

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM093019\

Data File : BM022822.D

Acq On : 30 Sep 2019 13:15

Operator : JU

Sample : K4958-08

Misc :

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Oct 01 00:07:20 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Sep 27 18:03:27 2019

Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

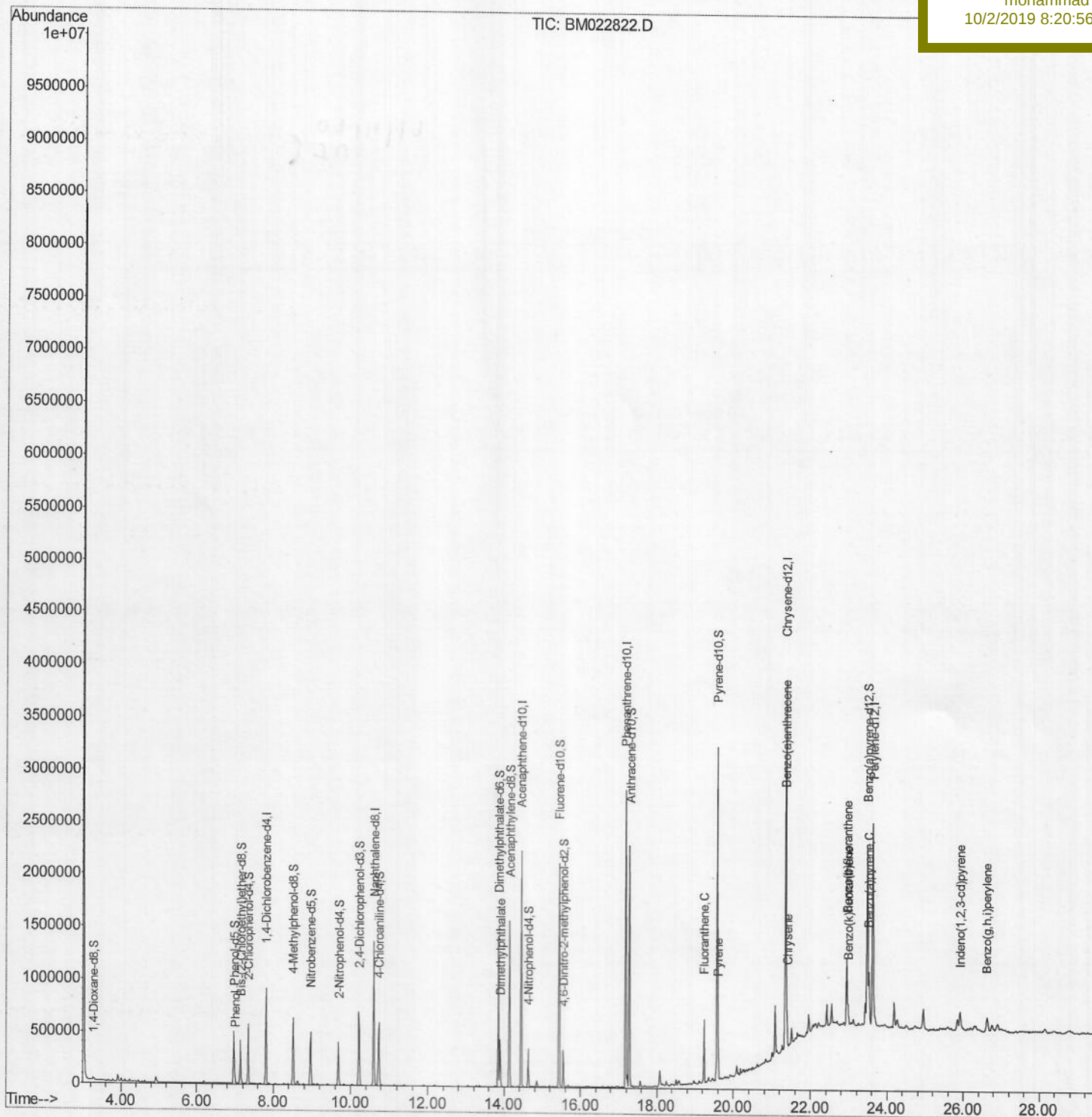
ESNY4

Manual Integrations

APPROVED

mohammad

10/2/2019 8:20:56 AM



Quantitation Report (Qedit)

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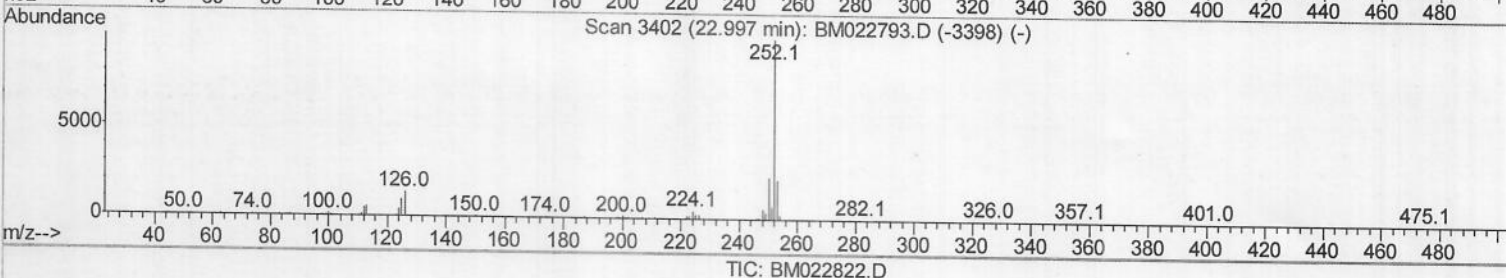
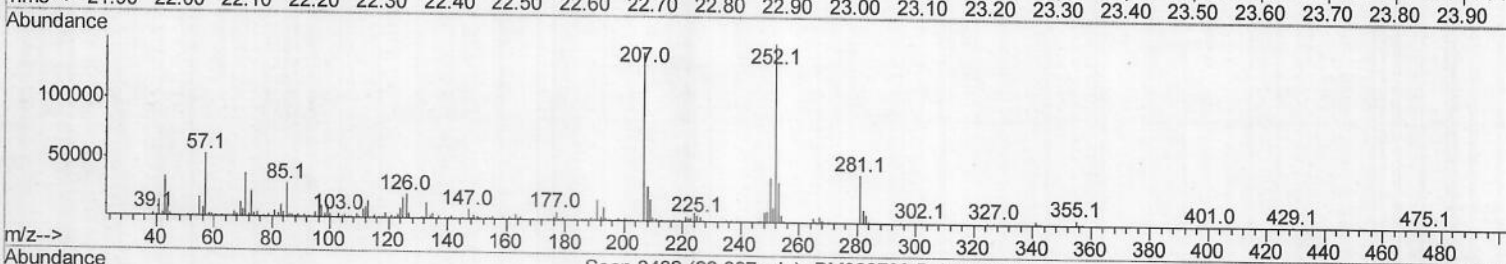
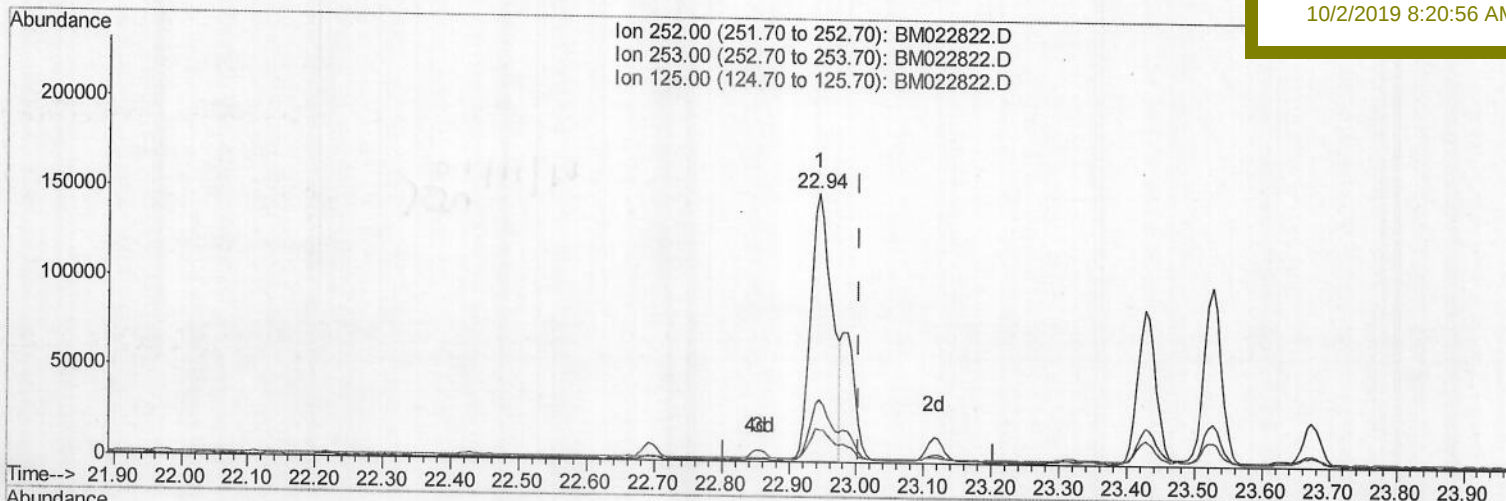
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10/2/2019 8:20:56 AM



TIC: BM022822.D

(89) Benzo(k)fluoranthene

22.944min (-0.059) 3.75ng/ul

response 328140

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	22.95
125.00	9.90	11.96#
0.00	0.00	0.00

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BNA_M

Client Sampled :

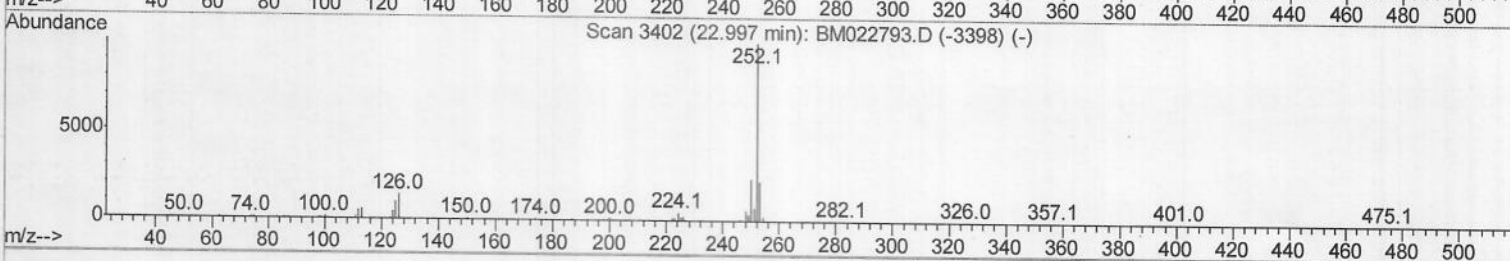
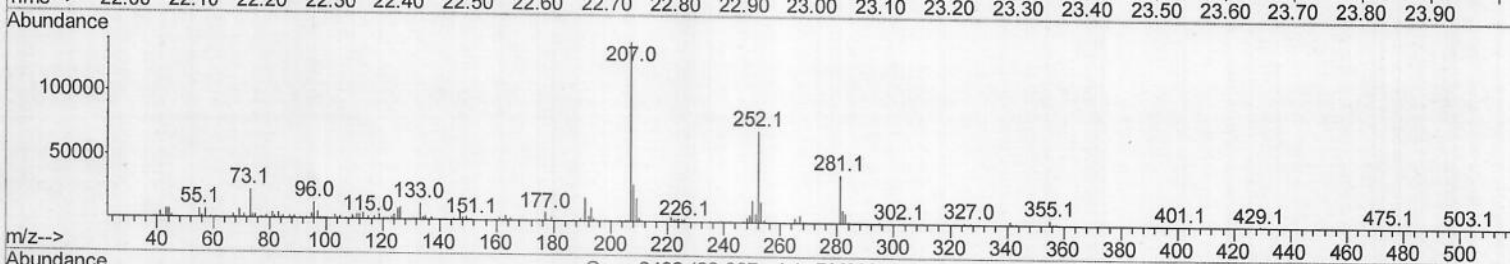
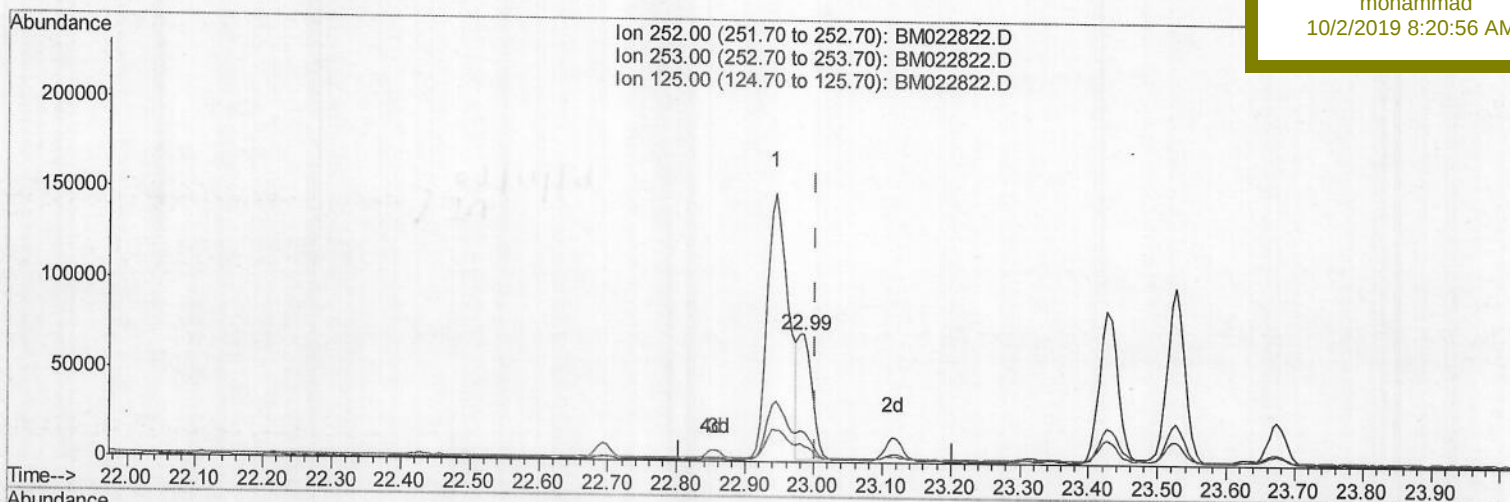
ESNY4

Manual Integrations

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mohammad

10/2/2019 8:20:56 AM



(89) Benzo(k)fluoranthene

22.986min (-0.018) 1.09ng/ul m

response 95133

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	24.01
125.00	9.90	12.53#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ESNY4

Quant Time: Oct 01 00:07:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
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 QLast Update : Fri Sep 27 18:03:27 2019
 Response via : Initial Calibration

Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	256898	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	1149704	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.45	164	744466	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.18	188	1563740	20.00	ng/ul	0.00
78) Chrysene-d12	21.36	240	1542798	20.00	ng/ul	0.00
86) Perylene-d12	23.63	264	1379071	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	15431	2.60	ng/uL	0.00
5) Phenol-d5	6.97	99	295979	13.04	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.14	67	189001	15.21	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	263892	14.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.51	113	246894	13.28	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	130187	14.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.68	143	137417	14.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	273067	14.15	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.73	131	348443	14.87	ng/ul	0.00
44) Dimethylphthalate-d6	13.85	166	927812	16.06	ng/ul	0.00
47) Acenaphthylene-d8	14.13	160	1057239	15.85	ng/ul	0.00
52) 4-Nitrophenol-d4	14.63	143	90227	7.90	ng/ul	-0.01
58) Fluorene-d10	15.43	176	816255	16.58	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.55	200	76357	7.50	ng/ul	0.00
71) Anthracene-d10	17.28	188	1277864	16.73	ng/ul	0.00
79) Pyrene-d10	19.57	212	1539582	18.54	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.48	264	1271998	17.83	ng/ul	-0.02

Target Compounds

					Ovalue
6) Phenol	7.00	94	31324	1.316 ng/ul	97
45) Dimethylphthalate	13.90	163	285160	4.804 ng/ul	100
77) Fluoranthene	19.24	202	355248	3.384 ng/ul	99
80) Pyrene	19.60	202	312180	2.854 ng/ul	97
83) Benzo(a)anthracene	21.34	228	137582	1.242 ng/ul	99
85) Chrysene	21.40	228	210318	2.059 ng/ul	99
88) Benzo(b)fluoranthene	22.94	252	328140	3.586 ng/ul	97
89) Benzo(k)fluoranthene	22.99	252	95133m	1.087 ng/ul	
91) Benzo(a)pyrene	23.53	252	177398	2.066 ng/ul	96
92) Indeno(1,2,3-cd)pyrene	25.92	276	154235	1.691 ng/ul	96
94) Benzo(a,h,i)perylene	26.62	276	157565	2.142 ng/ul	98

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