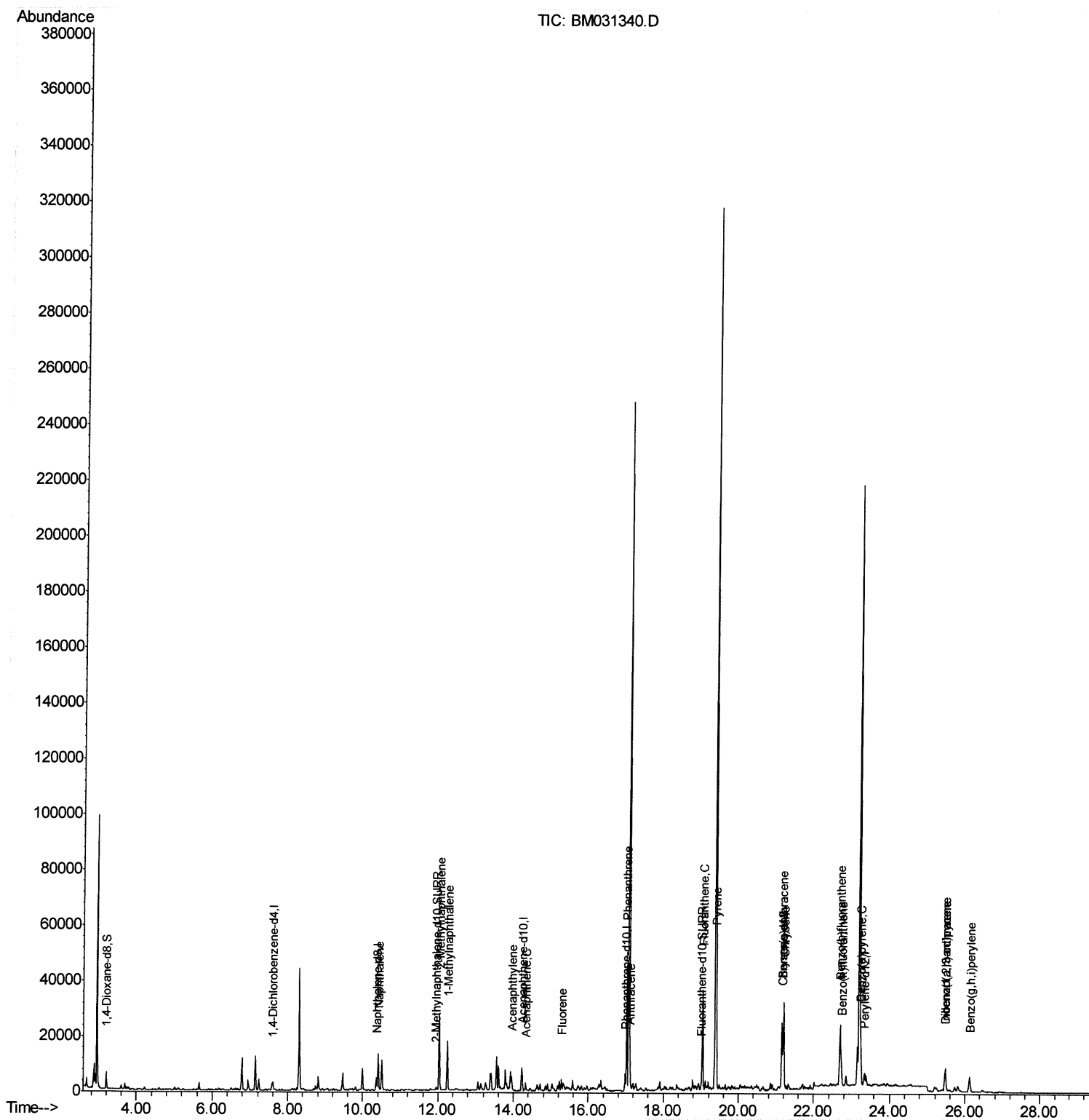
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM072921\
Data File : BM031340.D
Acq On : 01 Aug 2021 08:54
Operator : CG/JU
Sample : M3091-11
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0049

Manual Integrations
APPROVED

mohammad
8/2/2021 3:43:01 PM

Quant Time: Aug 02 04:53:58 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM072821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 30 14:16:52 2021
Response via : Initial Calibration



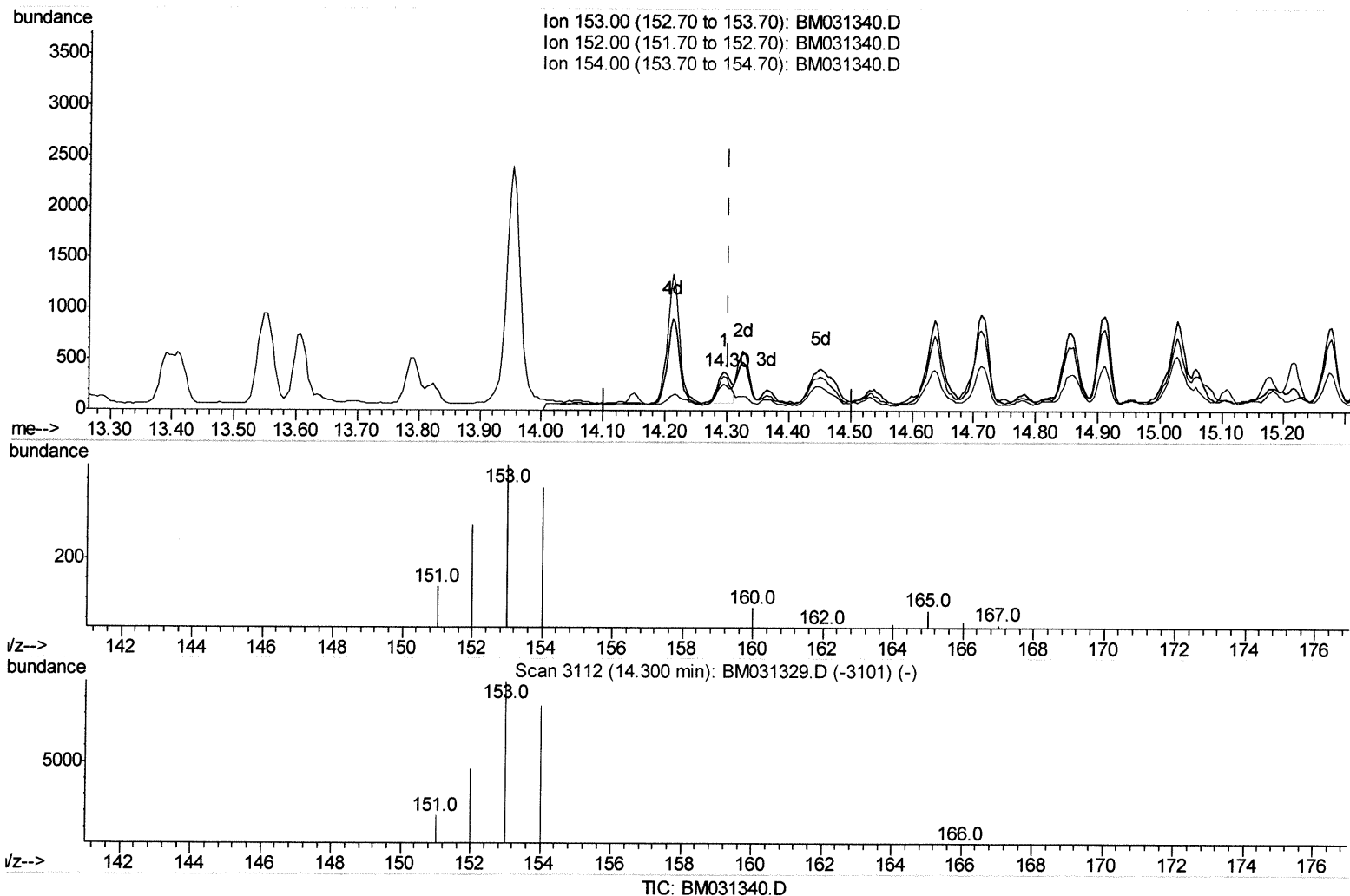
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(11) Acenaphthene (C)

14.295min (-0.008) 0.04ng/ul

response 473

Ion	Exp%	Act%
153.00	100	100
152.00	50.90	68.73#
154.00	87.70	88.37
0.00	0.00	0.00

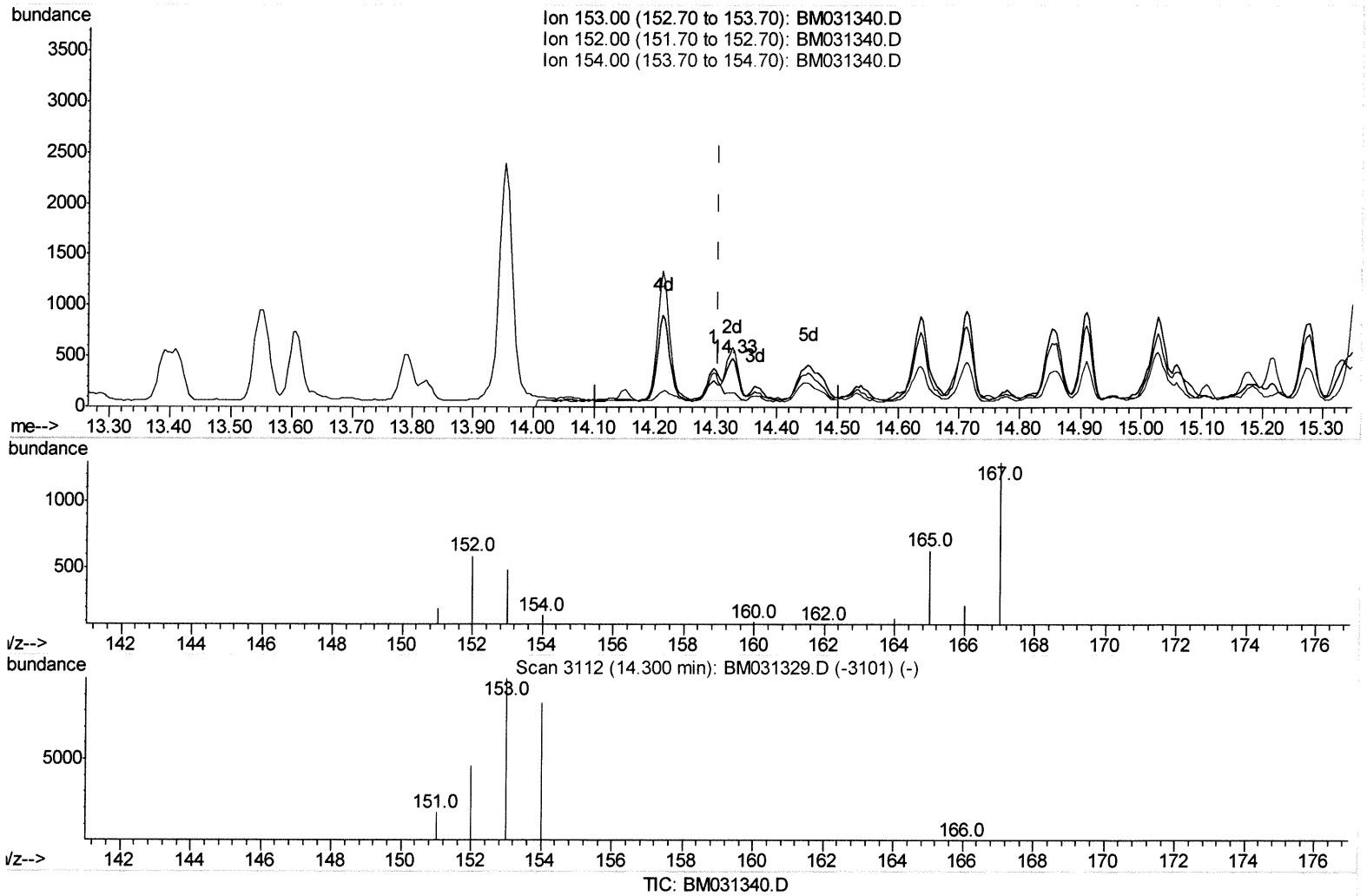
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(11) Acenaphthene (C)

14.326min (+0.023) 0.08ng/ul m

response 1027

Ion	Exp%	Act%
153.00	100	100
152.00	50.90	121.44#
154.00	87.70	28.87#
0.00	0.00	0.00

ju 8/3/21

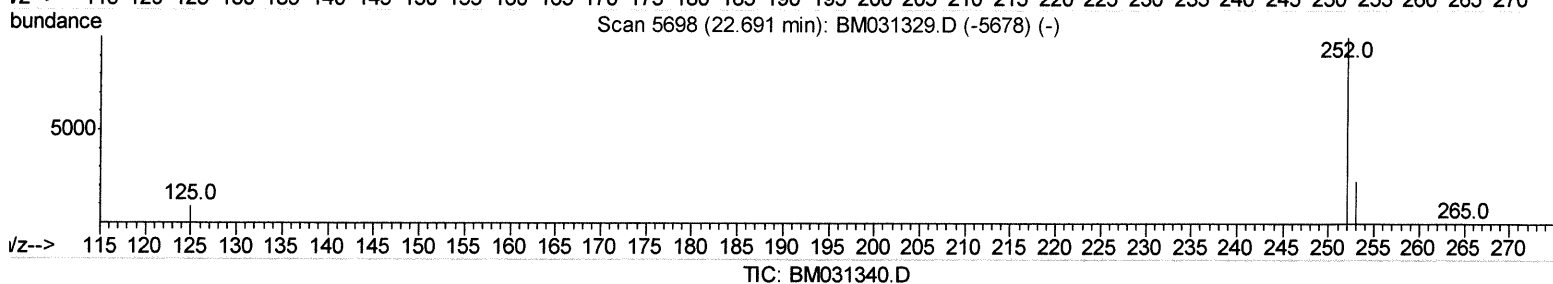
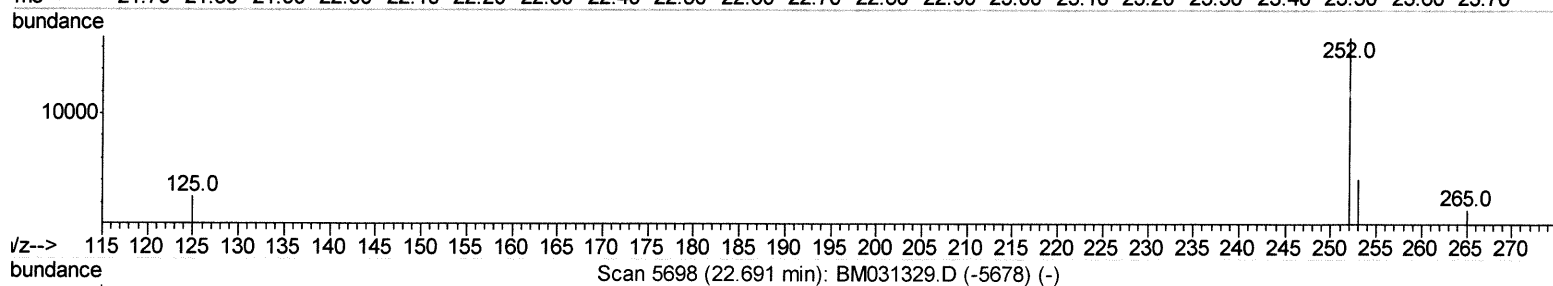
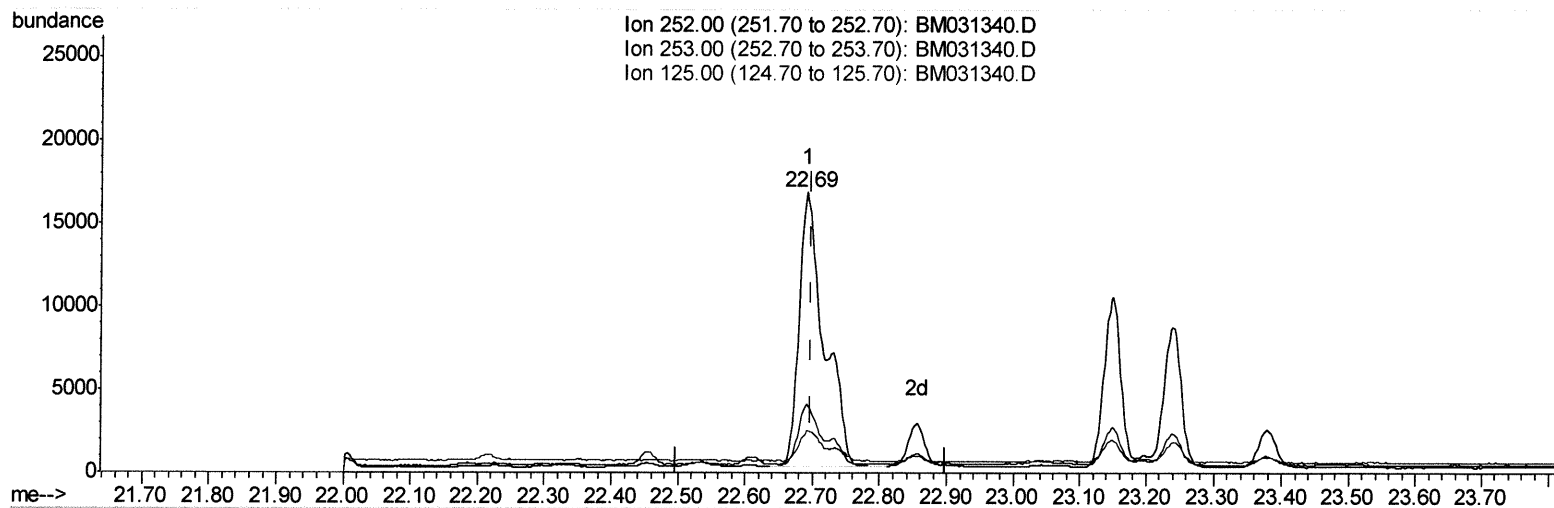
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Instrument :
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Quant Time: Aug 02 04:51:12 2021
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(24) Benzo(b)fluoranthene

22.691min (-0.007) 1.81ng/ul

response 42691

Ion	Exp%	Act%
252.00	100	100
253.00	26.50	24.50
125.00	11.40	14.90
0.00	0.00	0.00

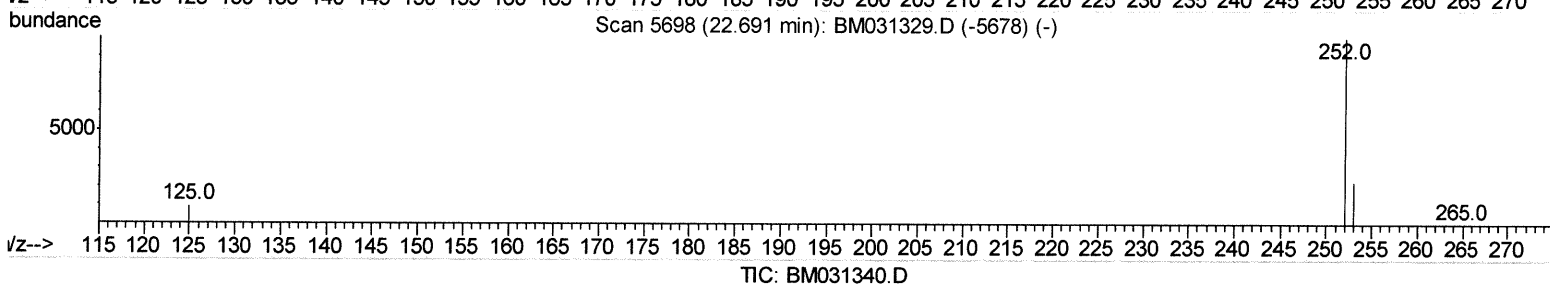
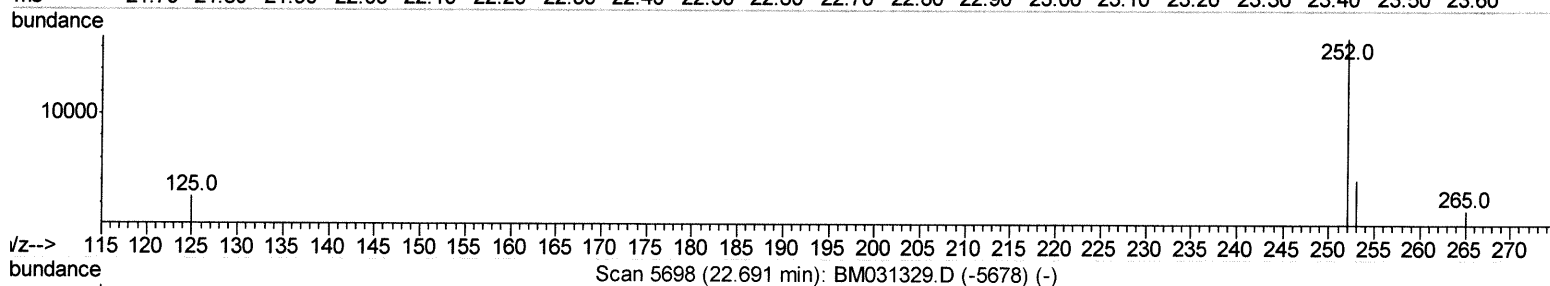
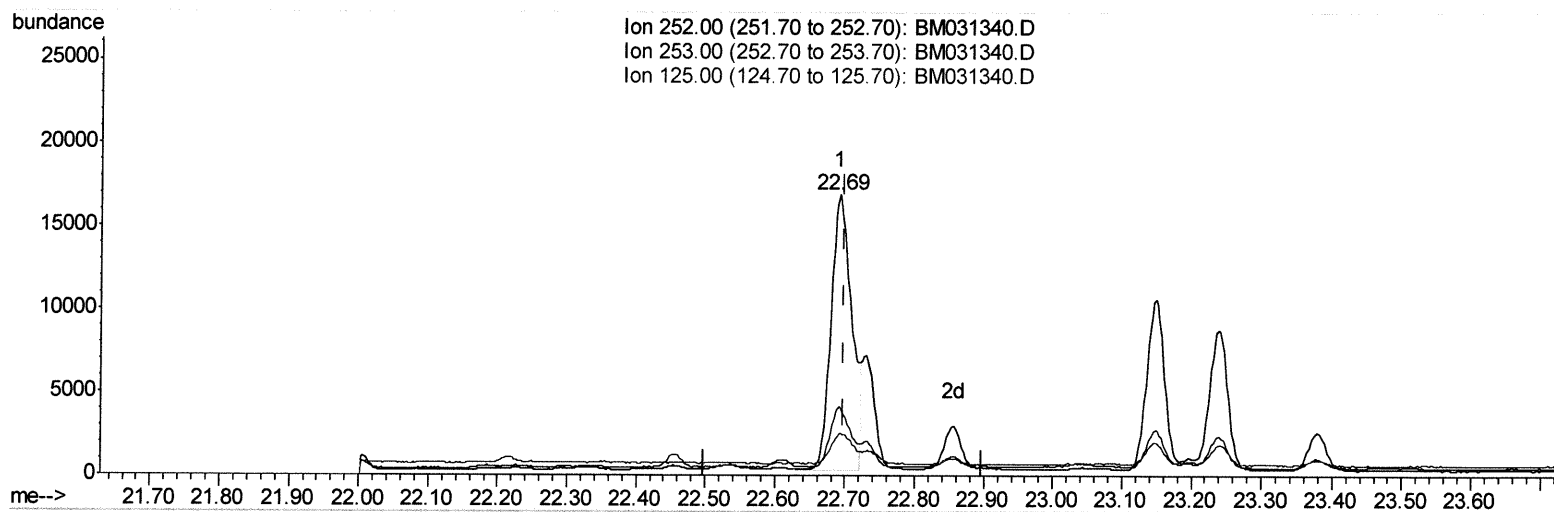
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 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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(24) Benzo(b)fluoranthene

22.691min (-0.007) 1.42ng/ul m

response 33466

Ion	Exp%	Act%
252.00	100	100
253.00	26.50	24.50
125.00	11.40	14.90
0.00	0.00	0.00

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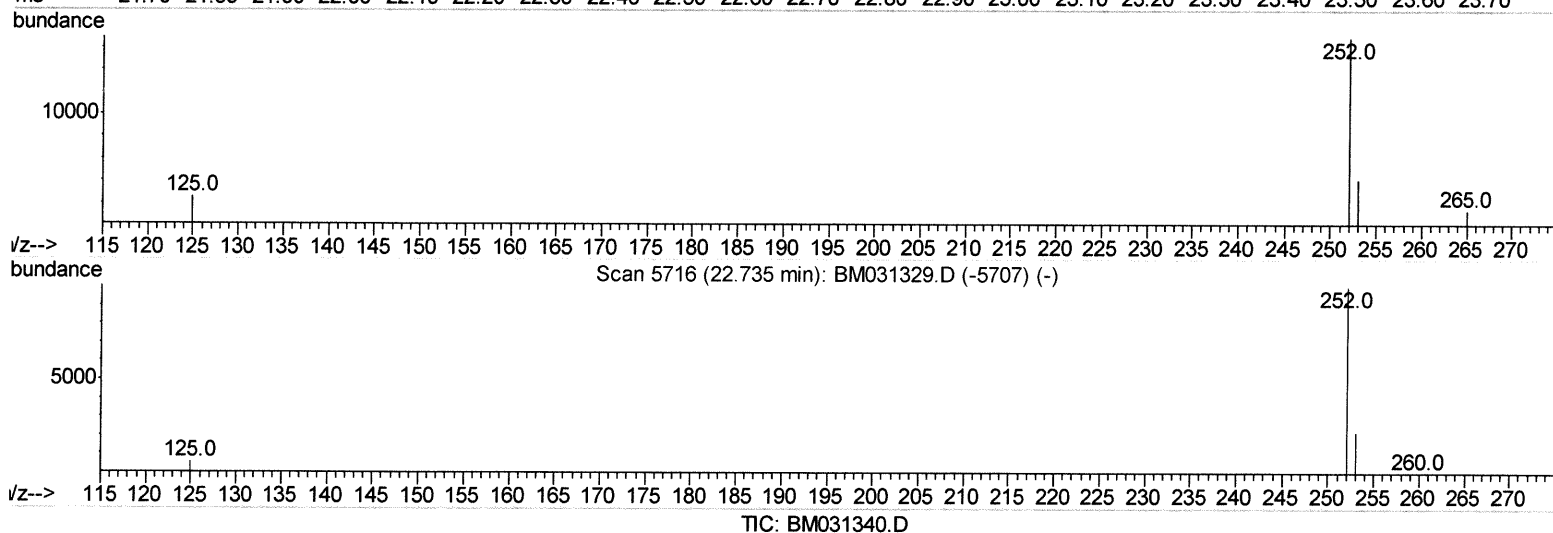
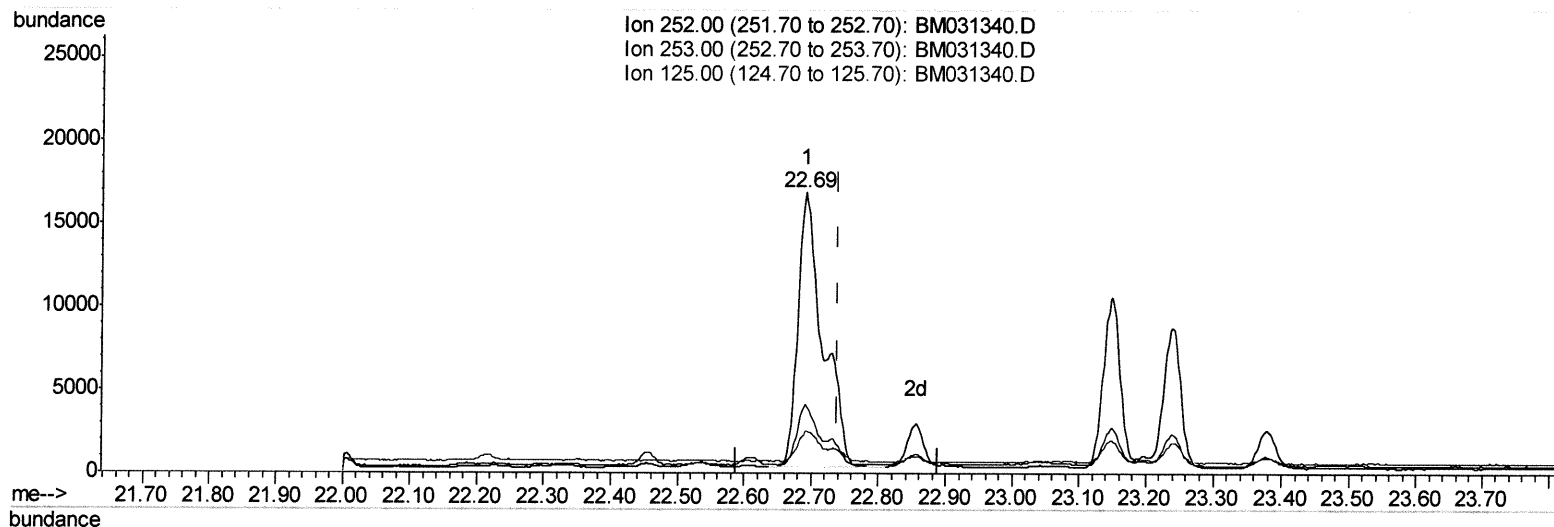
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(25) Benzo(k)fluoranthene

22.691min (-0.048) 1.75ng/ul

response 42691

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	24.50
125.00	11.00	14.90#
0.00	0.00	0.00

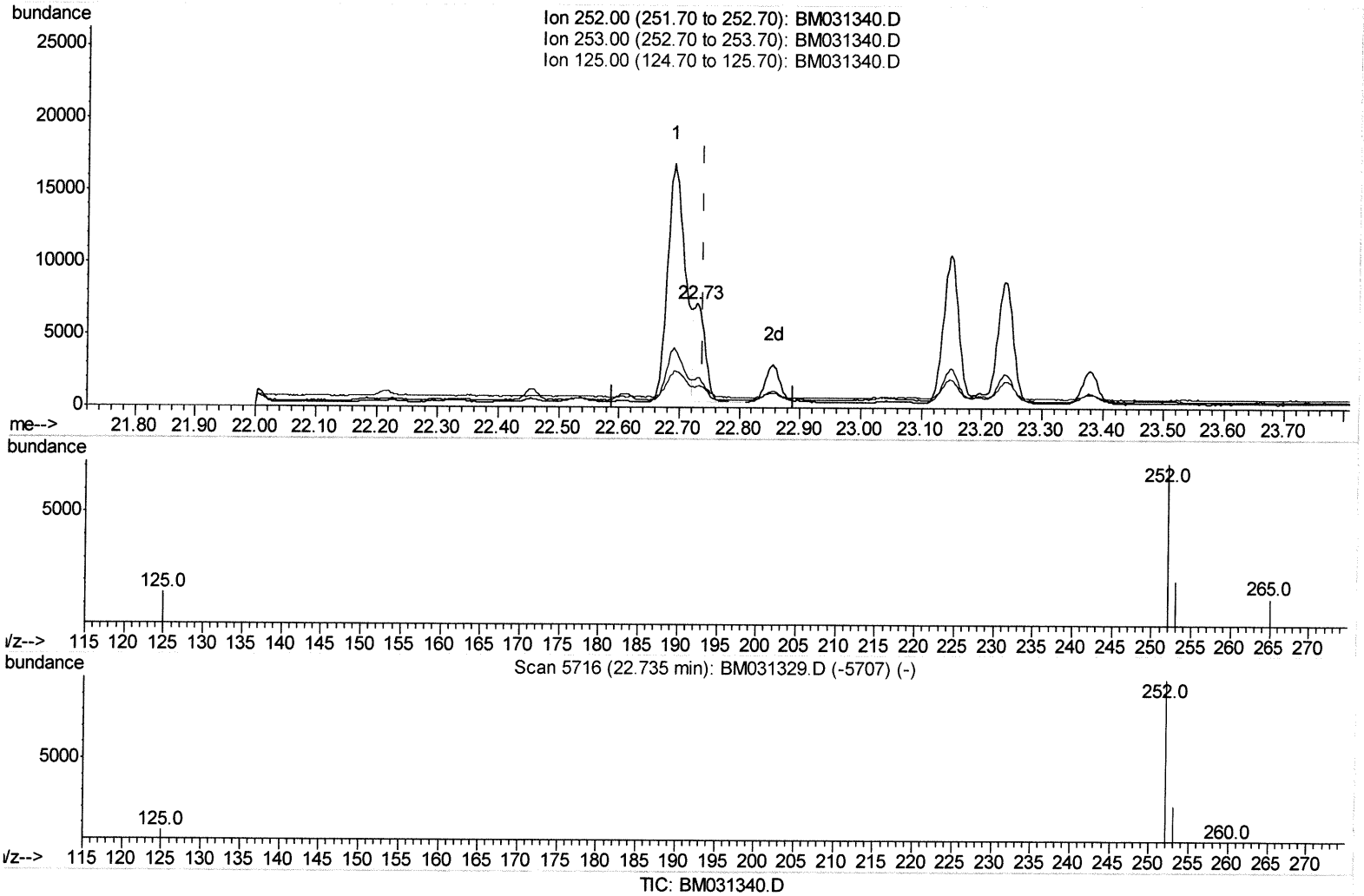
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Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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C0049

Manual Integrations
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Response via : Initial Calibration



(25) Benzo(k)fluoranthene

22.730min (-0.010) 0.39ng/ul m

response 9499

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	28.35
125.00	11.00	20.49#
0.00	0.00	0.00

rest 21

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1380	0.40	ng/ul	0.00
4) Naphthalene-d8	10.36	136	5375	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.23	164	3631	0.40	ng/ul	-0.01
13) Phenanthrene-d10	16.98	188	6858	0.40	ng/ul	0.00
17) Chrysene-d12	21.17	240	5605	0.40	ng/ul	0.00
23) Perylene-d12	23.33	264	5226	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.20	96	3842	2.51	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.96	152	1546	0.17	ng/ul	0.00
18) Fluoranthene-d10	19.01	212	3625	0.19	ng/ul	0.00
Target Compounds						
					Ovalue	
5) Naphthalene	10.41	128	17690	1.105	ng/ul	97
7) 2-Methylnaphthalene	12.03	142	18830	1.721	ng/ul	100
8) 1-Methylnaphthalene	12.25	142	11658	1.079	ng/ul	100
10) Acenaphthylene	13.95	152	3687	0.238	ng/ul	96
11) Acenaphthene	14.33	153	1027m	0.079	ng/ul	96 8/3/21
12) Fluorene	15.29	166	1243	0.083	ng/ul#	67
15) Phenanthrene	17.02	178	46120	2.118	ng/ul	98
16) Anthracene	17.11	178	4707	0.231	ng/ul	96
19) Fluoranthene	19.04	202	31082	1.299	ng/ul#	96
20) Pyrene	19.40	202	28912	1.191	ng/ul#	94
21) Benzo(a)anthracene	21.15	228	18528	0.819	ng/ul#	91
22) Chrysene	21.21	228	33380	1.425	ng/ul	96
24) Benzo(b)fluoranthene	22.69	252	33466m	1.421	ng/ul	96 8/3/21
25) Benzo(k)fluoranthene	22.73	252	9499m	0.390	ng/ul	96
26) Benzo(a)pyrene	23.24	252	15157	0.735	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.47	276	11862	0.445	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.48	278	4014	0.187	ng/ul#	73
29) Benzo(a,h,i)perylene	26.12	276	10711	0.471	ng/ul#	95

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