

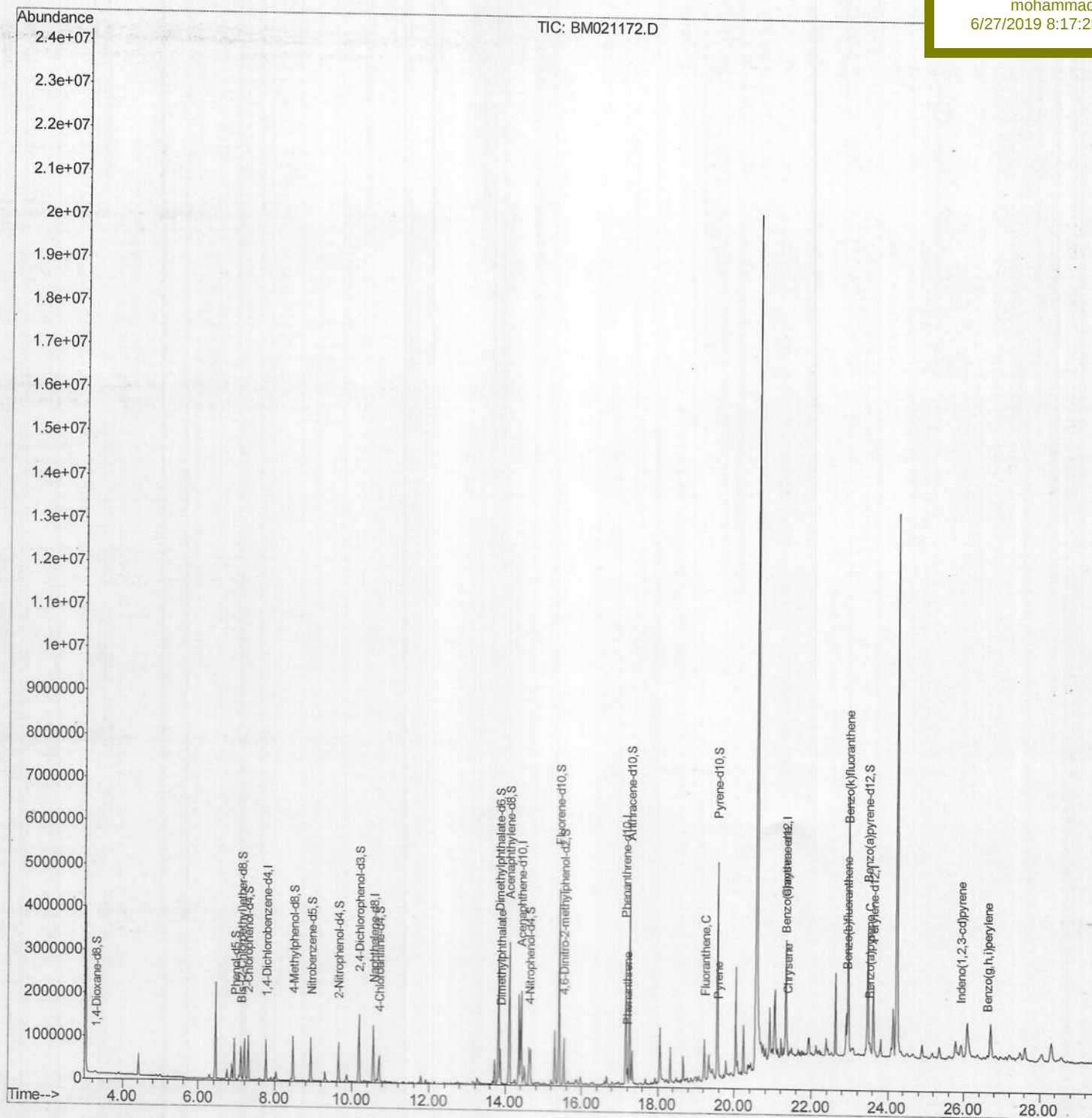
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062519\
Data File : BM021172.D
Acq On : 25 Jun 2019 21:40
Operator : HP/JU
Sample : K3498-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Jun 26 02:08:47 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 26 01:32:57 2019
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
C5AA2

Manual Integrations
APPROVED

mohammad
6/27/2019 8:17:28 AM



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062519\

Data File : BM021172.D

Acq On : 25 Jun 2019 21:40

Operator : HP/JU

Sample : K3498-03

Misc :

ALS Vial : 21 Sample Multiplier: 1

Quant Time: Jun 26 01:39:09 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M

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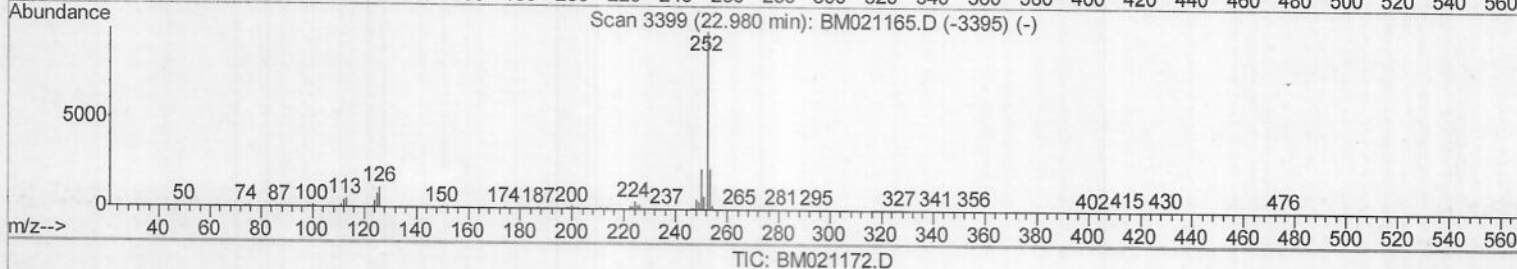
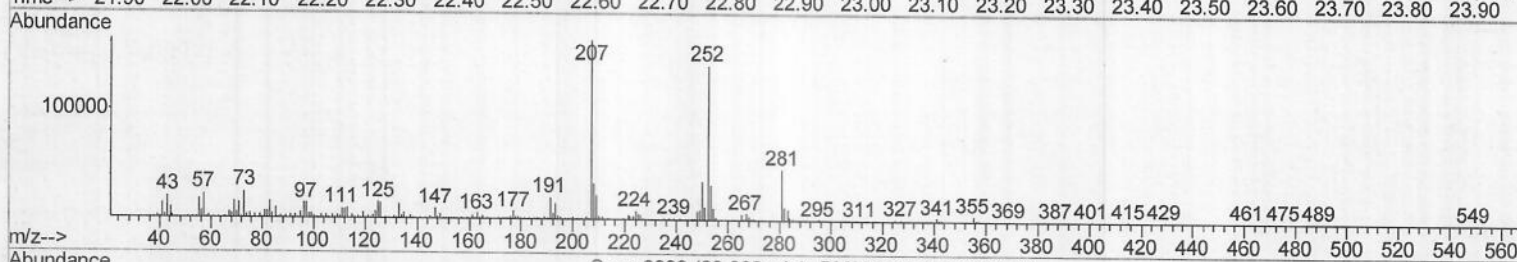
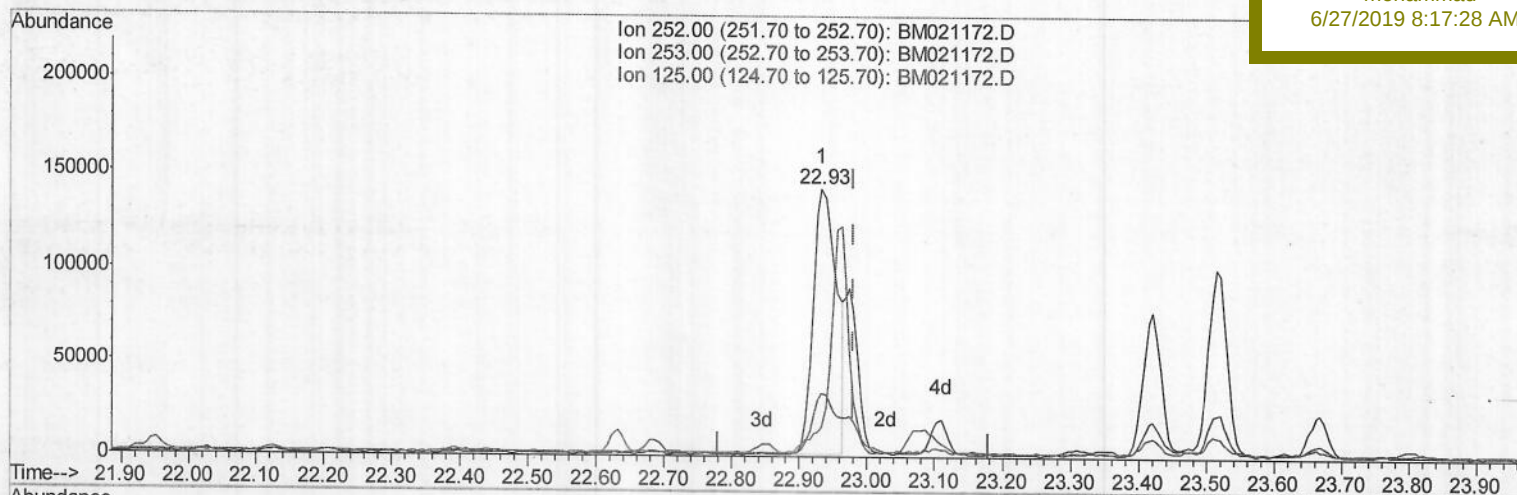
C5AA2

Manual Integrations

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mohammad

6/27/2019 8:17:28 AM



(88) Benzo(k)fluoranthene

22.933min (-0.047) 4.58ng/ul

response 337142

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	23.51
125.00	8.20	11.06#
0.00	0.00	0.00

Quantitation Report (Qedit)

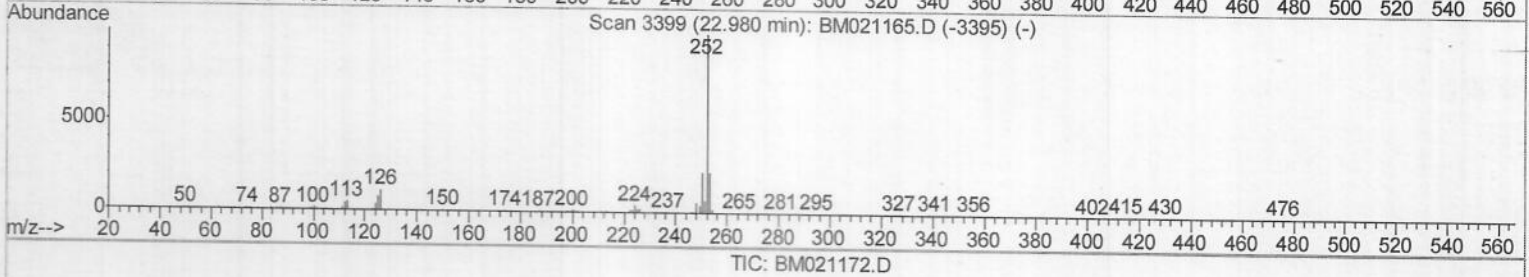
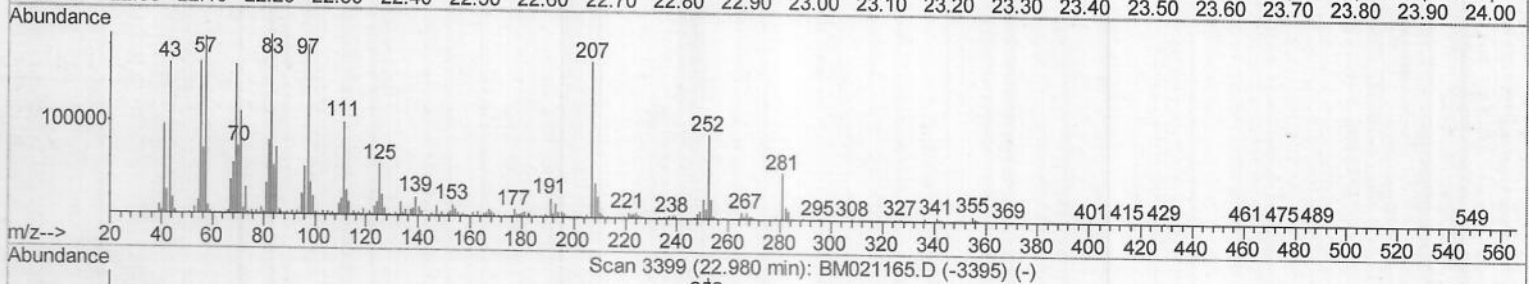
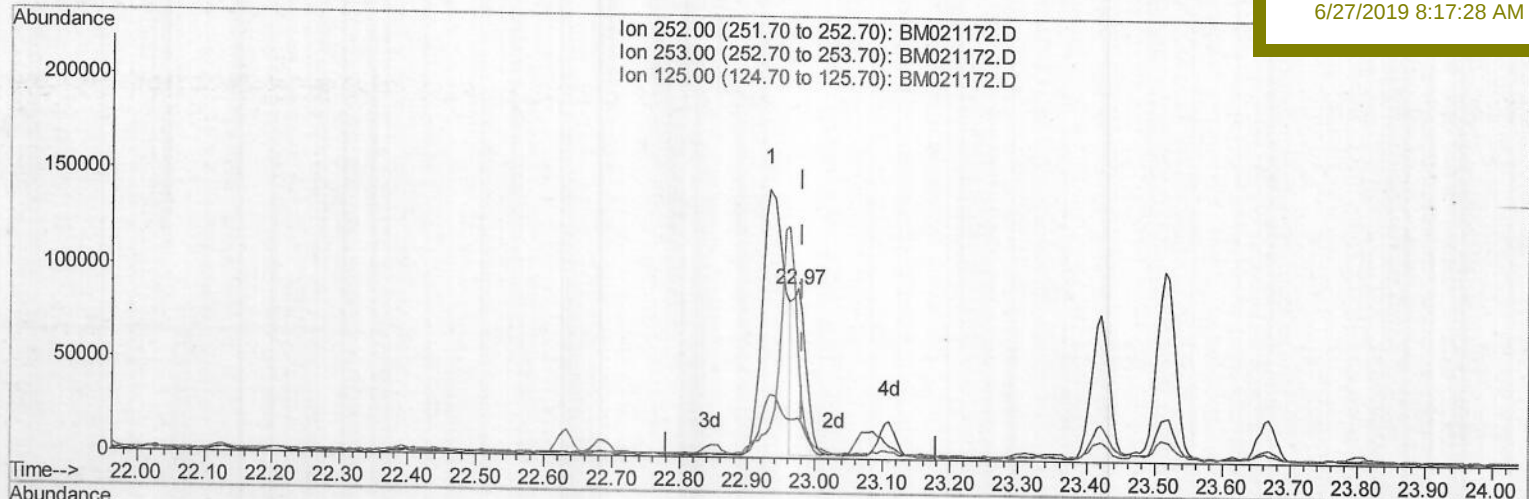
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 BNA_M
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 C5AA2

Manual Integrations
 APPROVED

mohammad
 6/27/2019 8:17:28 AM



(88) Benzo(k)fluoranthene

22.974min (-0.006) 1.66ng/ul m > JU 06/24/19

response 122040

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	23.56
125.00	8.20	60.66#
0.00	0.00	0.00

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Manual Integrations
 APPROVED

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 6/27/2019 8:17:28 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	251700	20.00	ng/ul	0.00
18) Naphthalene-d8	10.55	136	1103836	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.41	164	708971	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	1621079	20.00	ng/ul	0.00
77) Chrysene-d12	21.34	240	1181494	20.00	ng/ul	0.00
85) Perylene-d12	23.61	264	1209109	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.29	96	22487	4.24	ng/uL	0.00
5) Phenol-d5	6.93	99	551964	25.19	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.11	67	330353	25.31	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	463407	26.03	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	393402	21.51	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	249615	27.85	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	297920	29.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.17	165	608290	30.07	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	353117	16.73	ng/ul	0.00
43) Dimethylphthalate-d6	13.82	166	2066520	32.26	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	2334834	29.77	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	237662	22.81	ng/ul	0.00
57) Fluorene-d10	15.40	176	1755908	31.51	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	238415	23.92	ng/ul	0.00
70) Anthracene-d10	17.26	188	2605026	30.92	ng/ul	0.00
78) Pyrene-d10	19.55	212	2477825	34.95	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.47	264	2127700	29.83	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
44) Dimethylphthalate	13.87	163	504590	8.653	ng/ul	99
69) Phenanthrene	17.20	178	186579	2.081	ng/ul	99
76) Fluoranthene	19.22	202	525860	5.428	ng/ul	100
79) Pyrene	19.58	202	421854	5.051	ng/ul	99
82) Benzo(a)anthracene	21.33	228	217618	2.702	ng/ul	96
84) Chrysene	21.38	228	262543	3.480	ng/ul	97
87) Benzo(b)fluoranthene	22.93	252	337142	4.402	ng/ul#	95
88) Benzo(k)fluoranthene	22.97	252	122040m	1.656	ng/ul	
90) Benzo(a)pyrene	23.52	252	190784	2.647	ng/ul#	98
91) Indeno(1,2,3-cd)pyrene	25.90	276	144823	1.706	ng/ul	97
93) Benzo(g,h,i)perylene	26.61	276	114819	1.642	ng/ul	98

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

JU/06/24/19