

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081320\

Data File : BM027102.D

Acq On : 15 Aug 2020 00:33

Operator : JU/CG

Sample : L3631-20

Misc :

ALS Vial : 48 Sample Multiplier: 1

Quant Time: Aug 17 03:03:36 2020

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM080720MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Aug 13 11:47:07 2020

Response via : Initial Calibration

Instrument :

BNA_M

Client Sample Id :

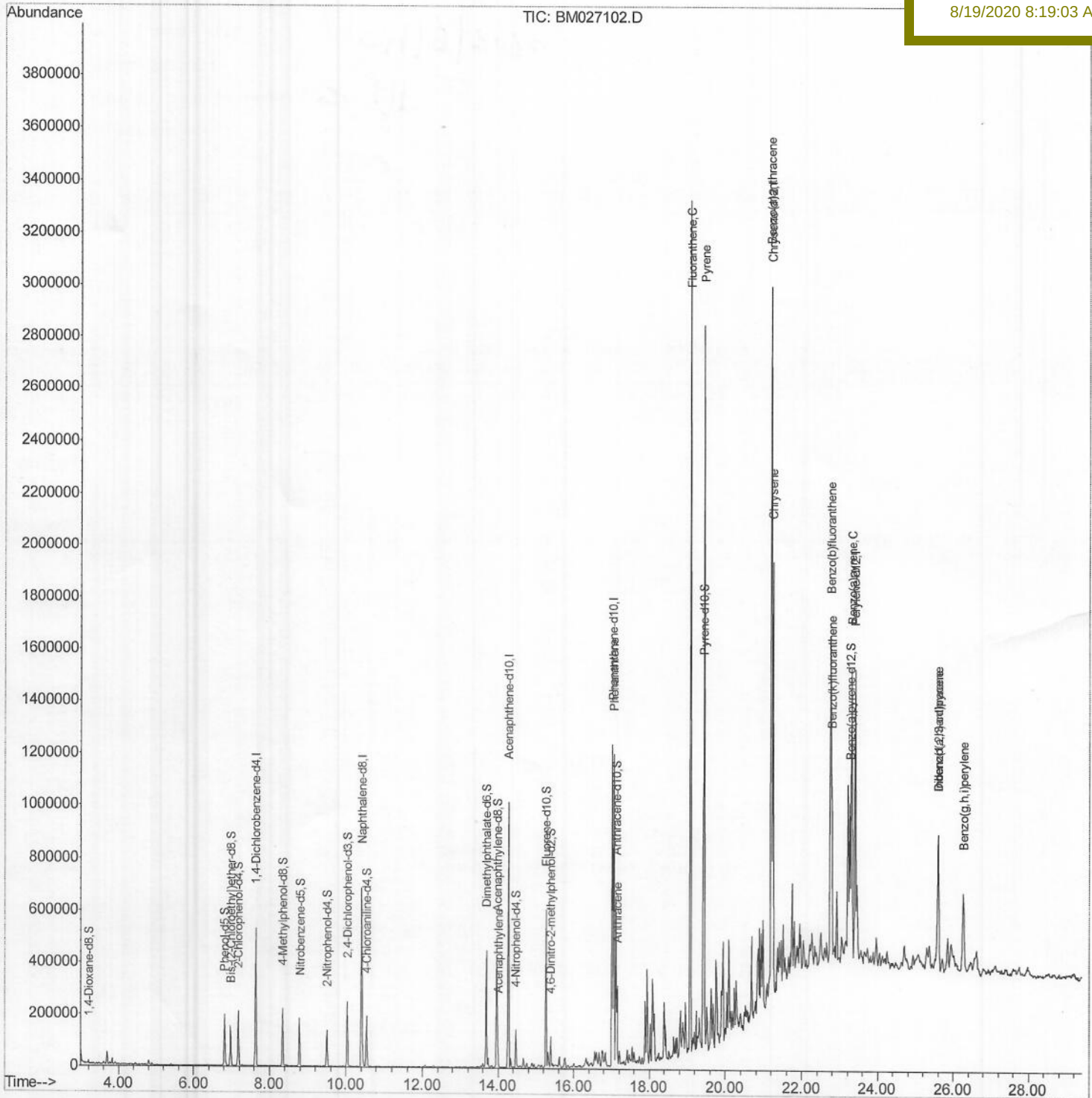
BFMN6

Manual Integrations

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mohammad

8/19/2020 8:19:03 AM



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Quant Title : SVOA CALIBRATION

QLast Update : Thu Aug 13 11:47:07 2020

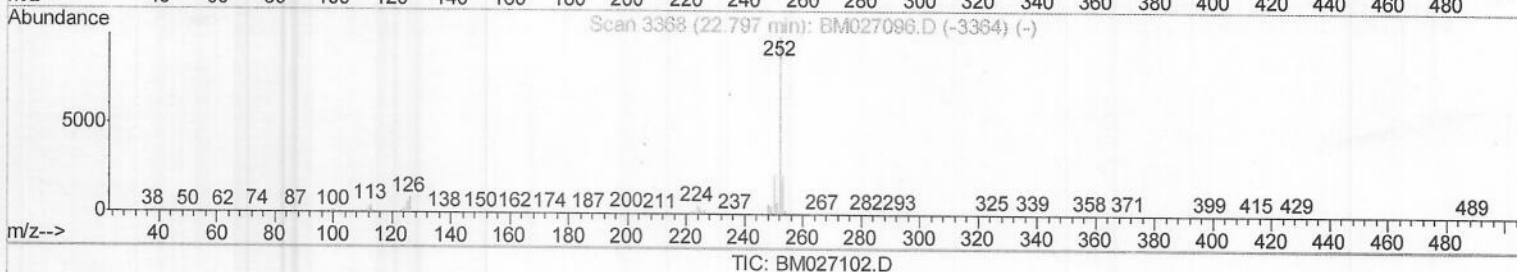
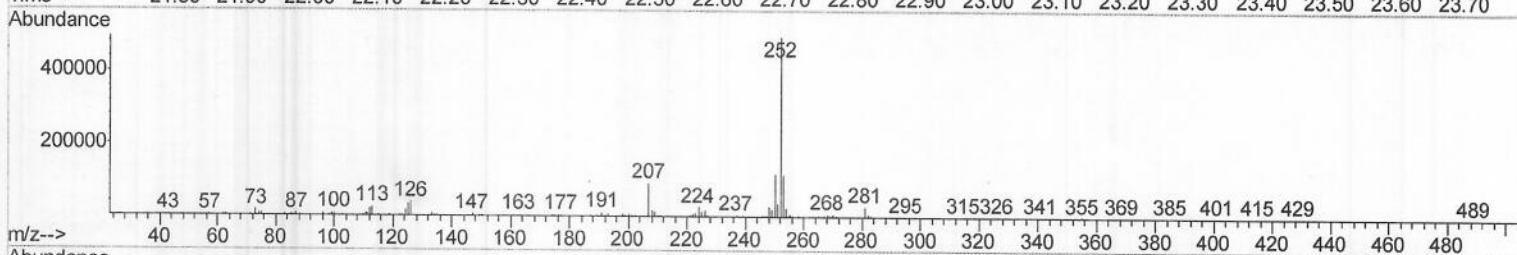
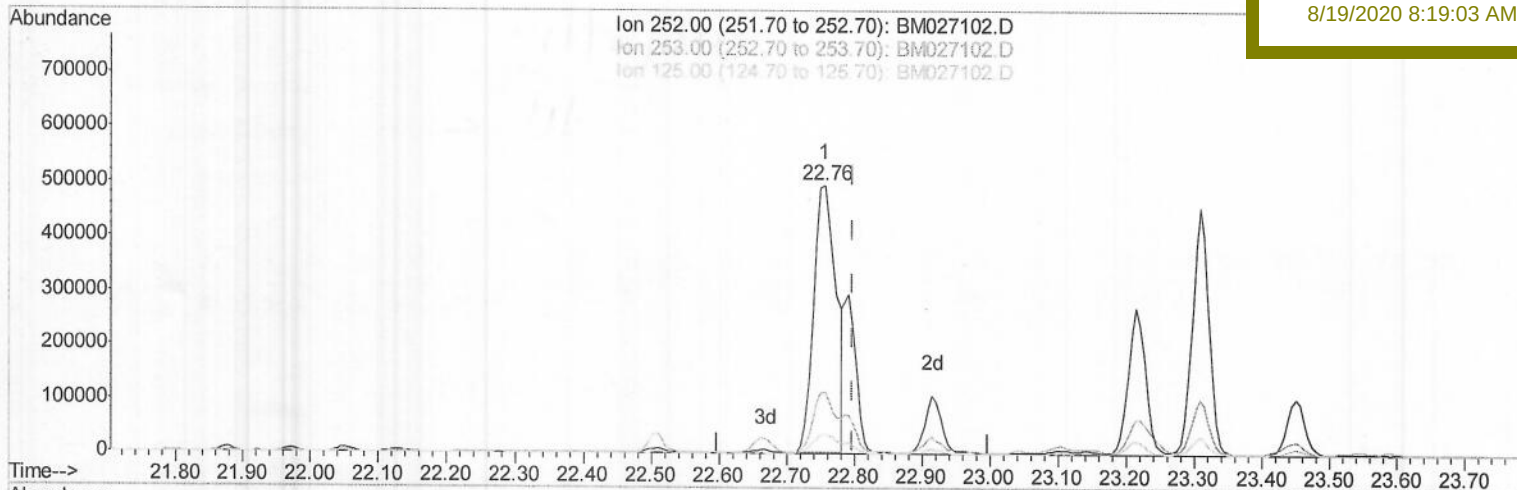
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Manual Integrations
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(89) Benzo(k)fluoranthene

22.756min (-0.041) 20.94ng/ul

response 1121962

Ion	Exp%	Act%
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252.00	100	100
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253.00	21.50	23.25
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125.00	6.90	7.09
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0.00	0.00	0.00
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Quantitation Report (Qedit)

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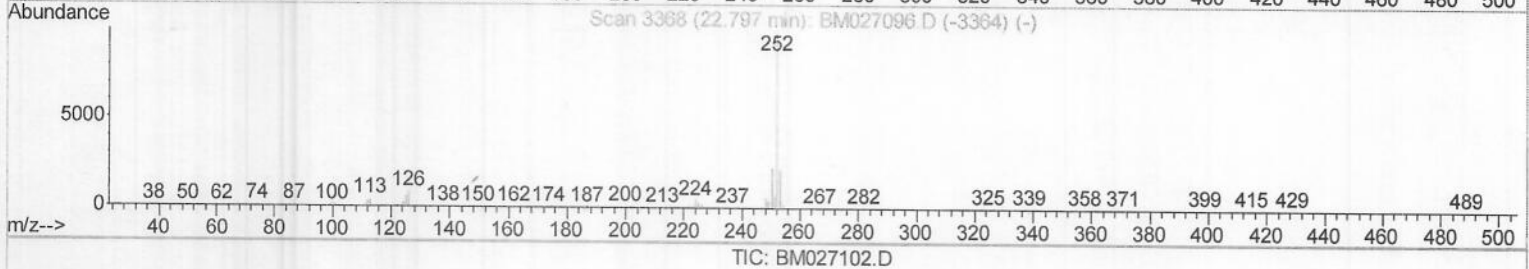
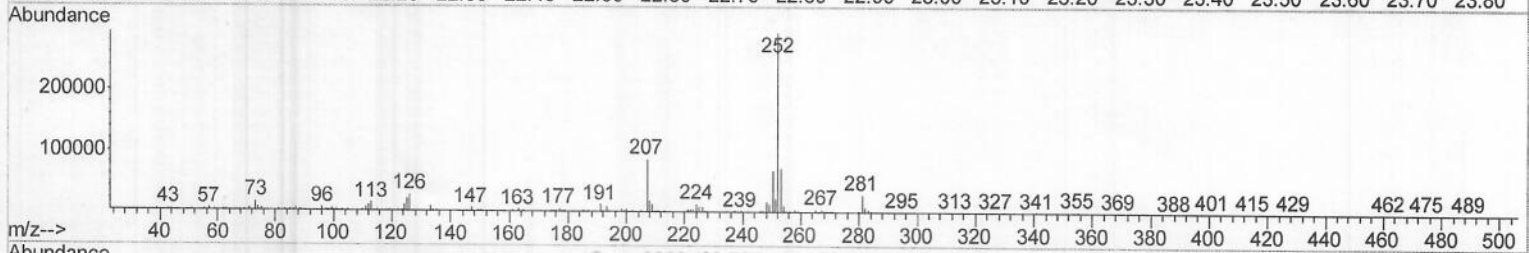
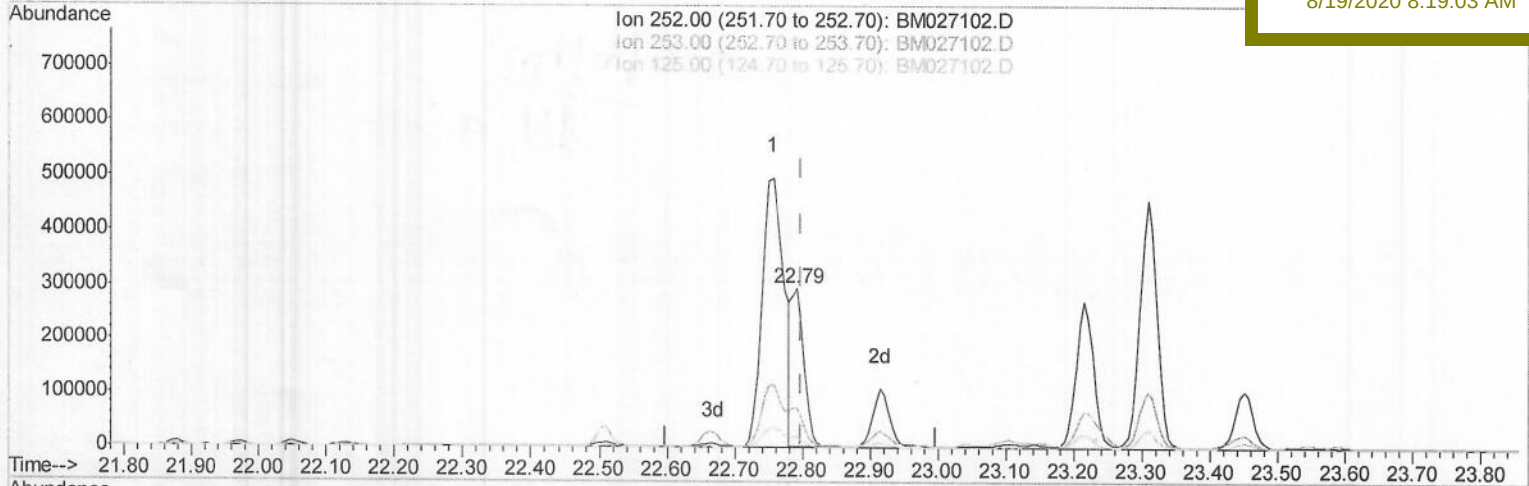
BFMN6

Manual Integrations

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(89) Benzo(k)fluoranthene

22.791min (-0.006) 6.90ng/ul m\ 08/25/20 JV

response 369575

Ion	Exp%	Act%
252.00	100	100
253.00	21.50	24.70
125.00	6.90	6.59
0.00	0.00	0.00

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 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BFMN6

Quant Time: Aug 17 03:03:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM080720MA.M
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 QLast Update : Thu Aug 13 11:47:07 2020
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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	7.64	152	139279	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	518286	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	333176	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	709151	20.00	ng/ul	0.00
78) Chrysene-d12	21.21	240	780807	20.00	ng/ul	0.00
86) Perylene-d12	23.40	264	778361	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.20	96	4730	1.54	ng/uL	0.00
5) Phenol-d5	6.82	99	99444	9.18	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.97	67	68735	9.27	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	79452	9.30	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	75678	8.96	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	38277	9.32	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	39193	9.21	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	83432	9.32	ng/ul	0.00
29) 4-Chloroaniline-d4	10.54	131	88298	8.49	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	273112	10.28	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	300704	9.50	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	30501	6.80	ng/ul	0.00
58) Fluorene-d10	15.27	176	223243	9.64	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	22917	5.92	ng/ul	0.00
71) Anthracene-d10	17.11	188	342388	9.80	ng/ul	0.00
79) Pyrene-d10	19.41	212	393714	10.11	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.26	264	428211	9.86	ng/ul	0.00

Target Compounds

					Ovalue	
48) Acenaphthylene	13.99	152	78735	2.446	ng/ul	98
70) Phenanthrene	17.06	178	672502	15.818	ng/ul	98
72) Anthracene	17.14	178	182257	4.213	ng/ul	99
77) Fluoranthene	19.08	202	1968325	38.180	ng/ul	97
80) Pyrene	19.44	202	1708292	32.233	ng/ul	99
83) Benzo(a)anthracene	21.19	228	1060968	20.184	ng/ul	96
85) Chrysene	21.24	228	877335	17.004	ng/ul	97
88) Benzo(b)fluoranthene	22.76	252	1121962	20.758	ng/ul	98
89) Benzo(k)fluoranthene	22.79	252	369575m	6.899	ng/ul	97
91) Benzo(a)pyrene	23.31	252	787875	16.017	ng/ul	97
92) Indeno(1,2,3-cd)pyrene	25.60	276	476732	7.522	ng/ul	98
93) Dibenzo(a,h)anthracene	25.60	278	136091	2.535	ng/ul#	90
94) Benzo(g,h,i)perylene	26.26	276	358236	6.843	ng/ul	98

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

> 08/25/20 JU