

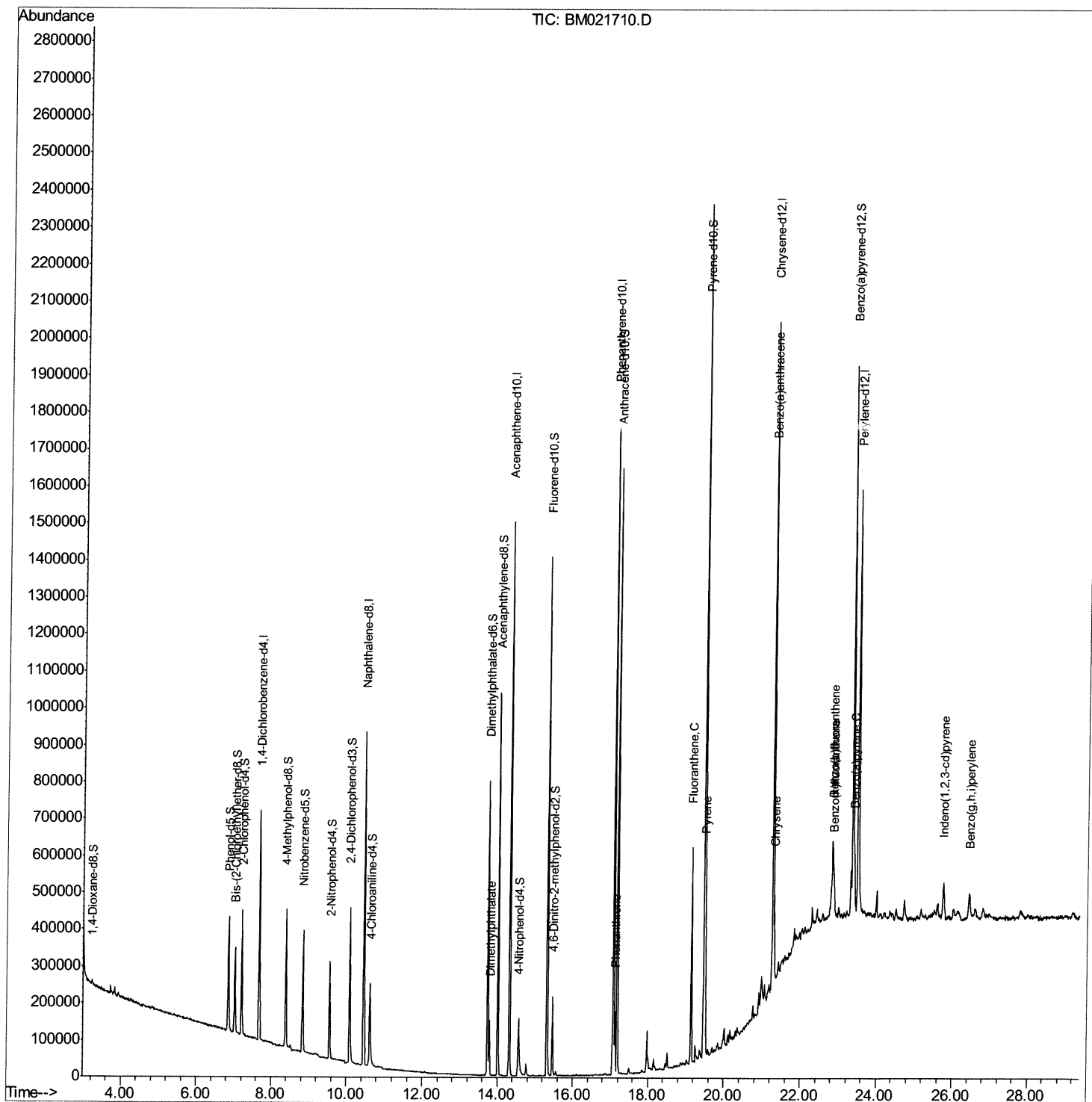
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM072919\
Data File : BM021710.D
Acq On : 30 Jul 2019 01:14
Operator : HP/JU
Sample : K3863-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleID :
A52F2

Manual Integrations
APPROVED

mohammad
7/30/2019 4:51:44 PM

Quant Time: Jul 30 05:28:45 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072619MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 26 18:42:09 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

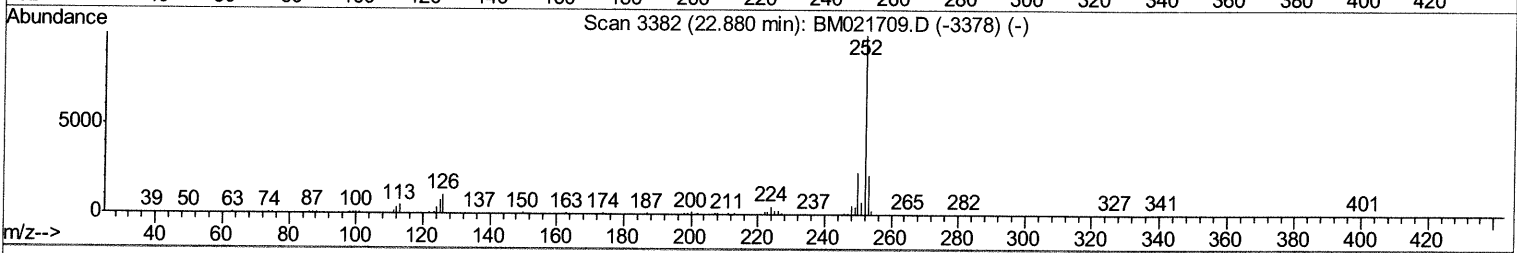
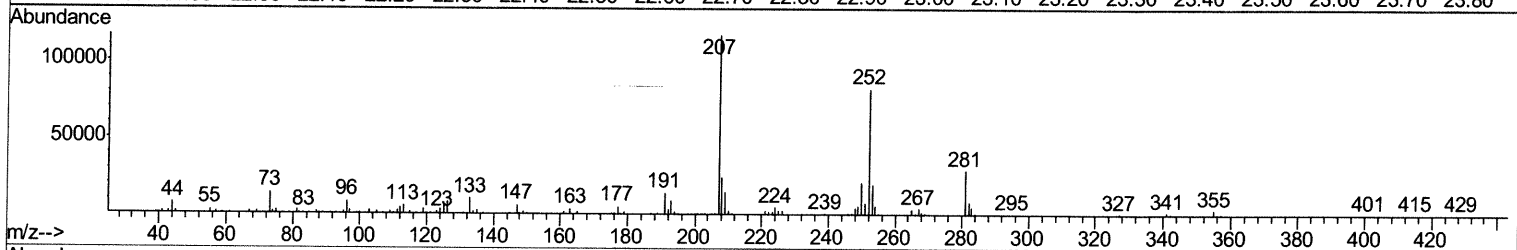
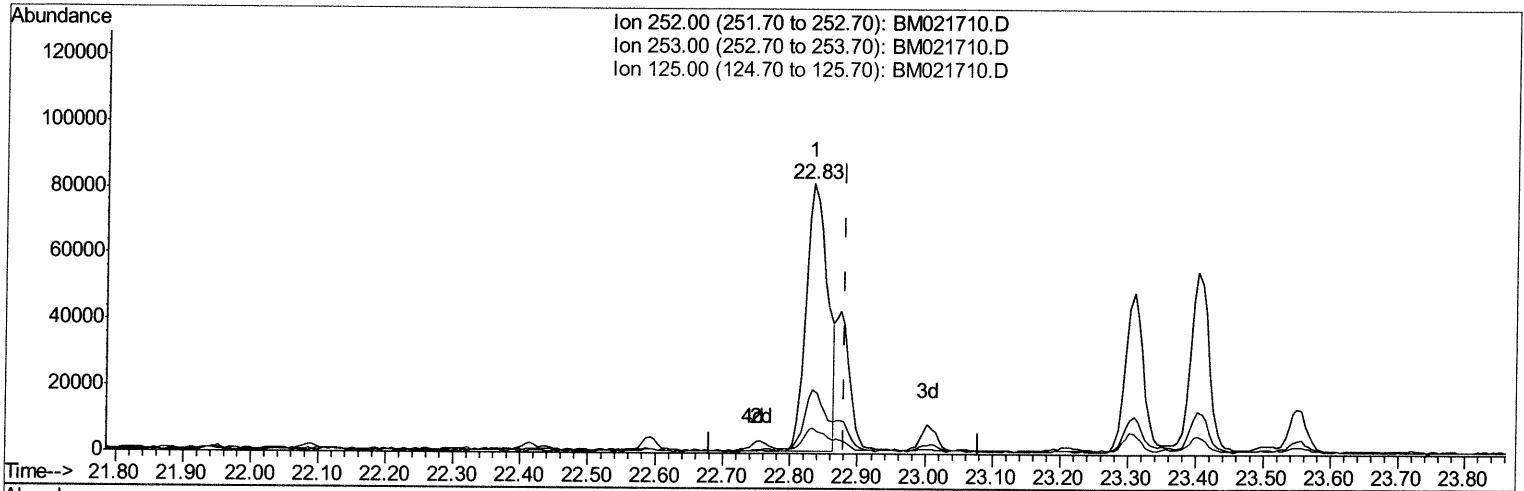
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TIC: BM021710.D

(88) Benzo(k)fluoranthene

22.833min (-0.047) 3.21ng/ul

response 179344

Ion	Exp%	Act%
252.00	100	100
253.00	22.10	23.49
125.00	7.90	9.57#
0.00	0.00	0.00

Quantitation Report (Qedit)

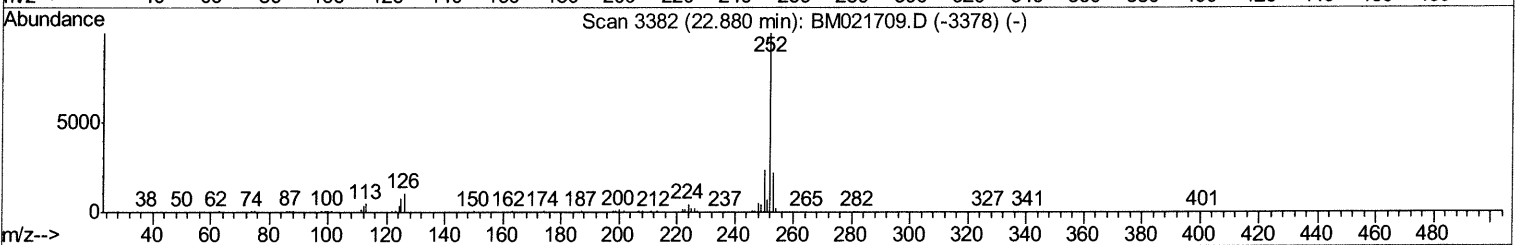
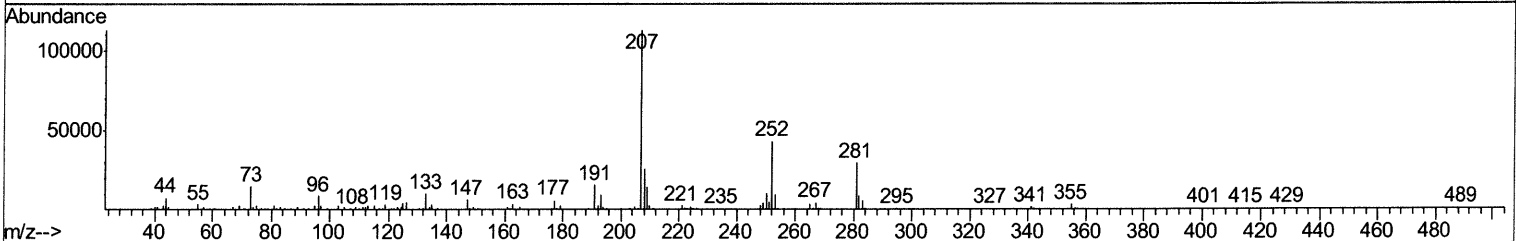
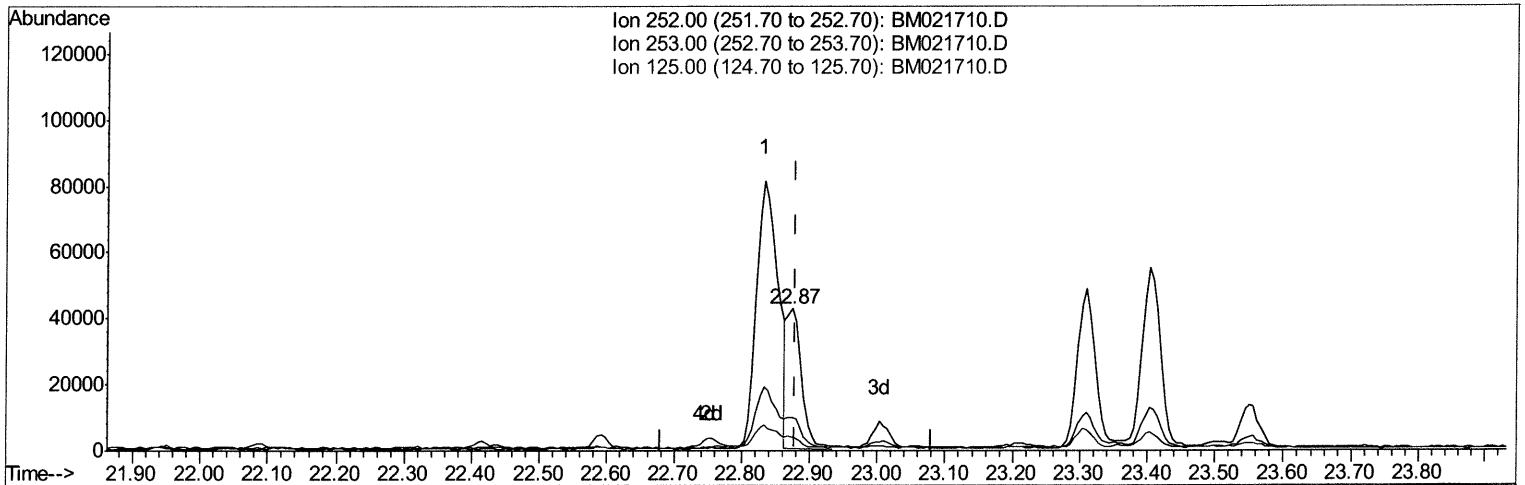
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Instrument :
 BNA_M
 ClientSampled :
 A52F2

Manual Integrations
 APPROVED

mohammad
 7/30/2019 4:51:44 PM

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 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072619MA.M
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 QLast Update : Fri Jul 26 18:42:09 2019
 Response via : Initial Calibration



TIC: BM021710.D

(88) Benzo(k)fluoranthene

22.874min (-0.006) 1.13ng/ul m 07/31/19

response 63202

Ion	Exp%	Act%
252.00	100	100
253.00	22.10	23.24
125.00	7.90	9.90#
0.00	0.00	0.00

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 Operator : HP/JU
 Sample : K3863-12
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A52F2

Manual Integrations
 APPROVED

mohammad
 7/30/2019 4:51:44 PM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	175046	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	733696	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.32	164	475890	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.07	188	1000836	20.00	ng/ul	0.00
77) Chrysene-d12	21.27	240	778792	20.00	ng/ul	0.00
85) Perylene-d12	23.50	264	839954	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.23	96	8176	2.25	ng/uL	0.00
5) Phenol-d5	6.86	99	181991	13.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.03	67	107382	14.24	ng/ul	0.00
9) 2-Chlorophenol-d4	7.21	132	151721	13.89	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	143243	12.96	ng/ul	0.00
19) Nitrobenzene-d5	8.85	128	76255	13.90	ng/ul	0.00
22) 2-Nitrophenol-d4	9.56	143	83251	13.78	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	179458	13.87	ng/ul	0.00
29) 4-Chloroaniline-d4	10.62	131	151223	12.36	ng/ul	0.00
43) Dimethylphthalate-d6	13.74	166	550862	15.00	ng/ul	0.00
46) Acenaphthylene-d8	14.02	160	674265	15.45	ng/ul	0.00
51) 4-Nitrophenol-d4	14.56	143	51143	9.17	ng/ul	0.00
57) Fluorene-d10	15.32	176	543542	16.87	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.46	200	47120	7.57	ng/ul	0.00
70) Anthracene-d10	17.17	188	957052	19.00	ng/ul	0.00
78) Pyrene-d10	19.47	212	1160149	26.56	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.36	264	1036176	23.32	ng/ul	0.00
Target Compounds						
44) Dimethylphthalate	13.79	163	99777	2.617	ng/ul	99
69) Phenanthrene	17.12	178	101848	1.690	ng/ul	99
76) Fluoranthene	19.14	202	328245	4.668	ng/ul	99
79) Pyrene	19.50	202	272733	4.693	ng/ul	98
82) Benzo(a)anthracene	21.26	228	95445	1.683	ng/ul	96
84) Chrysene	21.30	228	137919	2.580	ng/ul	99
87) Benzo(b)fluoranthene	22.83	252	179344	3.138	ng/ul#	96
88) Benzo(k)fluoranthene	22.87	252	63202m	1.131	ng/ul →	96
90) Benzo(a)pyrene	23.40	252	102064	1.907	ng/ul	96
91) Indeno(1,2,3-cd)pyrene	25.74	276	88671	1.429	ng/ul	98
93) Benzo(g,h,i)perylene	26.43	276	89671	1.738	ng/ul	97

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