

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062319\

Data File : BM021112.D

Acq On : 22 Jun 2019 23:30

Operator : HP/JU

Sample : K3335-08

Misc :

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Jun 25 04:44:28 2019

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

Quant Title : SVOA CALIBRATION

QLast Update : Thu Jun 20 03:45:02 2019

Response via : Initial Calibration

Instrument :

BNA_M

Client Sample Id :

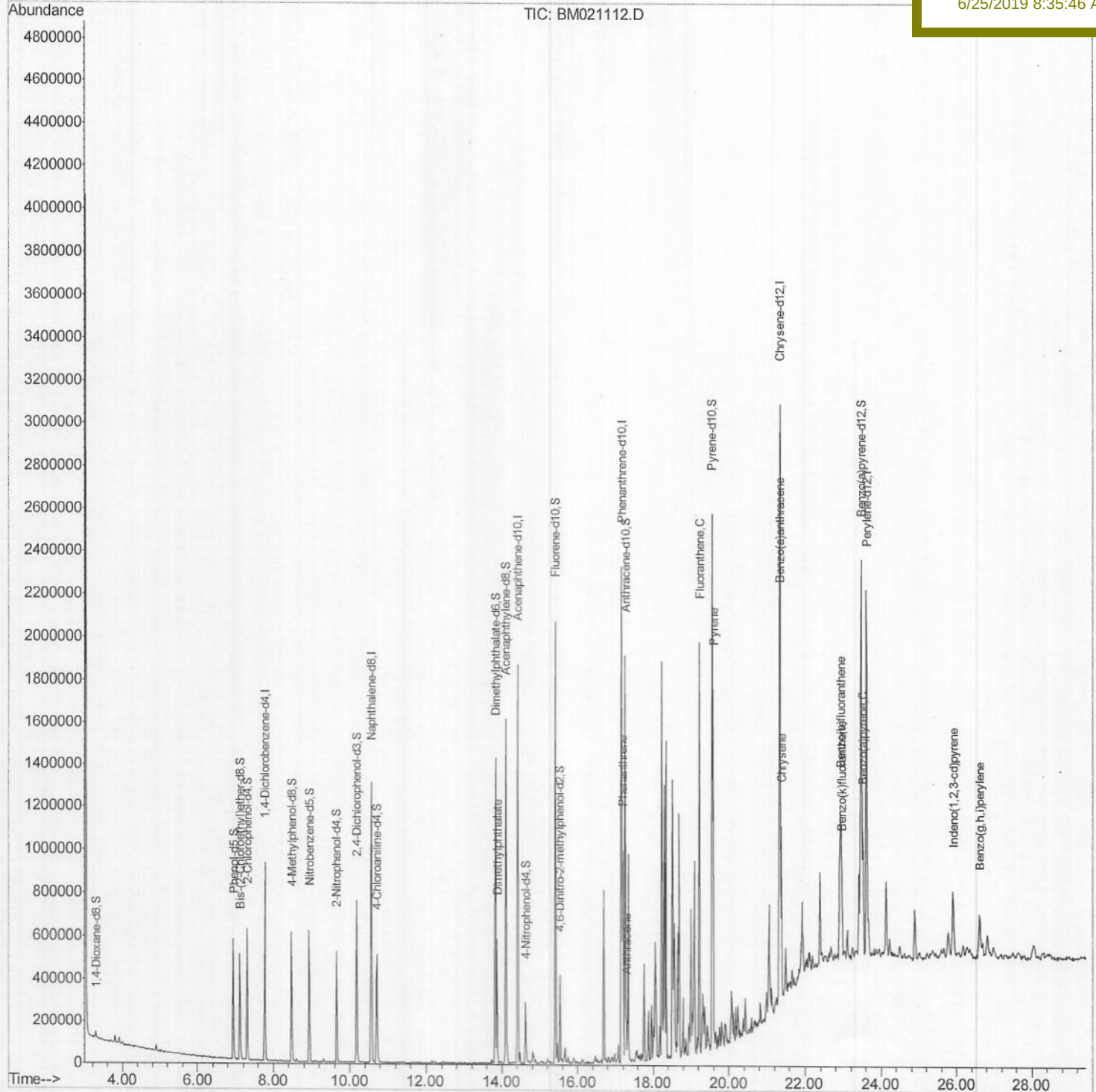
A4235

Manual Integrations

APPROVED

mohammad

6/25/2019 8:35:46 AM



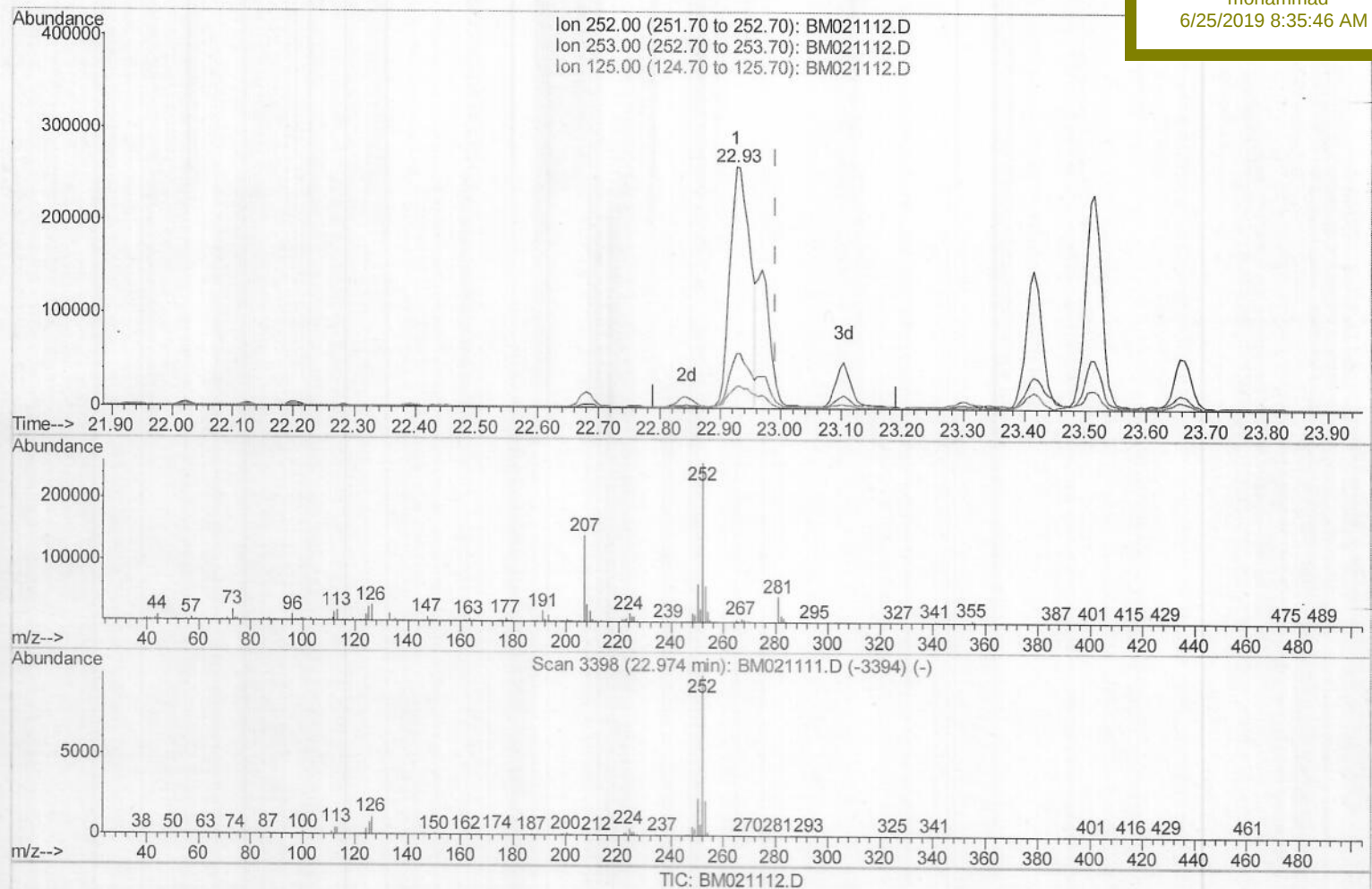
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\
Data File : BM021112.D
Acq On : 22 Jun 2019 23:30
Operator : HP/JU
Sample : K3335-08
Misc :
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Jun 24 00:52:53 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 20 03:45:02 2019
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
A4235

Manual Integrations
APPROVED

mohammad
6/25/2019 8:35:46 AM



(88) Benzo(k)fluoranthene

22.927min (-0.064) 7.51ng/ul

response 592985

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	22.73
125.00	8.20	9.18
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062319\

Data File : BM021112.D

Acq On : 22 Jun 2019 23:30

Operator : HP/JU

Sample : K3335-08

Misc :

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Jun 24 00:52:53 2019

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

Quant Title : SVOA CALIBRATION

QLast Update : Thu Jun 20 03:45:02 2019

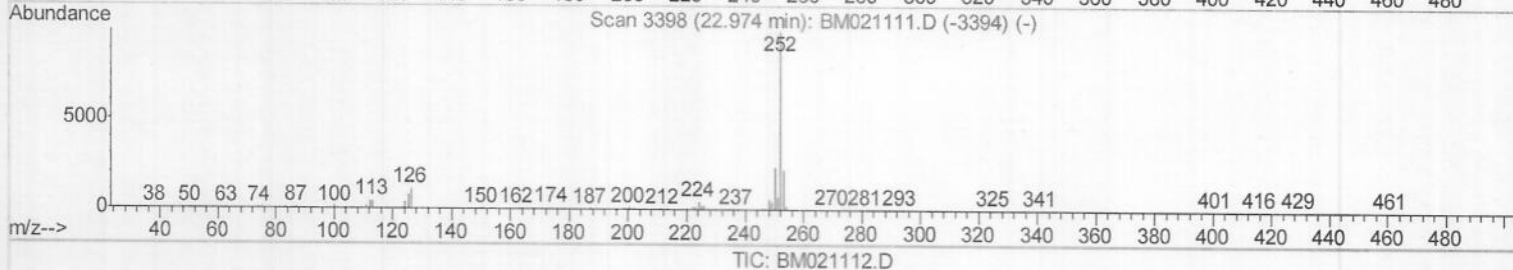
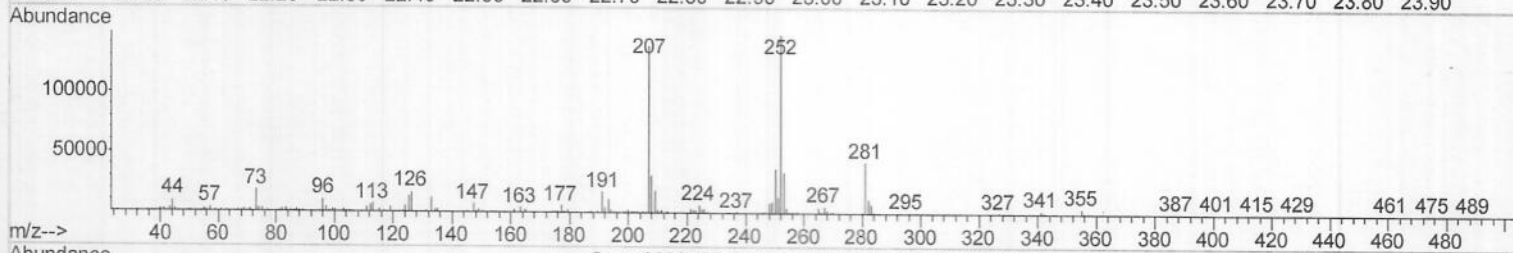
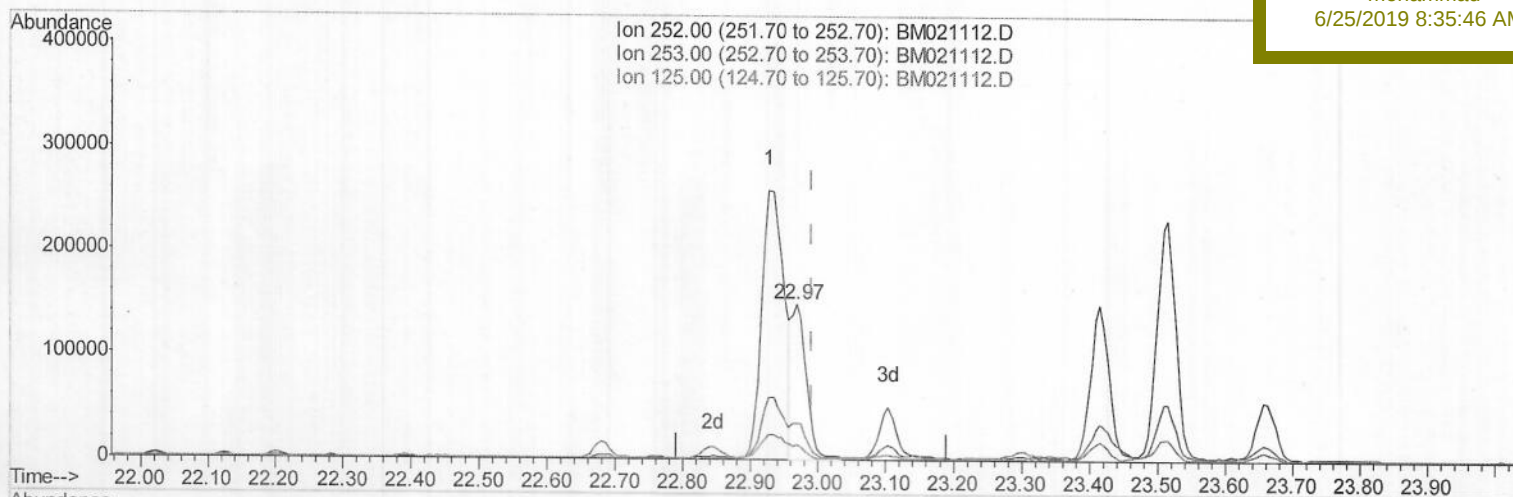
Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

A4235

Manual Integrations
APPROVEDmohammad
6/25/2019 8:35:46 AM

(88) Benzo(k)fluoranthene

22.968min (-0.023) 2.80ng/ul m > JU 06/25/19

response 221382

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	23.06
125.00	8.20	9.37
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062319\

Data File : BM021112.D

Acq On : 22 Jun 2019 23:30

Operator : HP/JU

Sample : K3335-08

Misc :

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Jun 25 04:44:28 2019

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

Quant Title : SVOA CALIBRATION

QLast Update : Thu Jun 20 03:45:02 2019

Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

A4235

Manual Integrations
APPROVEDmohammad
6/25/2019 8:35:46 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	255256	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.56	136	1088369	20.00	ng/ul	-0.02
35) Acenaphthene-d10	14.42	164	632799	20.00	ng/ul	-0.01
61) Phenanthrene-d10	17.16	188	1335876	20.00	ng/ul	-0.01
77) Chrysene-d12	21.34	240	1127730	20.00	ng/ul	-0.02
85) Perylene-d12	23.61	264	1296070	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	15868	2.95	ng/uL	0.00
5) Phenol-d5	6.94	99	316381	14.24	ng/ul	-0.01
7) Bis-(2-Chloroethyl) ether-d	7.12	67	223861	16.91	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.30	132	275630	15.27	ng/ul	-0.02
13) 4-Methylphenol-d8	8.48	113	223835	12.07	ng/ul	-0.02
19) Nitrobenzene-d5	8.93	128	151004	17.09	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.65	143	164976	16.82	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.18	165	298511	14.96	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.71	131	295997	14.23	ng/ul	-0.01
43) Dimethylphthalate-d6	13.83	166	987508	17.27	ng/ul	-0.01
46) Acenaphthylene-d8	14.10	160	1138690	16.27	ng/ul	-0.02
51) 4-Nitrophenol-d4	14.62	143	82578	8.88	ng/ul	-0.01
57) Fluorene-d10	15.41	176	824115	16.57	ng/ul	-0.01
62) 4,6-Dinitro-2-methylphenol	15.53	200	86628	10.55	ng/ul	-0.01
70) Anthracene-d10	17.26	188	1037580	14.95	ng/ul	-0.02
78) Pyrene-d10	19.55	212	1138855	16.83	ng/ul	-0.02
89) Benzo(a)pyrene-d12	23.47	264	1248249	16.32	ng/ul	-0.02

Target Compounds

					Ovalue
44) Dimethylphthalate	13.87	163	376465	7.233 ng/ul	99
69) Phenanthrene	17.20	178	555349	7.517 ng/ul	99
71) Anthracene	17.29	178	126030	1.674 ng/ul	98
76) Fluoranthene	19.22	202	1161029	14.544 ng/ul	99
79) Pyrene	19.58	202	1021178	12.809 ng/ul	99
82) Benzo(a)anthracene	21.32	228	504364	6.560 ng/ul	98
84) Chrysene	21.37	228	449963	6.249 ng/ul	97
87) Benzo(b)fluoranthene	22.93	252	592985	7.223 ng/ul	98
88) Benzo(k)fluoranthene	22.97	252	221382m	2.803 ng/ul	98
90) Benzo(a)pyrene	23.52	252	428234	5.542 ng/ul	98
91) Indeno(1,2,3-cd)pyrene	25.90	276	290233	3.190 ng/ul	96
93) Benzo(a,h,i)perylene	26.60	276	244006	3.256 ng/ul	98

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