

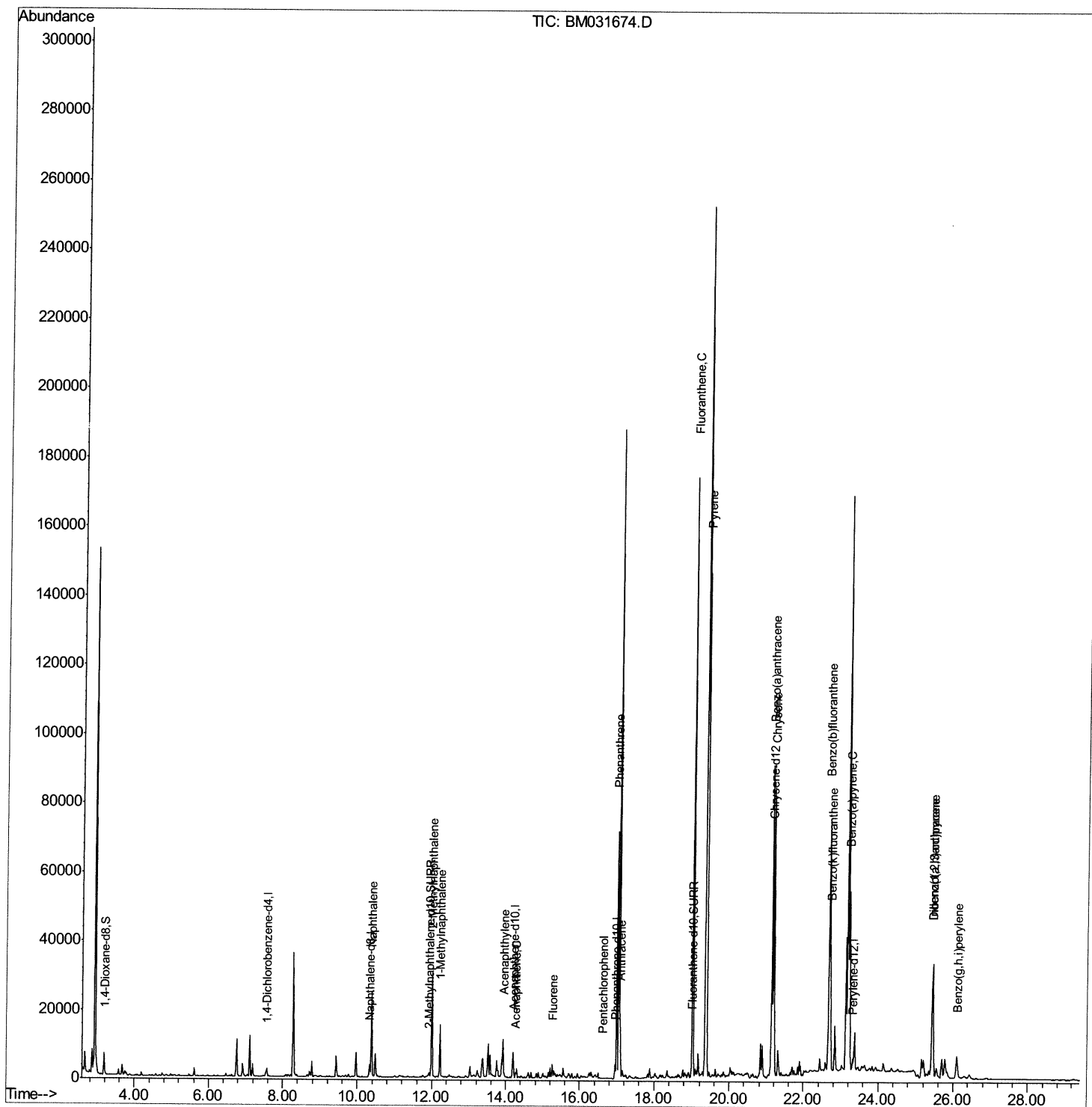
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\
Data File : BM031674.D
Acq On : 12 Aug 2021 13:08
Operator : CG/JU
Sample : M3182-13
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C00D1

Manual Integrations
APPROVED

mohammad
8/13/2021 5:13:48 PM

Quant Time: Aug 12 13:39:55 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 : Thu Aug 12 11:28:10 2021
Response via : Initial Calibration



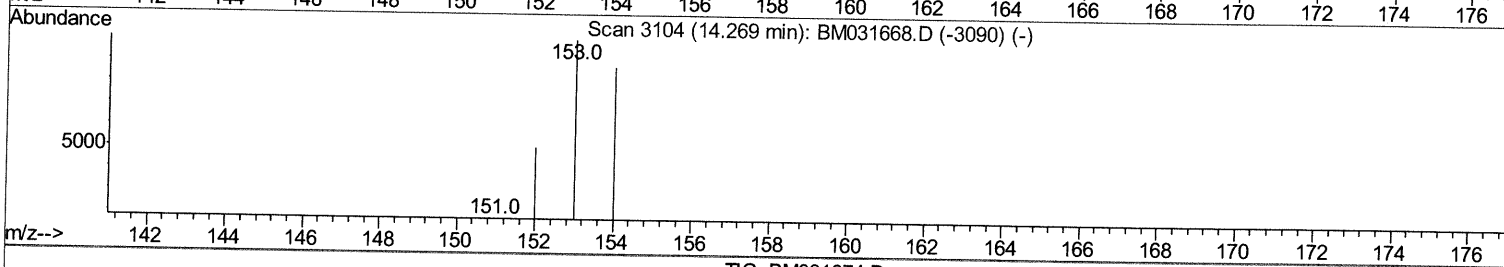
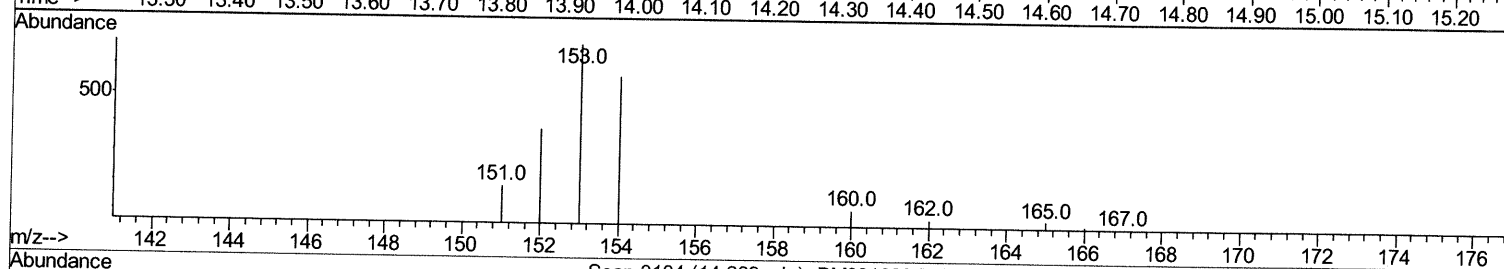
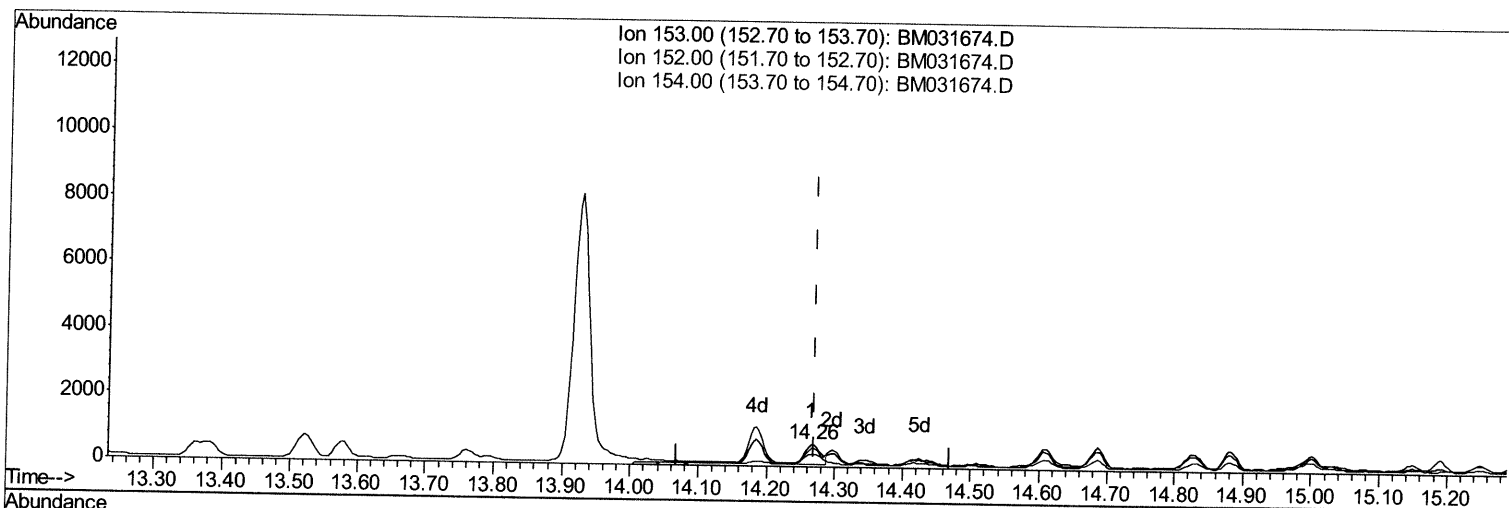
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TIC: BM031674.D

(11) Acenaphthene (C)

14.265min (-0.004) 0.10ng/ul

response 996

Ion	Exp%	Act%
153.00	100	100
152.00	50.90	57.12
154.00	87.70	83.70
0.00	0.00	0.00

Quantitation Report (Qedit)

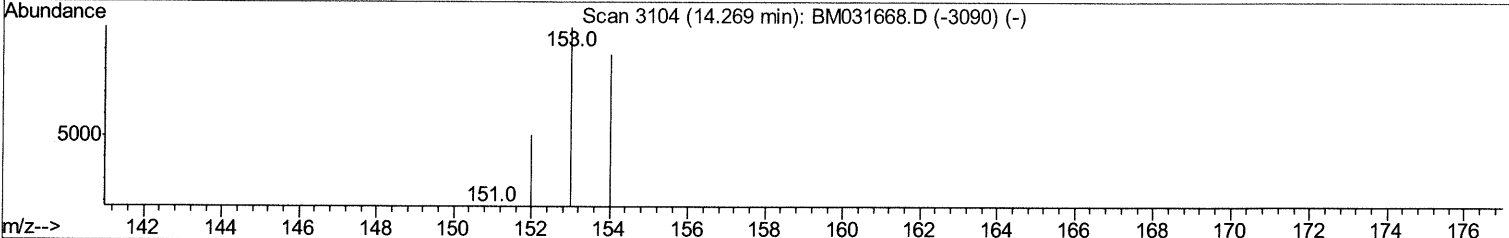
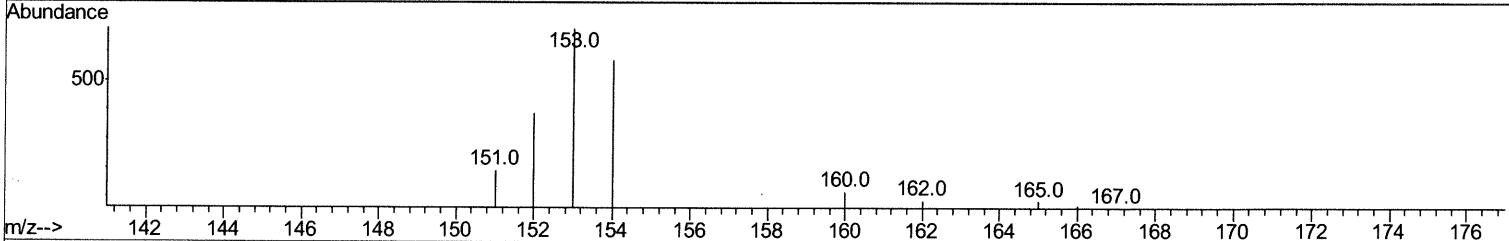
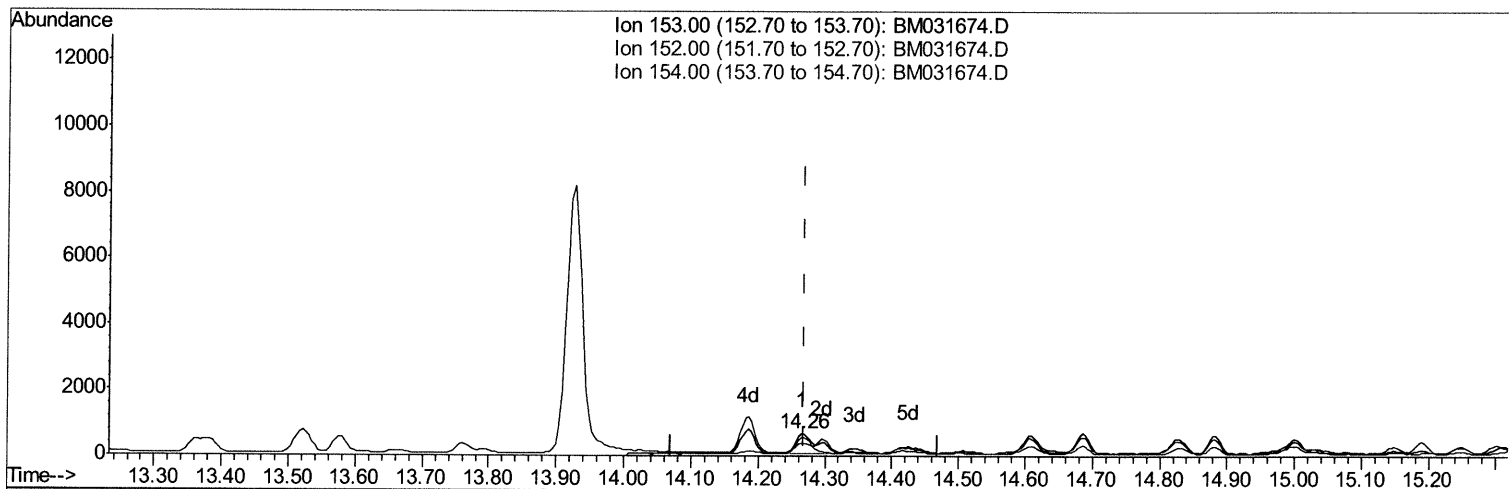
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TIC: BM031674.D

(11) Acenaphthene (C)

14.265min (-0.004) 0.14ng/ul m 08/16/21 JU

response 1421

Ion	Exp%	Act%
153.00	100	100
152.00	50.90	57.12
154.00	87.70	83.70
0.00	0.00	0.00

Quantitation Report (Qedit)

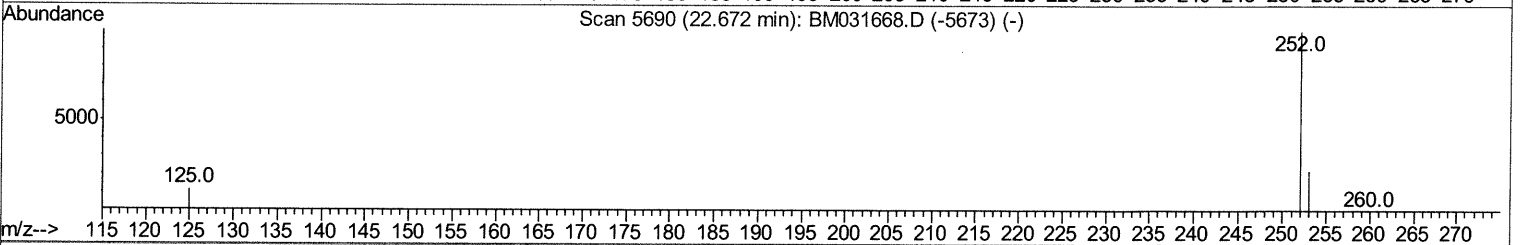
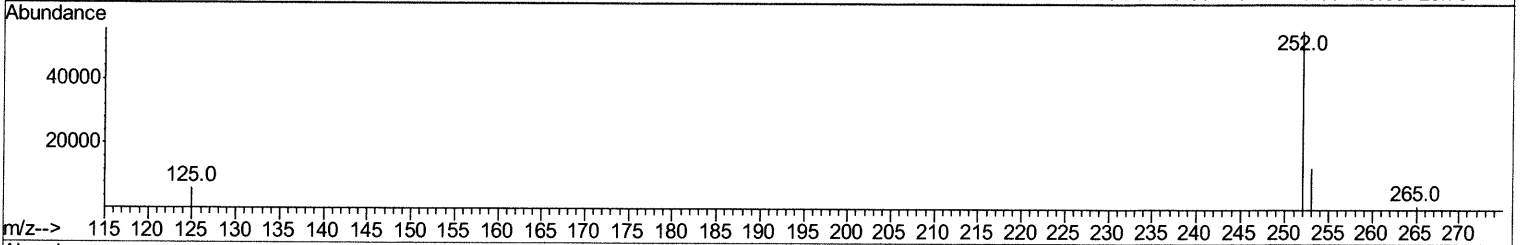
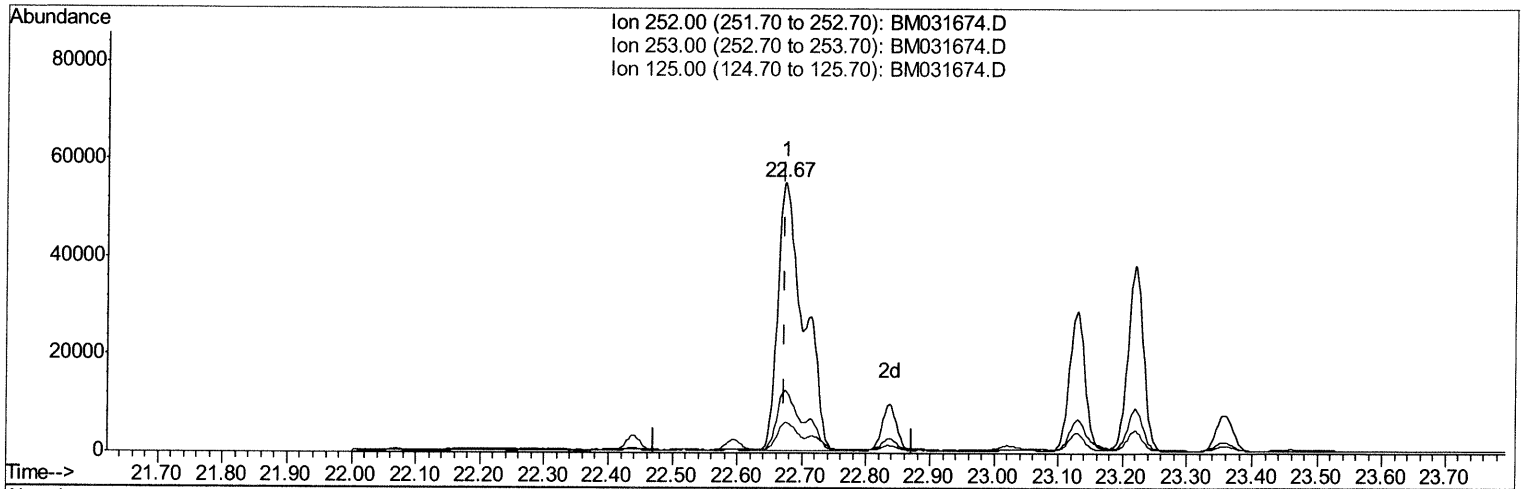
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TIC: BM031674.D

(24) Benzo(b)fluoranthene
 22.674min (+0.002) 8.99ng/ul
 response 153980

Ion	Exp%	Act%
252.00	100	100
253.00	26.50	23.23
125.00	11.40	11.20
0.00	0.00	0.00

Quantitation Report (Qedit)

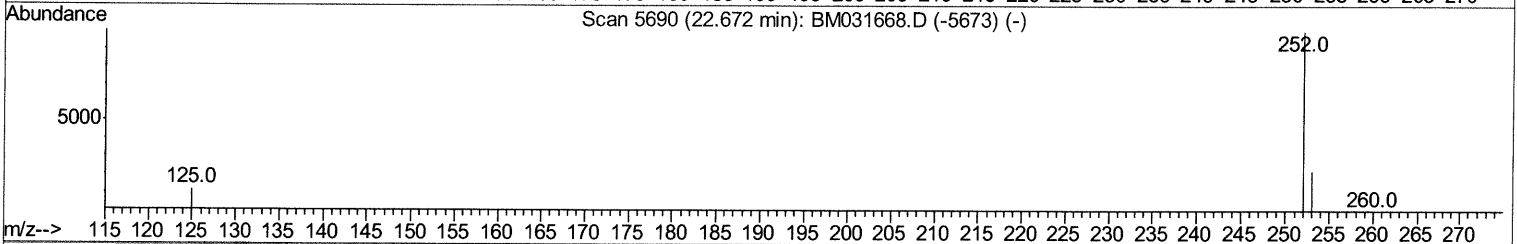
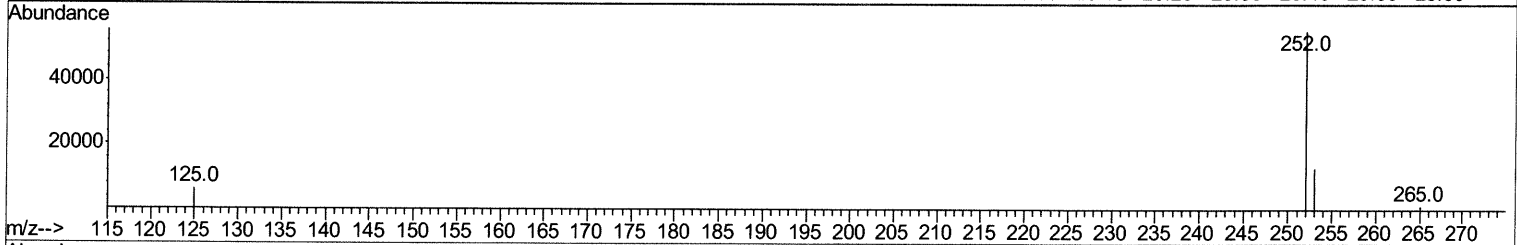
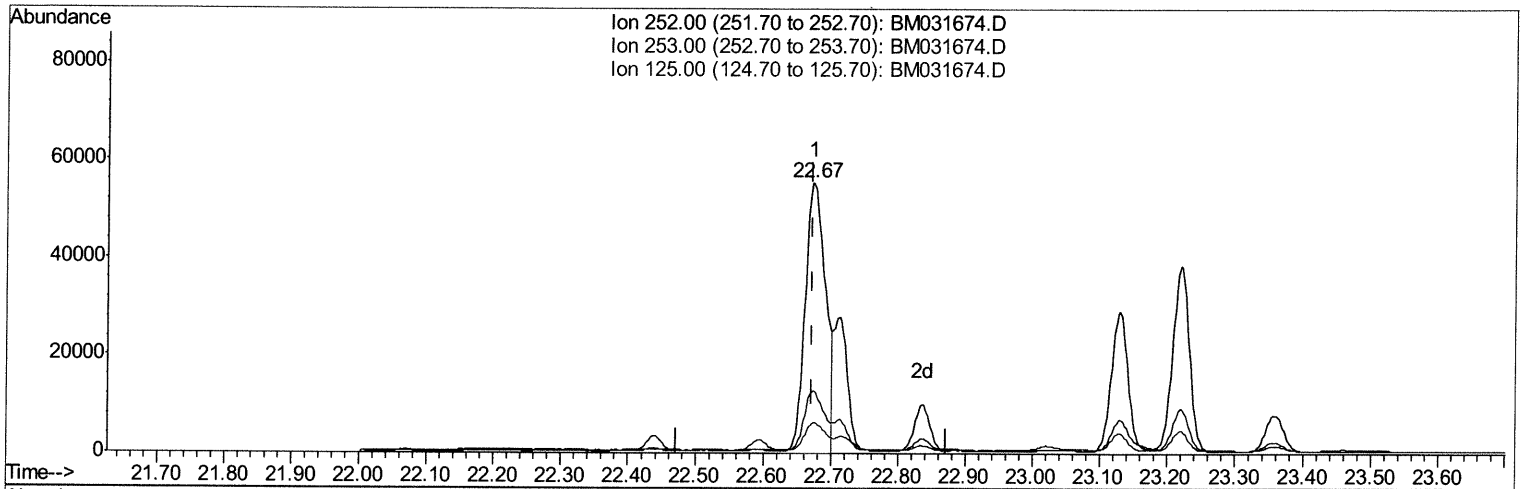
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 Data File : BM031674.D
 Acq On : 12 Aug 2021 13:08
 Operator : CG/JU
 Sample : M3182-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00D1

Manual Integrations
 APPROVED

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TIC: BM031674.D

(24) Benzo(b)fluoranthene

22.674min (+0.002) 6.75ng/ul m *CG/JU*

response 115666

Ion	Exp%	Act%
252.00	100	100
253.00	26.50	23.23
125.00	11.40	11.20
0.00	0.00	0.00

Quantitation Report (Qedit)

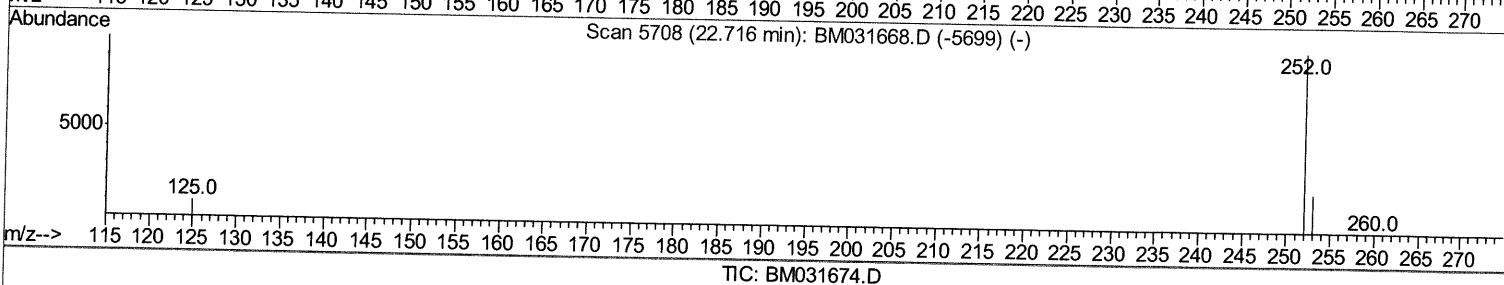
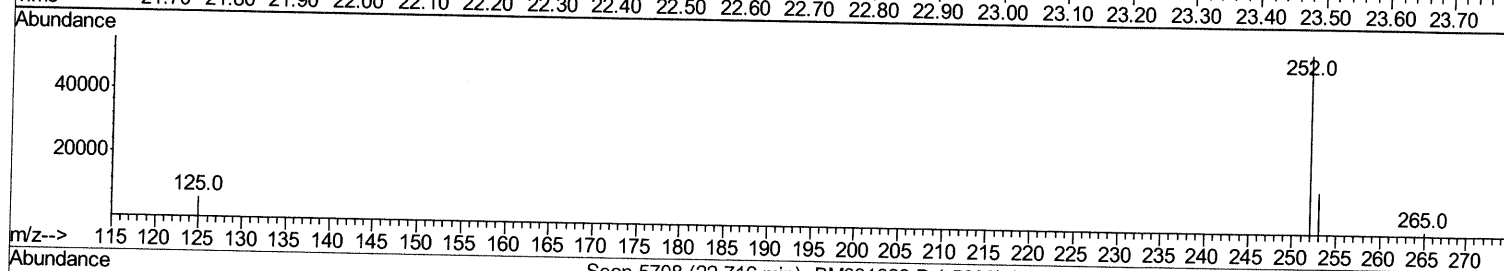
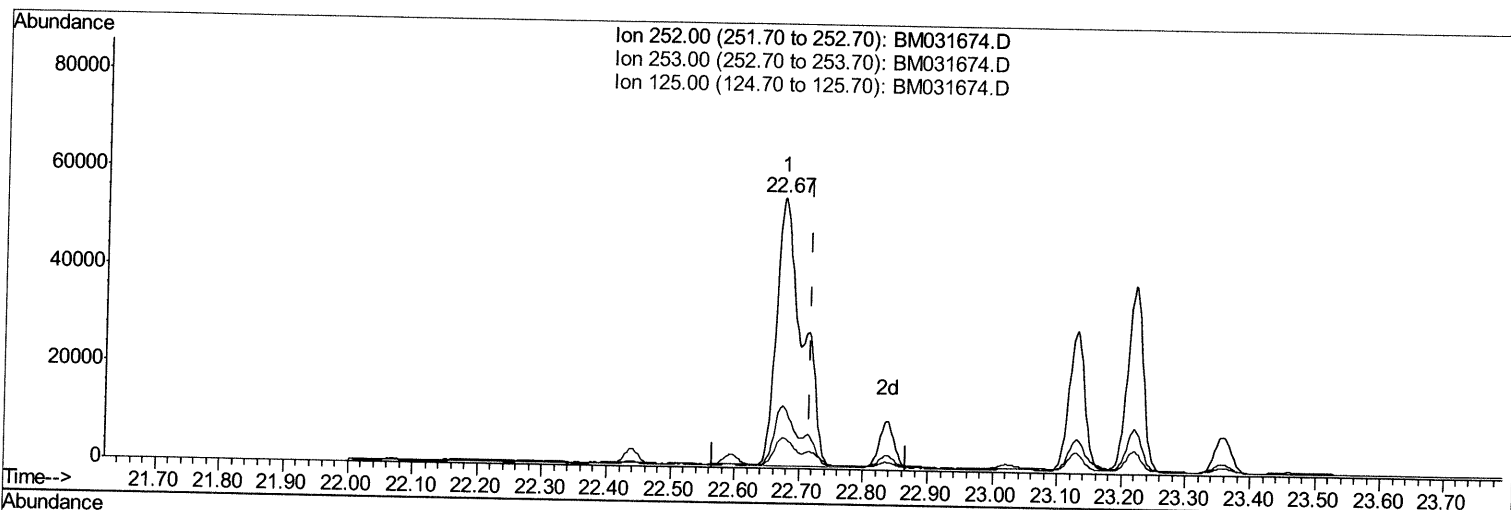
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\
 Data File : BM031674.D
 Acq On : 12 Aug 2021 13:08
 Operator : CG/JU
 Sample : M3182-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Manual Integrations
 APPROVED

mohammad
 8/13/2021 5:13:48 PM

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(25) Benzo(k)fluoranthene
 22.674min (-0.042) 8.70ng/ul
 response 153980

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	23.23
125.00	11.00	11.20
0.00	0.00	0.00

Quantitation Report (Qedit)

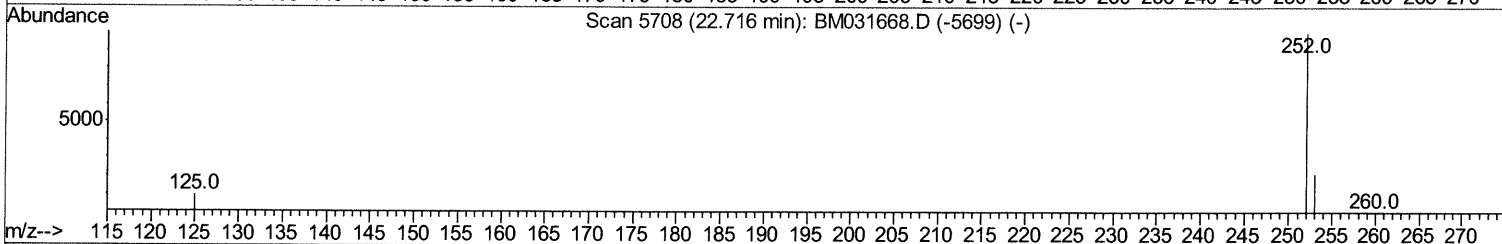
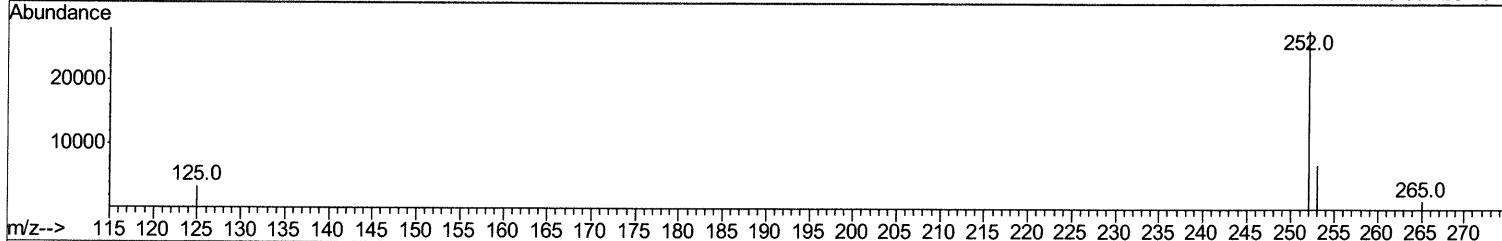
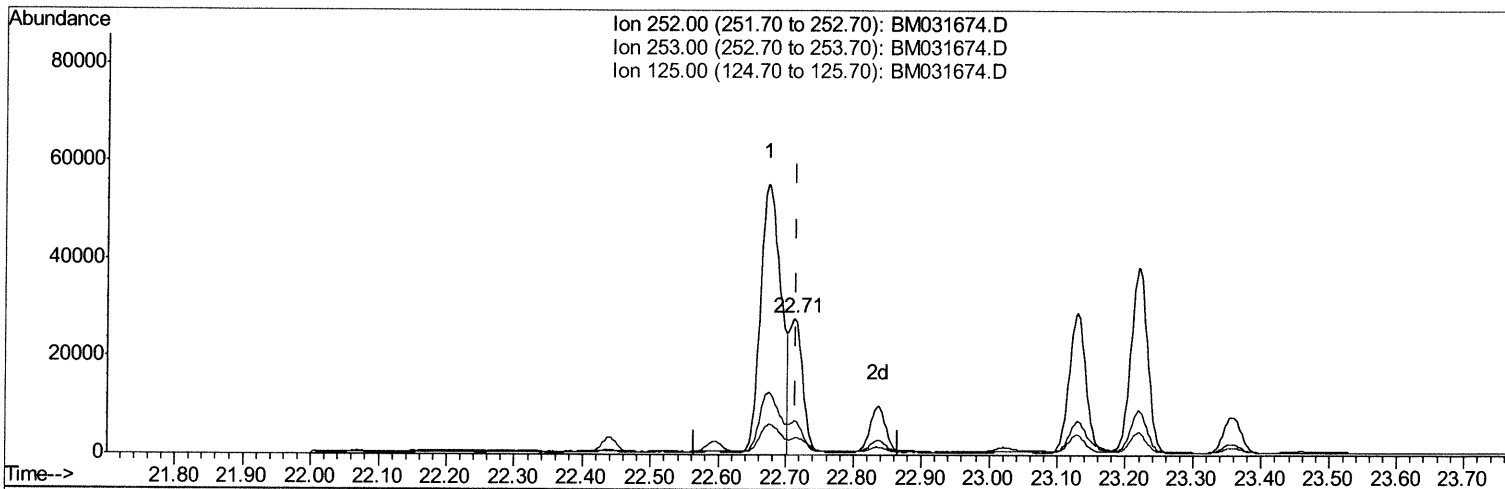
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\
 Data File : BM031674.D
 Acq On : 12 Aug 2021 13:08
 Operator : CG/JU
 Sample : M3182-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00D1

Manual Integrations
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TIC: BM031674.D

(25) Benzo(k)fluoranthene

22.713min (-0.003) 2.34ng/ul m

08/16/21 JU

response 41462

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	24.80
125.00	11.00	12.25
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\
 Data File : BM031674.D
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 Operator : CG/JU
 Sample : M3182-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1234	0.40	ng/ul	0.00
4) Naphthalene-d8	10.33	136	4579	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.20	164	2810	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.96	188	4950	0.40	ng/ul	0.00
17) Chrysene-d12	21.16	240	3928	0.40	ng/ul	0.00
23) Perylene-d12	23.31	264	3801	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.18	96	3943	2.88	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.92	152	1618	0.21	ng/ul	0.00
18) Fluoranthene-d10	18.99	212	3191	0.24	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
5) Naphthalene	10.38	128	33409	2.449	ng/ul	97
7) 2-Methylnaphthalene	12.00	142	21140	2.269	ng/ul	99
8) 1-Methylnaphthalene	12.22	142	10434	1.133	ng/ul	100
10) Acenaphthylene	13.93	152	13107	1.095	ng/ul	98
11) Acenaphthene	14.26	153	1421m>	0.141	ng/ul>	08/16/21 JU
12) Fluorene	15.26	166	2107	0.182	ng/ul#	90
14) Pentachlorophenol	16.61	266	45	0.028	ng/ul#	89
15) Phenanthrene	17.00	178	73260	4.661	ng/ul	98
16) Anthracene	17.09	178	14667	0.997	ng/ul	99
19) Fluoranthene	19.02	202	146584	8.739	ng/ul	97
20) Pyrene	19.39	202	123390	7.253	ng/ul	98
21) Benzo(a)anthracene	21.14	228	72286	4.557	ng/ul	96
22) Chrysene	21.19	228	83400	5.082	ng/ul	98
24) Benzo(b)fluoranthene	22.67	252	115666m>	6.754	ng/ul>	08/16/21 JU
25) Benzo(k)fluoranthene	22.71	252	41462m>	2.343	ng/ul>	08/16/21 JU
26) Benzo(a)pyrene	23.22	252	65479	4.366	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	25.45	276	49426	2.549	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.45	278	12378	0.792	ng/ul#	93
29) Benzo(a,h,i)perylene	26.10	276	12895	0.780	ng/ul	94

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