

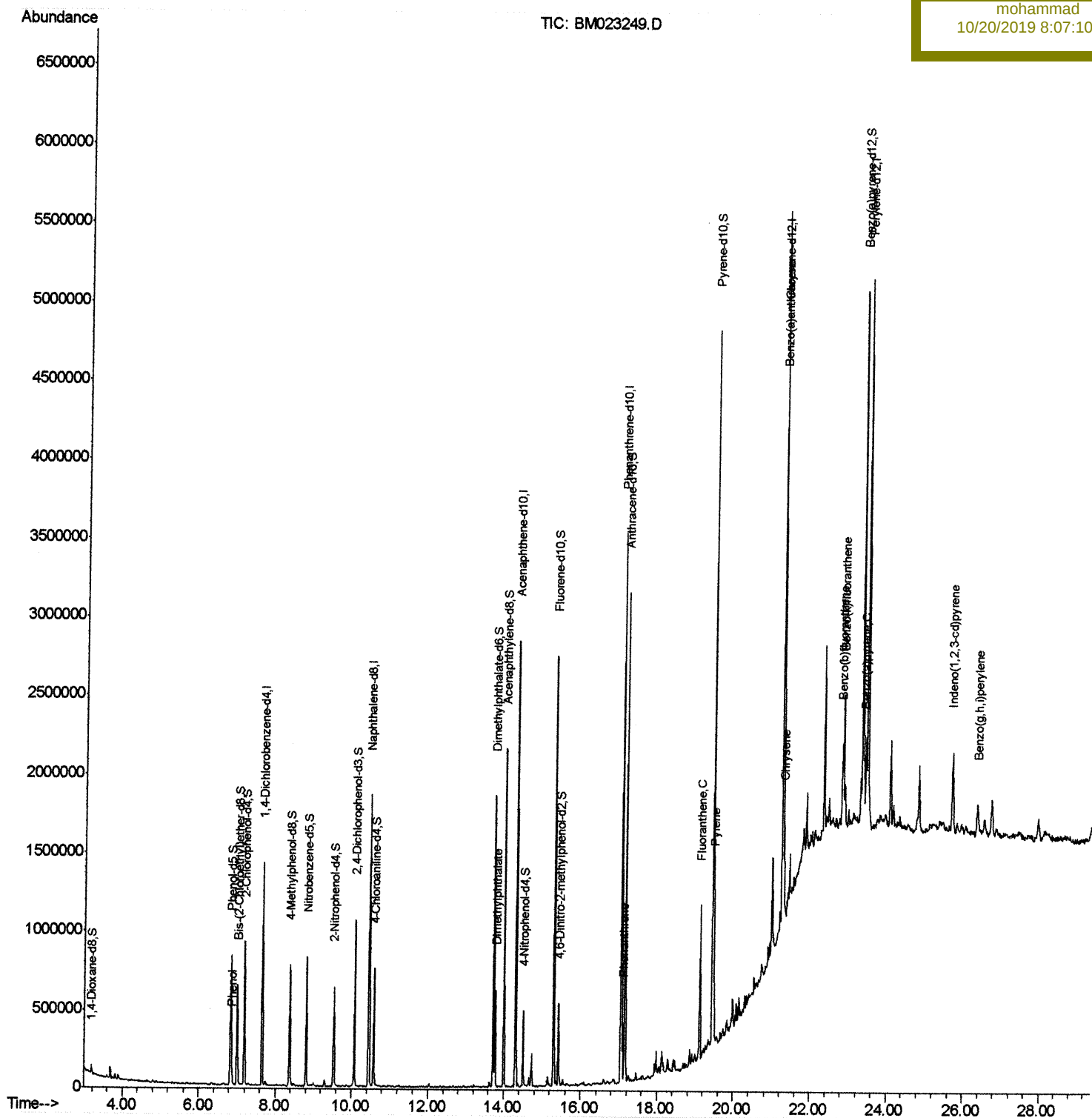
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101719\
Data File : BM023249.D
Acq On : 18 Oct 2019 09:27
Operator : JU
Sample : K5235-19
Misc :
ALS Vial : 28 Sample Multiplier: 1

Quant Time: Oct 18 15:14:58 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 07:48:41 2019
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
BEPB9

Manual Integrations
APPROVED

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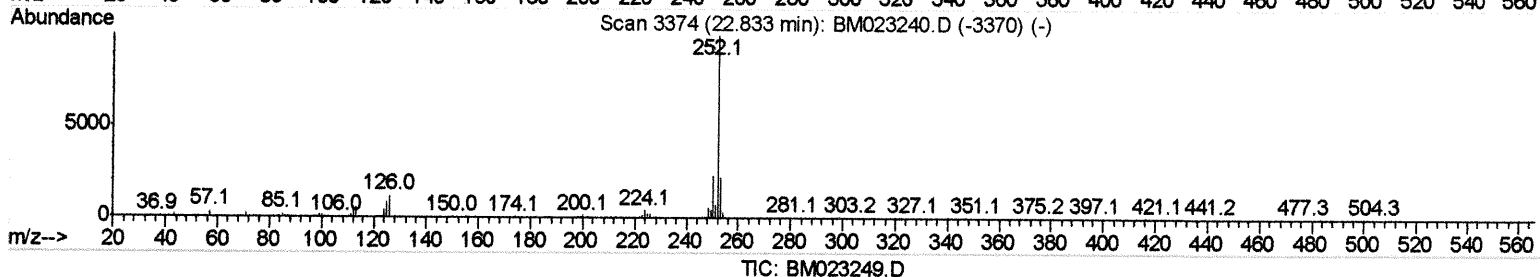
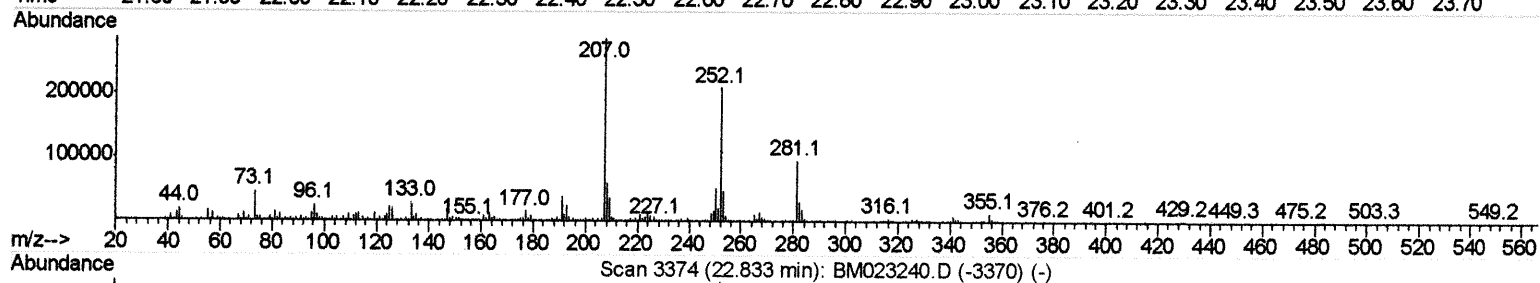
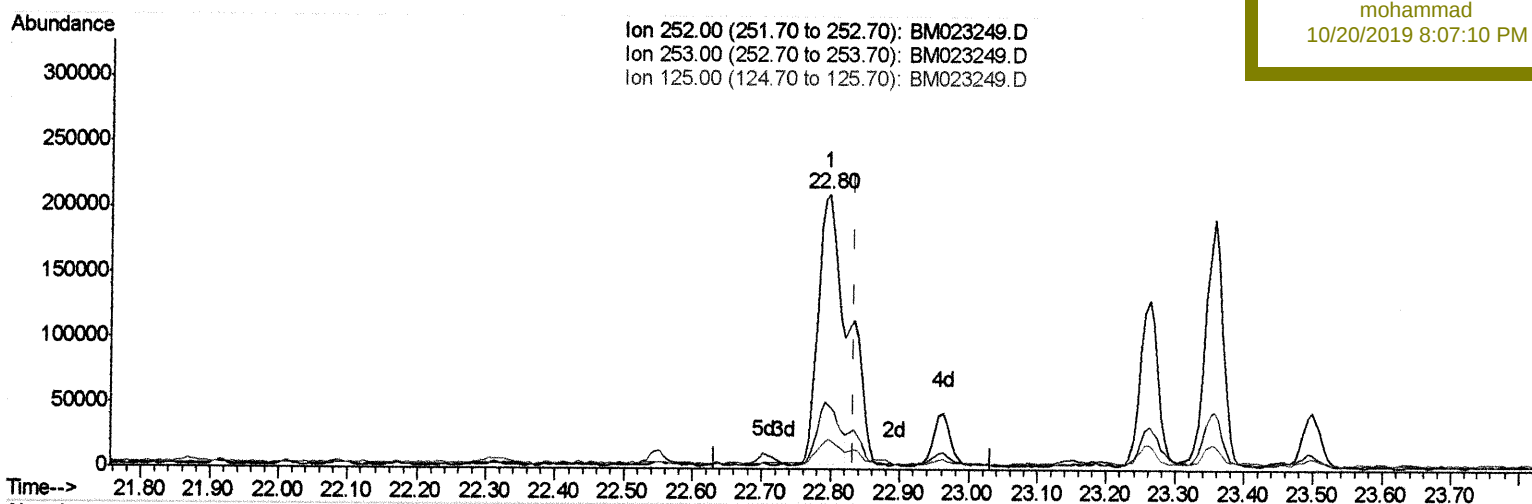
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(89) Benzo(k)fluoranthene

22.797min (-0.036) 3.24ng/ul

response 464416

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	23.09
125.00	8.10	10.72#
0.00	0.00	0.00

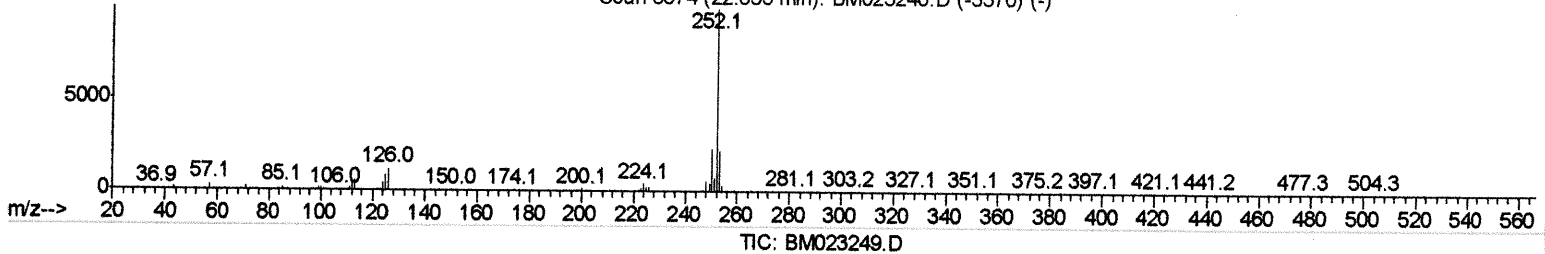
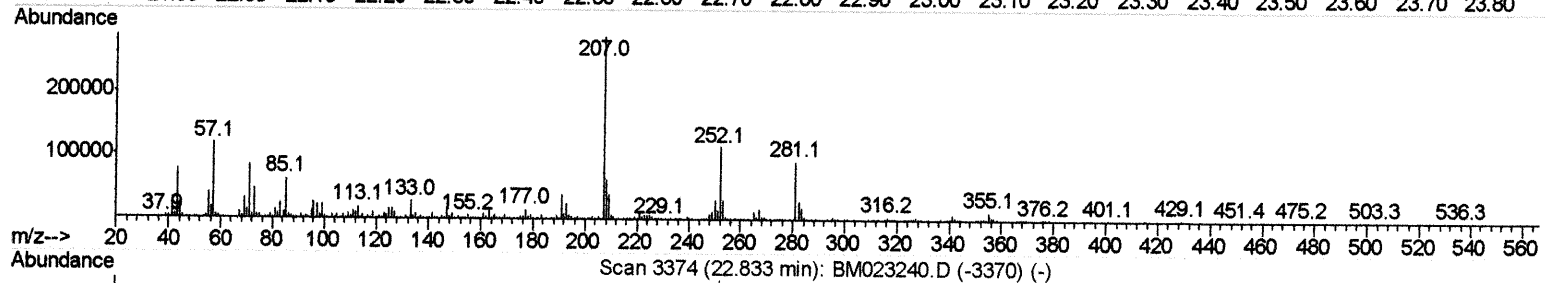
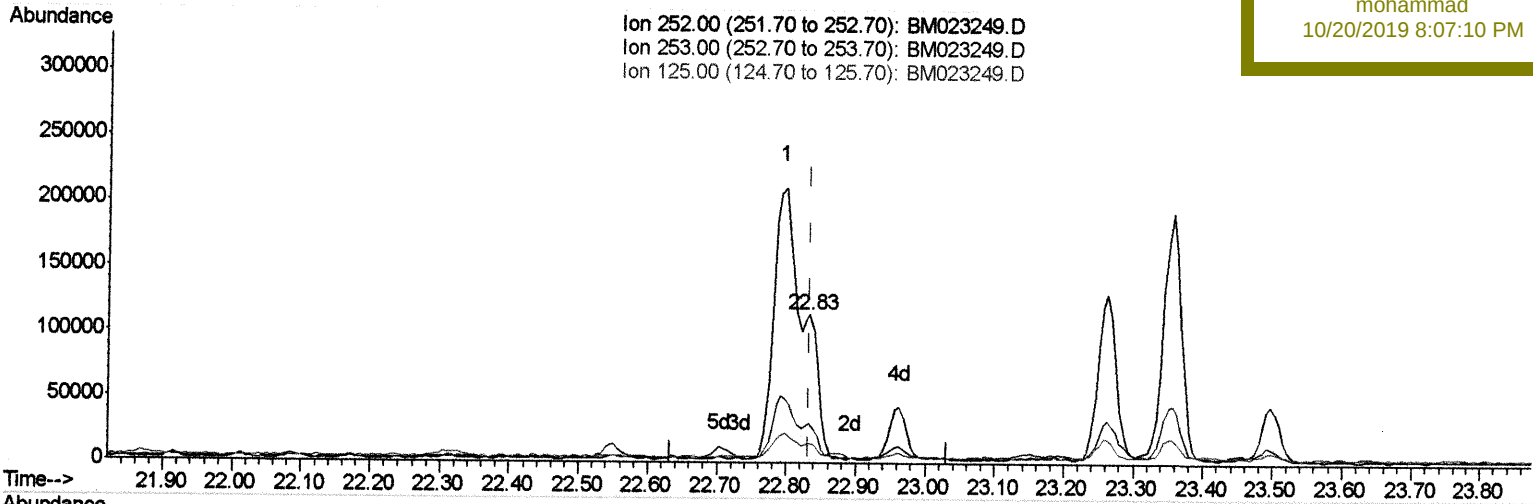
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Manual Integrations
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(89) Benzo(k)fluoranthene

22.833min (-0.000) 1.08ng/ul m

response 154104

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	26.87#
125.00	8.10	13.89#
0.00	0.00	0.00

ju 10/4/19

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 Sample : K5235-19
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Quant Time: Oct 18 15:14:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 18 07:48:41 2019
 Response via : Initial Calibration

Instrument :
 BNA_M
 ClientSampleId :
 BEPB9

Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	391672	20.00	ng/ul	0.00
18) Naphthalene-d8	10.44	136	1535408	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.30	164	951728	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.05	188	1993192	20.00	ng/ul	0.00
78) Chrysene-d12	21.24	240	2046471	20.00	ng/ul	0.00
86) Perylene-d12	23.45	264	2461190	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	27873	3.13	ng/uL	0.00
5) Phenol-d5	6.83	99	452797	13.27	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	284169	14.72	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	402399	15.41	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	292188	10.66	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	196085	18.88	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	211593	21.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	392229	15.59	ng/ul	0.00
29) 4-Chloroaniline-d4	10.57	131	403138	13.17	ng/ul	0.00
44) Dimethylphthalate-d6	13.72	166	1234528	16.93	ng/ul	0.00
47) Acenaphthylene-d8	13.99	160	1459825	17.07	ng/ul	0.00
52) 4-Nitrophenol-d4	14.50	143	114274	9.68	ng/ul	0.00
58) Fluorene-d10	15.30	176	1096576	17.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.42	200	115293	14.59	ng/ul	0.00
71) Anthracene-d10	17.14	188	1805119	18.78	ng/ul	0.00
79) Pyrene-d10	19.44	212	2237377	22.58	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.31	264	2621374	21.01	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	6.86	94	76516	2.175	ng/ul	95
45) Dimethylphthalate	13.77	163	391711	5.366	ng/ul	100
70) Phenanthrene	17.09	178	225828	2.019	ng/ul	99
77) Fluoranthene	19.11	202	596958	4.487	ng/ul	100
80) Pyrene	19.47	202	576164	4.494	ng/ul	99
83) Benzo(a)anthracene	21.23	228	324125	2.337	ng/ul	93
85) Chrysene	21.27	228	358195	2.782	ng/ul	97
88) Benzo(b)fluoranthene	22.80	252	464416	3.112	ng/ul#	96
89) Benzo(k)fluoranthene	22.83	252	154104m	1.075	ng/ul	96
91) Benzo(a)pyrene	23.36	252	339035	2.368	ng/ul	96
92) Indeno(1,2,3-cd)pyrene	25.66	276	240493	1.412	ng/ul	98
94) Benzo(a,h,i)perylene	26.34	276	230759	1.646	ng/ul	98

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