

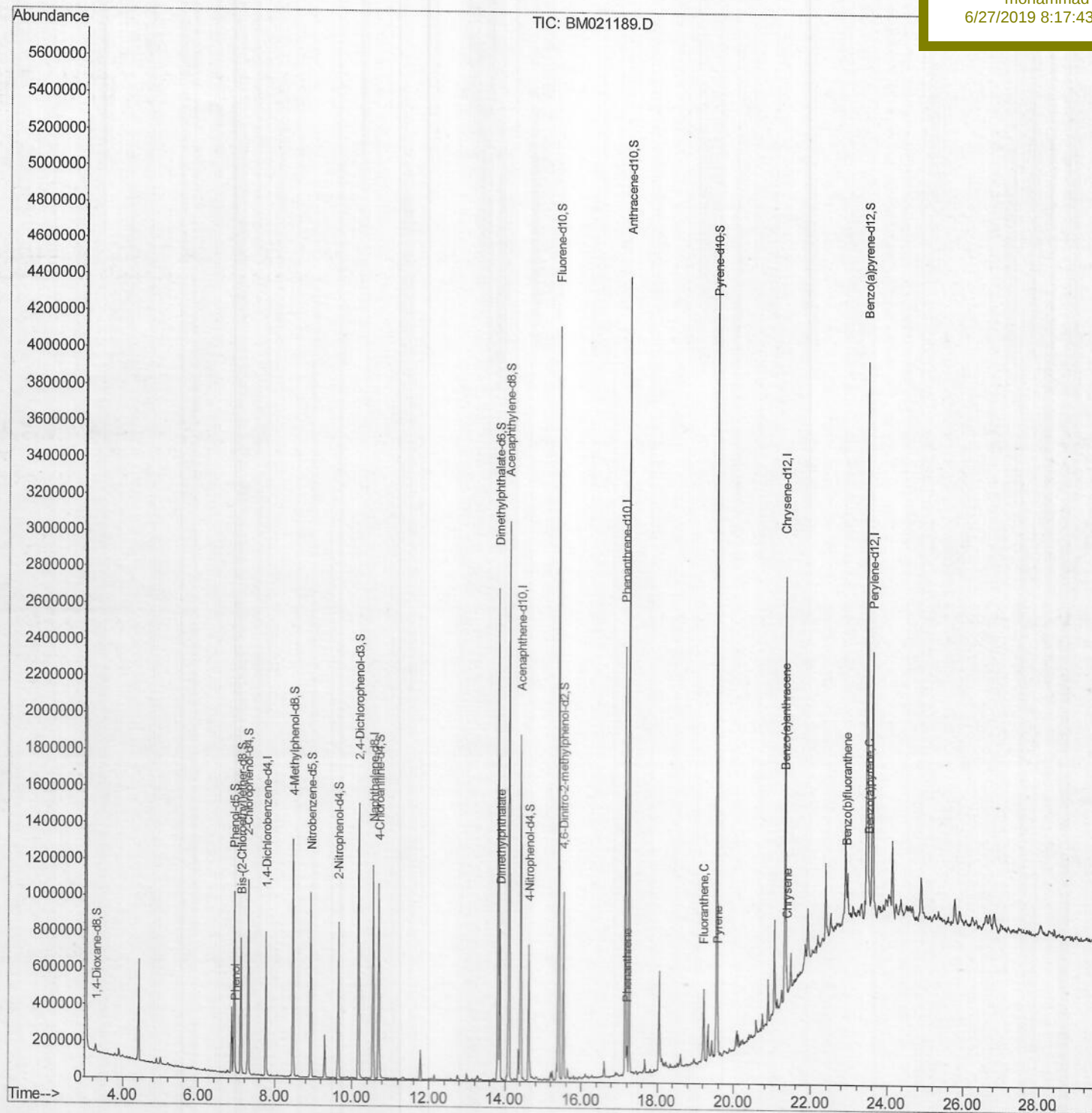
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062519\
Data File : BM021189.D
Acq On : 26 Jun 2019 13:27
Operator : HP/JU
Sample : K3501-05
Misc :
ALS Vial : 38 Sample Multiplier: 1

Quant Time: Jun 26 14:07:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 26 01:32:57 2019
Response via : Initial Calibration

Instrument :
BNA_M
Client Sample ID :
C5AC5

Manual Integrations
APPROVED

mohammad
6/27/2019 8:17:43 AM



Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062519\

Data File : BM021189.D

Acq On : 26 Jun 2019 13:27

Operator : HP/JU

Sample : K3501-05

Misc :

ALS Vial : 38 Sample Multiplier: 1

Quant Time: Jun 26 14:04:45 2019

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

Quant Title : SVOA CALIBRATION

QLast Update : Wed Jun 26 01:32:57 2019

Response via : Initial Calibration

Instrument :

BNA M

ClientSampleId :

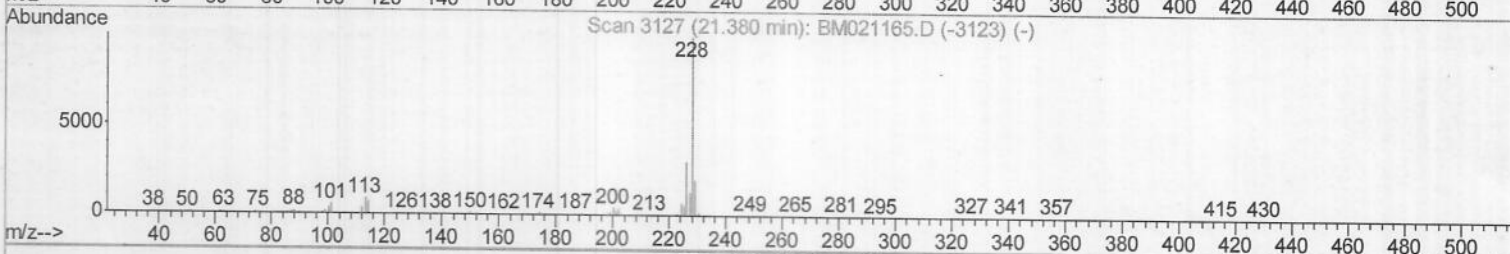
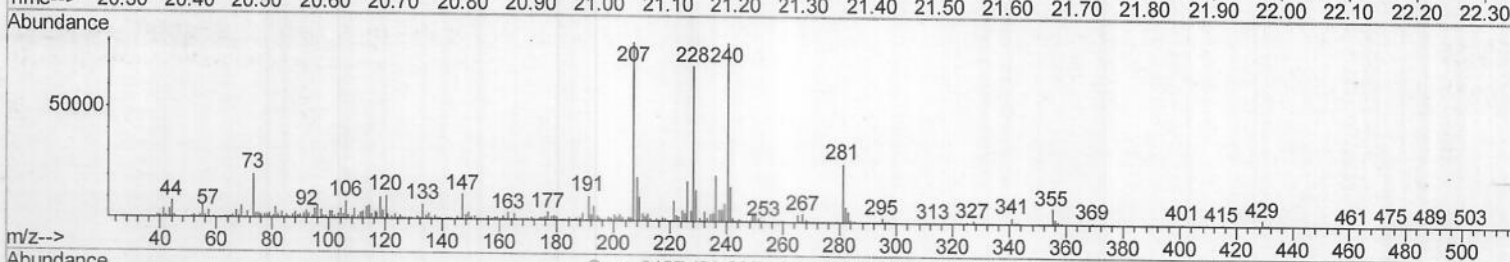
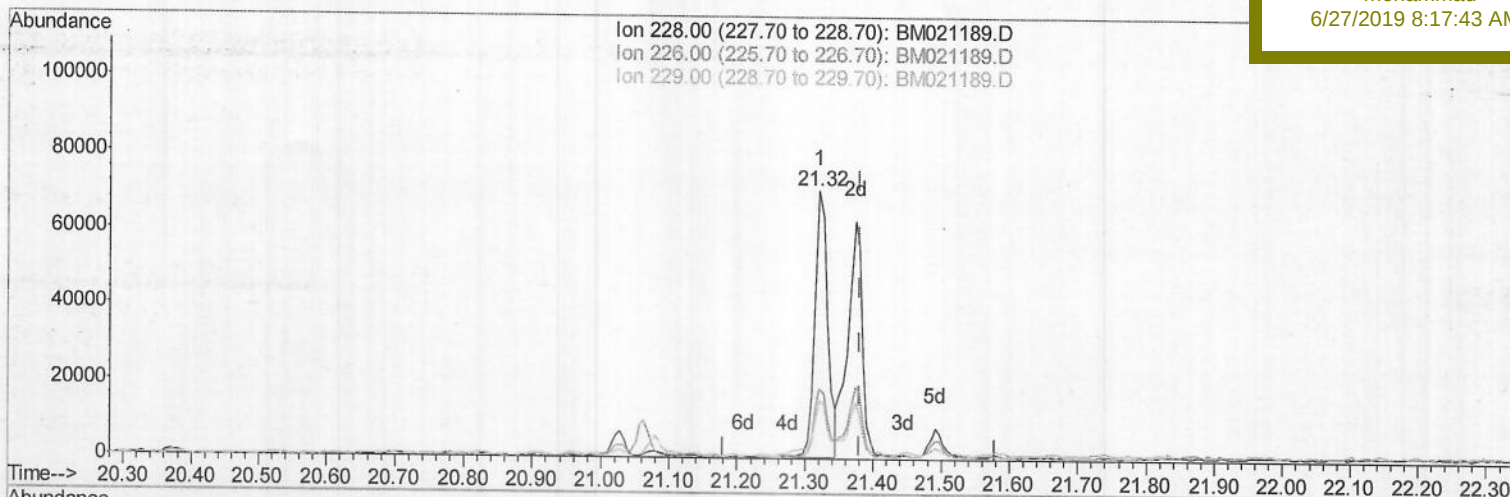
C5AC5

Manual Integrations

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mohammad

6/27/2019 8:17:43 AM



TIC: BM021189.D

(84) Chrysene

21.321min (-0.059) 1.42ng/ul

response 92611

Ion	Exp%	Act%
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228.00	100	100
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226.00	29.90	26.01
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229.00	19.30	21.16
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0.00	0.00	0.00
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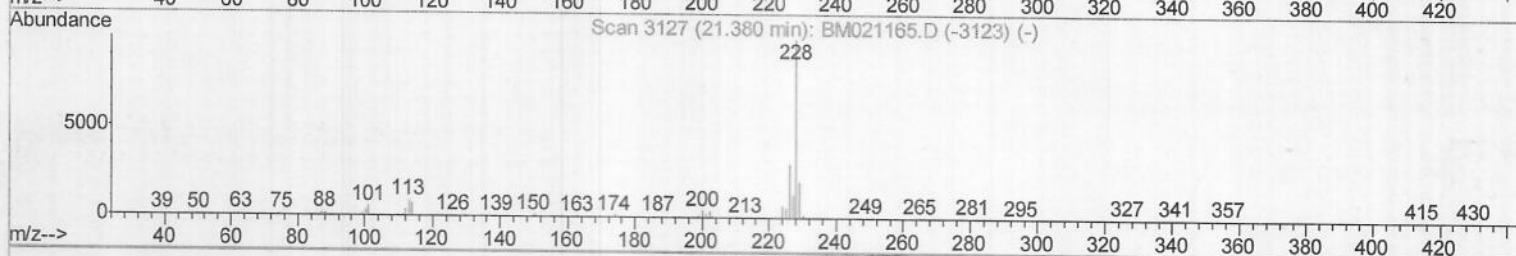
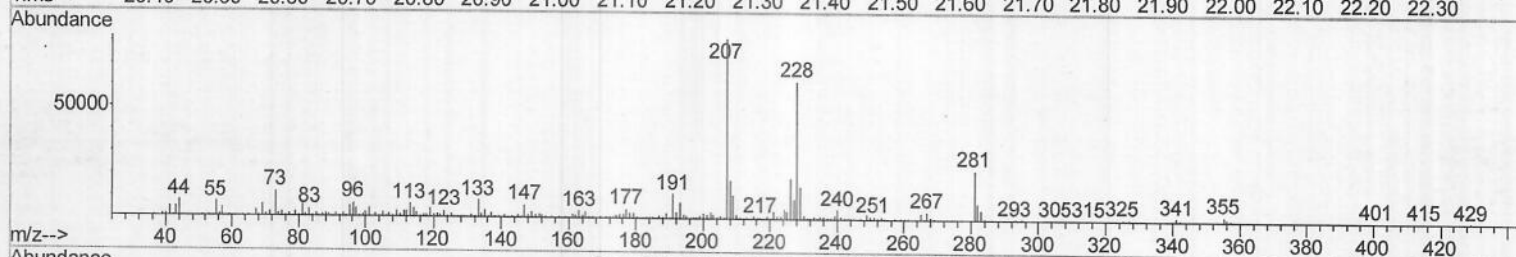
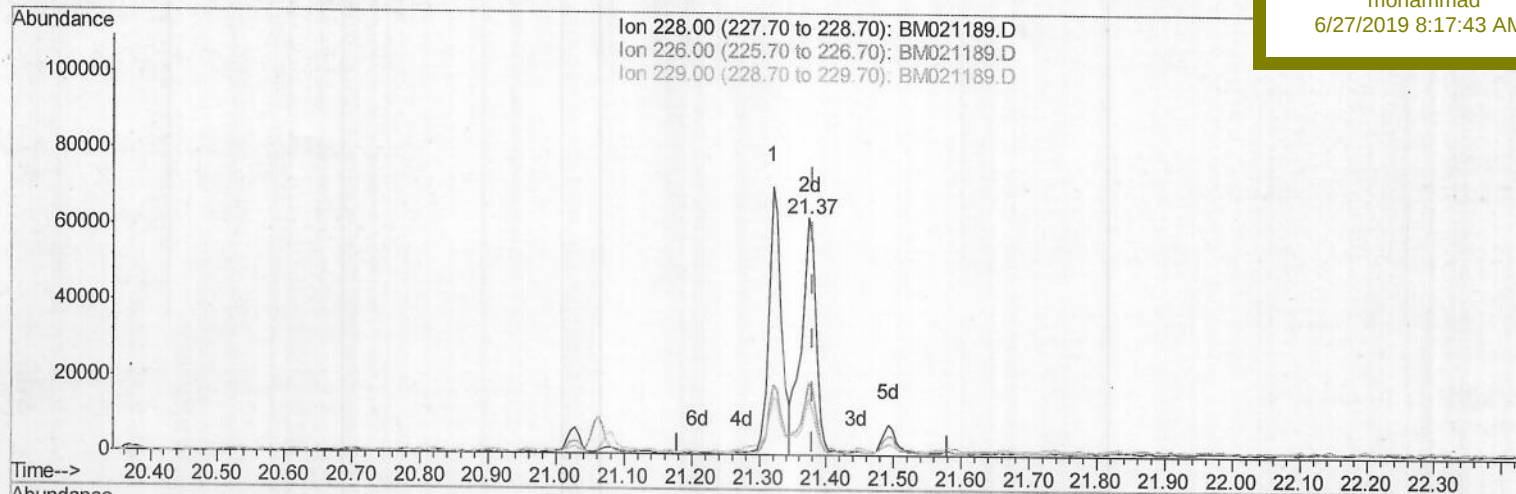
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Manual Integrations
APPROVED

mohammad
6/27/2019 8:17:43 AM



TIC: BM021189.D

(84) Chrysene

21.374min (-0.006) 1.46ng/ul m

JU 06/27/19

response 95640

Ion	Exp%	Act%
228.00	100	100
226.00	29.90	30.41
229.00	19.30	23.38#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062519\
 Data File : BM021189.D
 Acq On : 26 Jun 2019 13:27
 Operator : HP/JU
 Sample : K3501-05
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5AC5

Quant Time: Jun 26 14:07:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 26 01:32:57 2019
 Response via : Initial Calibration

Manual Integrations
 APPROVED

mohammad
 6/27/2019 8:17:43 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	226070	20.00	ng/ul	0.00
18) Naphthalene-d8	10.55	136	988554	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.40	164	620529	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	1407908	20.00	ng/ul	0.00
77) Chrysene-d12	21.34	240	1024855	20.00	ng/ul	0.00
85) Perylene-d12	23.61	264	1100939	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.28	96	22897	4.81	ng/uL	0.00
5) Phenol-d5	6.93	99	560896	28.50	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.11	67	329707	28.12	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	474276	29.66	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	489096	29.78	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	244721	30.49	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	278599	31.27	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.17	165	588500	32.48	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	595097	31.49	ng/ul	0.00
43) Dimethylphthalate-d6	13.82	166	1851895	33.03	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	2174887	31.69	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	198628	21.78	ng/ul	0.00
57) Fluorene-d10	15.40	176	1656728	33.97	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	222668	25.73	ng/ul	0.00
70) Anthracene-d10	17.26	188	2450495	33.49	ng/ul	0.00
78) Pyrene-d10	19.55	212	2318147	37.69	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.47	264	2103120	32.38	ng/ul	0.00
Target Compounds						
6) Phenol	6.96	94	27758	1.491	ng/ul	93
44) Dimethylphthalate	13.87	163	531197	10.407	ng/ul	99
69) Phenanthrene	17.20	178	91786	1.179	ng/ul	99
76) Fluoranthene	19.21	202	227859	2.708	ng/ul	98
79) Pyrene	19.58	202	181373	2.503	ng/ul	99
82) Benzo(a)anthracene	21.32	228	91727	1.313	ng/ul	98
84) Chrysene	21.37	228	95640m	1.462	ng/ul	91
87) Benzo(b)fluoranthene	22.93	252	118264	1.696	ng/ul#	91
90) Benzo(a)pyrene	23.51	252	78929	1.202	ng/ul#	90

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

JU 06/27/19