

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\

Data File : BN008359.D

Acq On : 23 Oct 2019 20:55

Operator : JU

Sample : K5223-17DL 2X

Misc :

ALS Vial : 52 Sample Multiplier: 1

Quant Time: Oct 24 02:03:02 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Oct 23 14:07:55 2019

Response via : Initial Calibration

Instrument :

BNA_N

ClientSampleId :

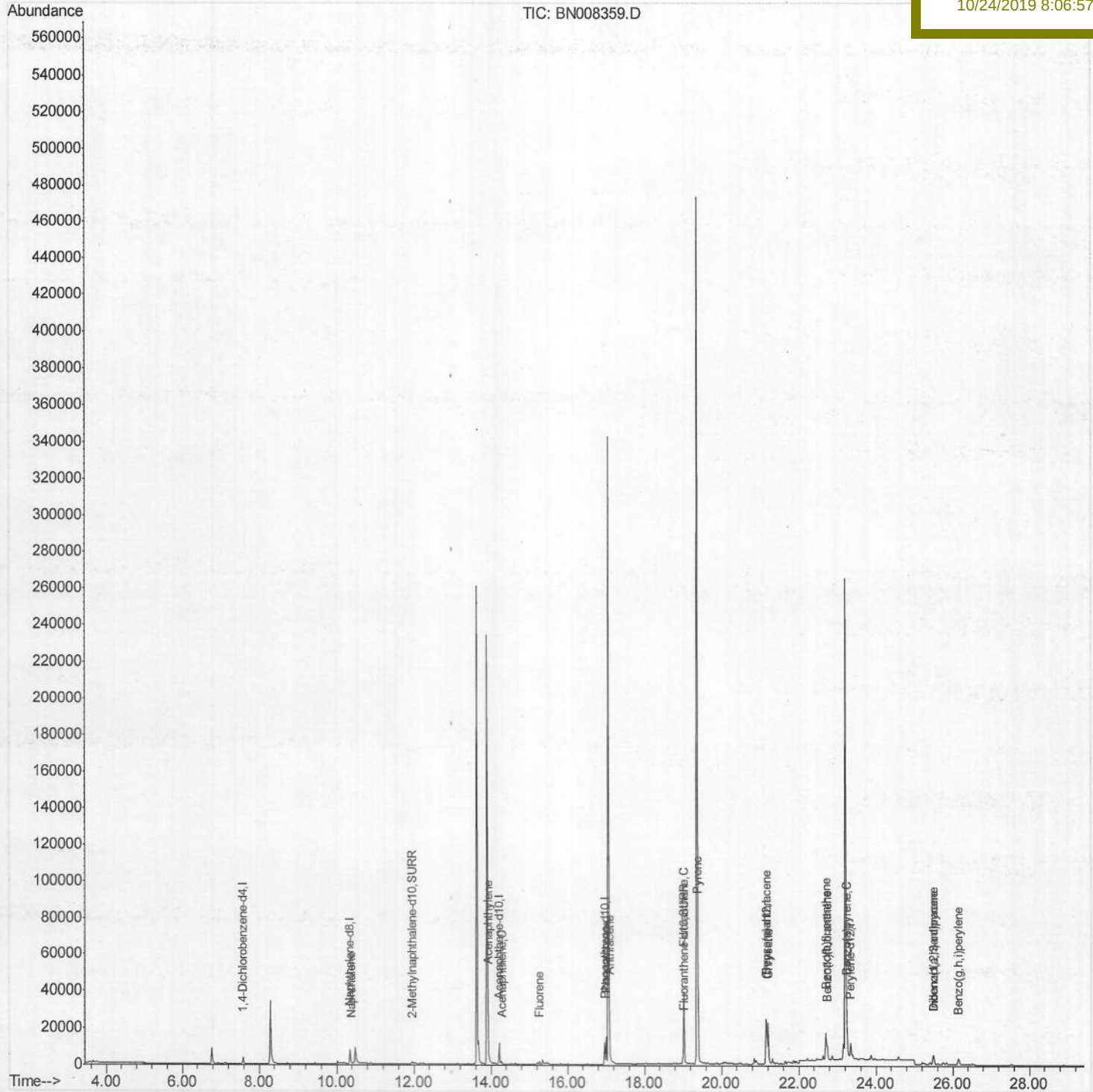
BEPA5DL

Manual Integrations

APPROVED

mohammad

10/24/2019 8:06:57 AM



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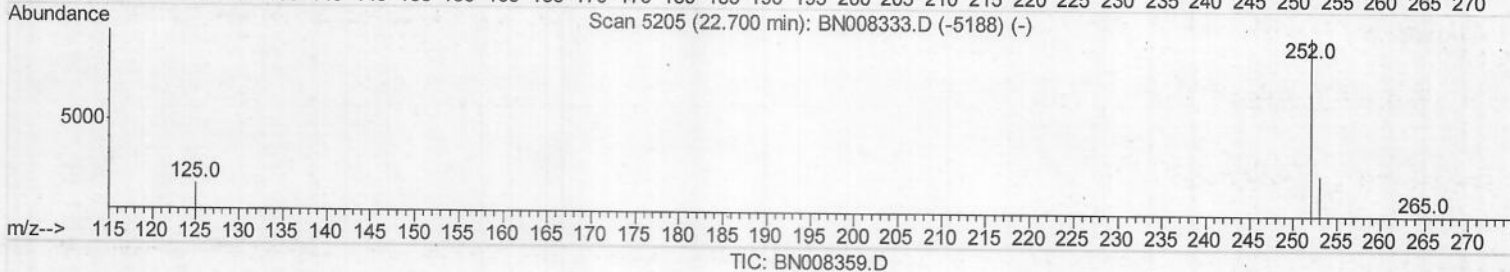
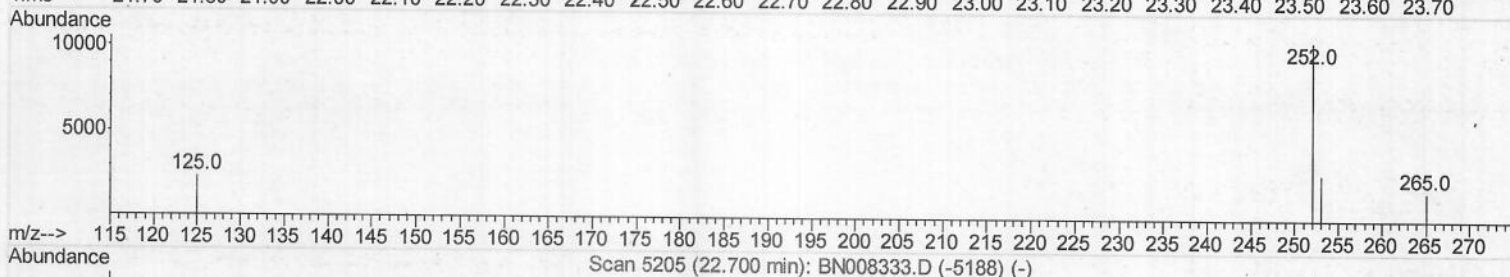
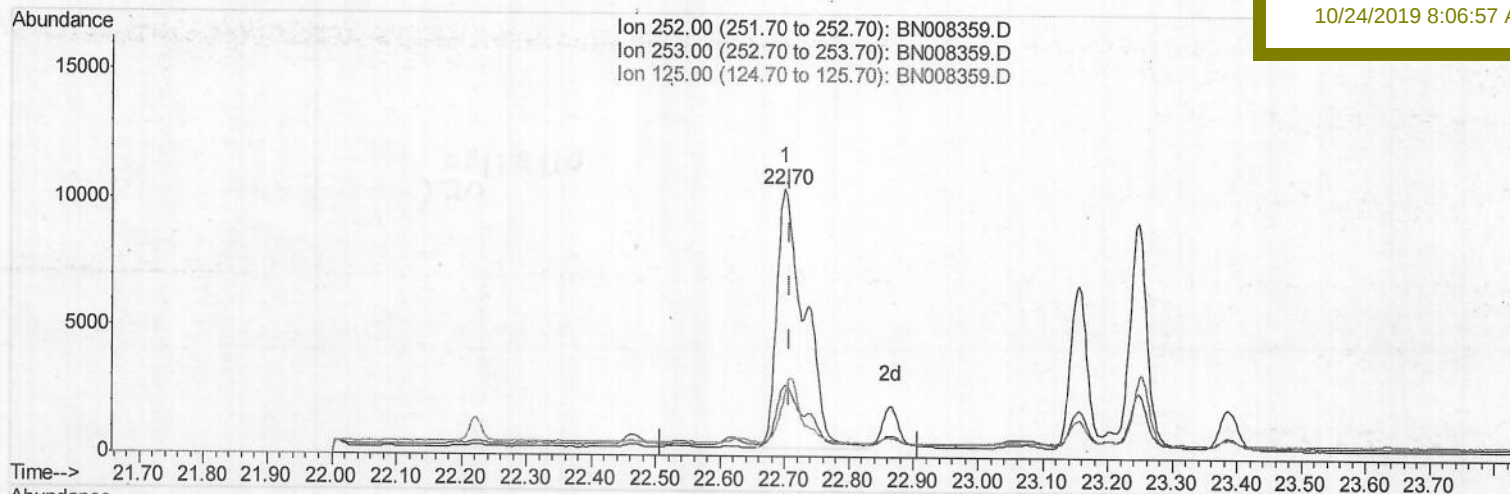
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(21) Benzo(b)fluoranthene

22.700min (-0.009) 0.94ng/ul

response 30764

Ion	Exp%	Act%
252.00	100	100
253.00	26.70	26.36
125.00	16.20	22.92
0.00	0.00	0.00

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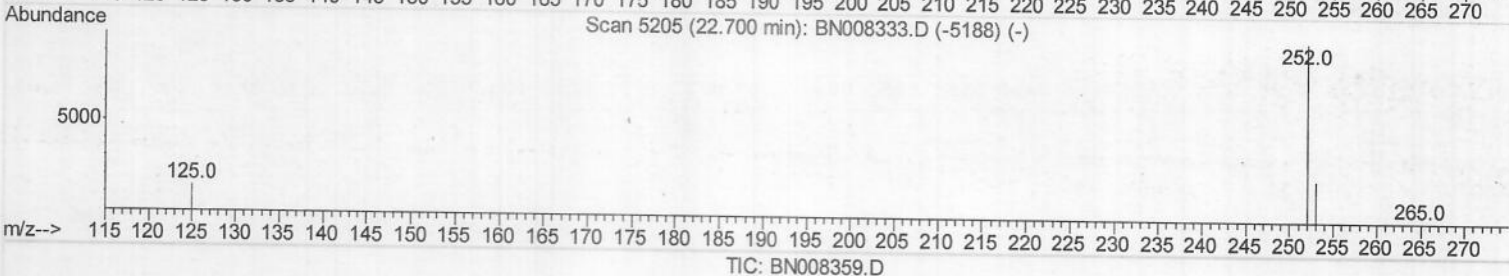
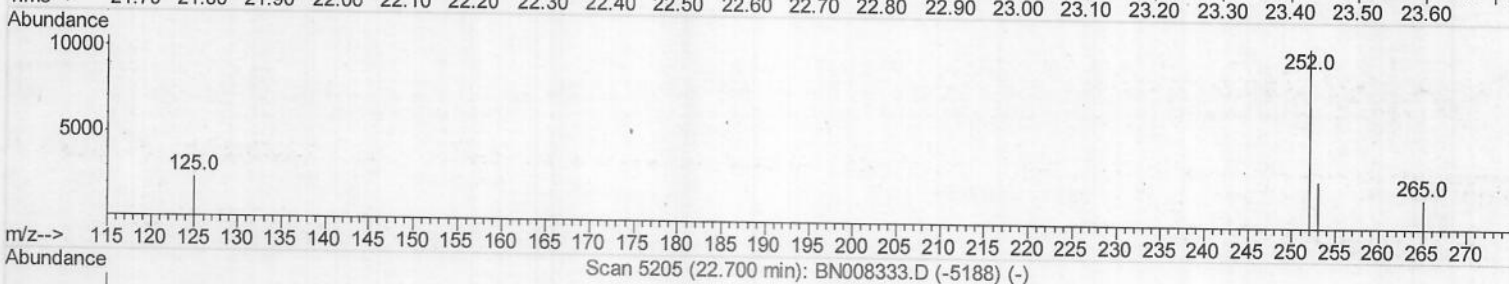
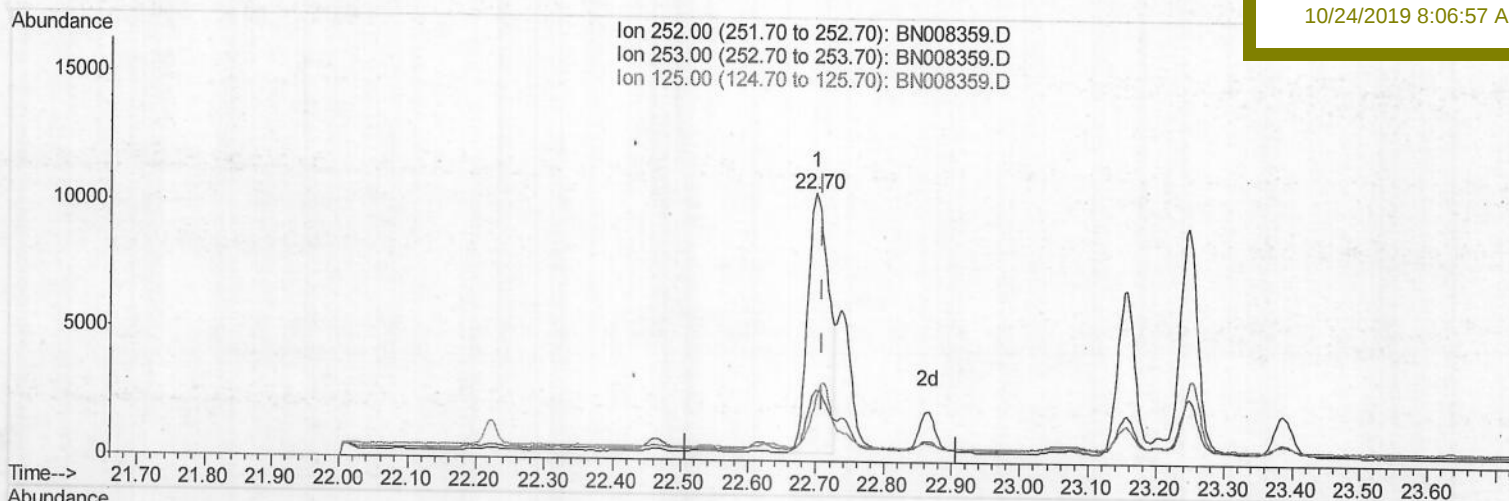
Response via : Initial Calibration

Instrument :

BNA_N

ClientSampleId :

BEPA5DL

Manual Integrations
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(21) Benzo(b)fluoranthene

22.700min (-0.009) 0.66ng/ul m

response 21588

Ion Exp% Act%

252.00 100 100

253.00 26.70 26.36

125.00 16.20 22.92

0.00 0.00 0.00

JU
10/24/19

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Misc :

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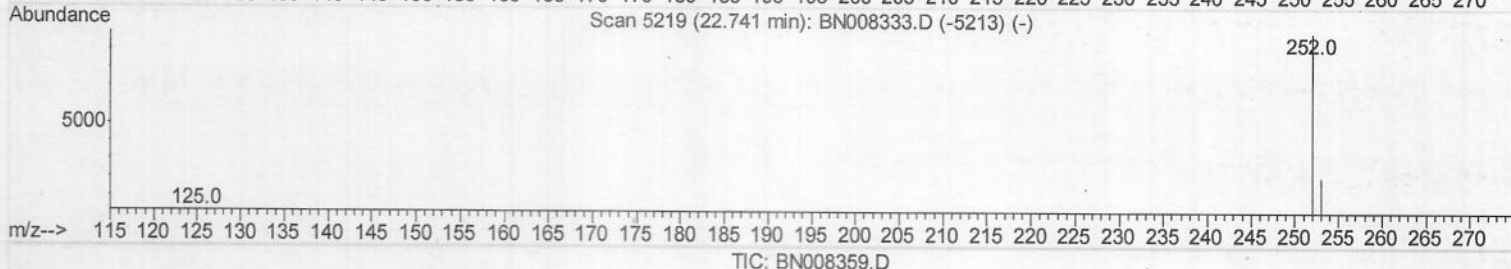
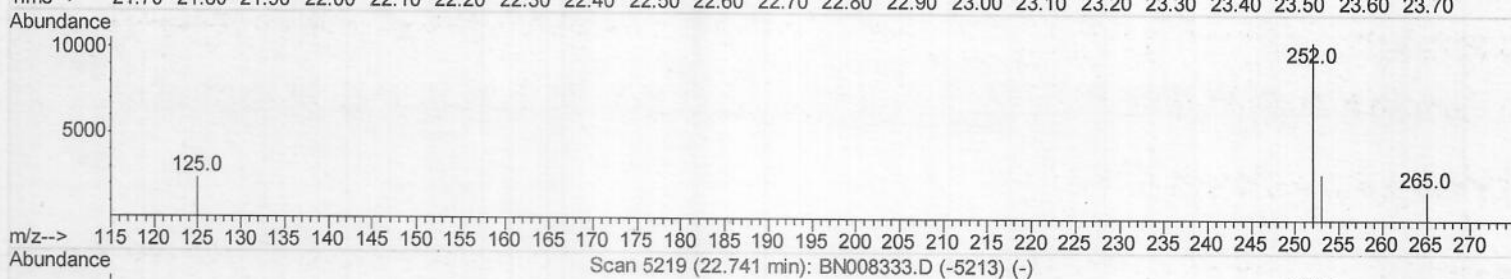
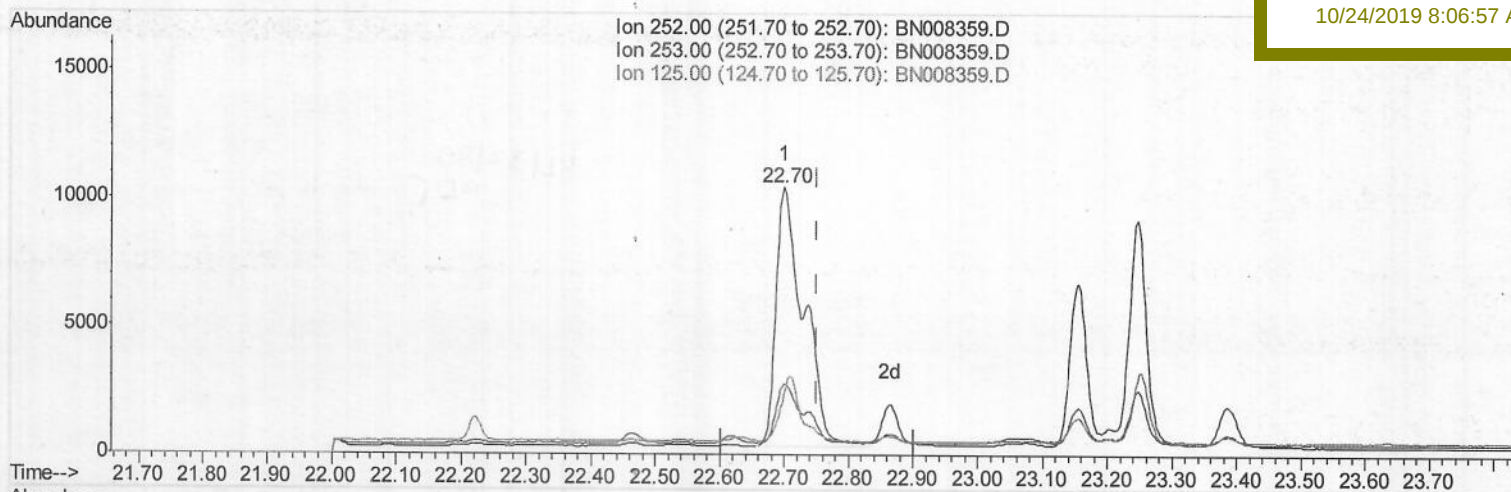
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Manual Integrations

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(22) Benzo(k)fluoranthene

22.700min (-0.052) 0.83ng/ul

response 30764

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	26.36
125.00	15.40	22.92#
0.00	0.00	0.00

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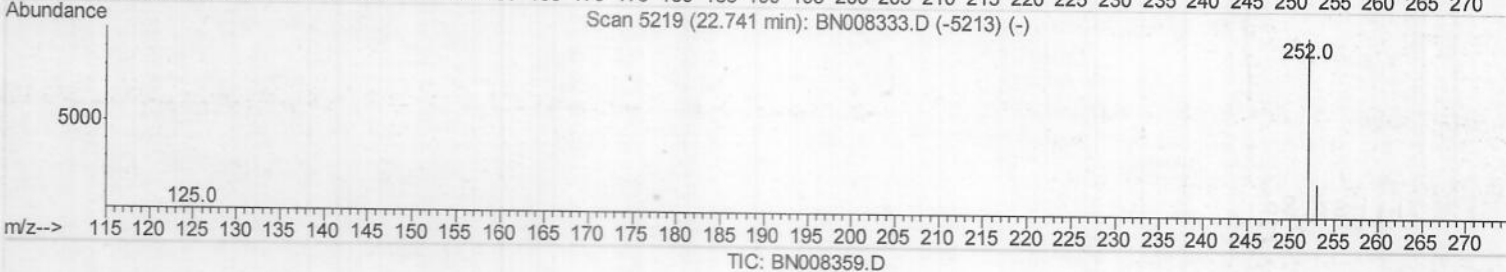
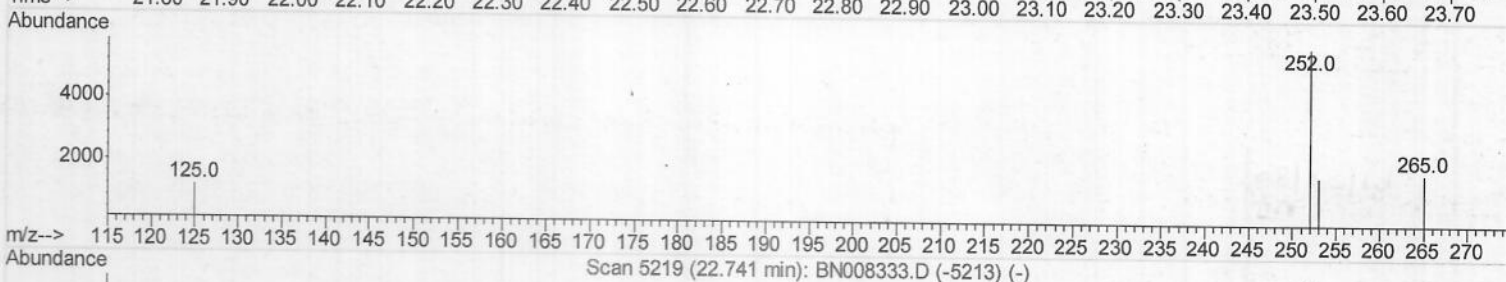
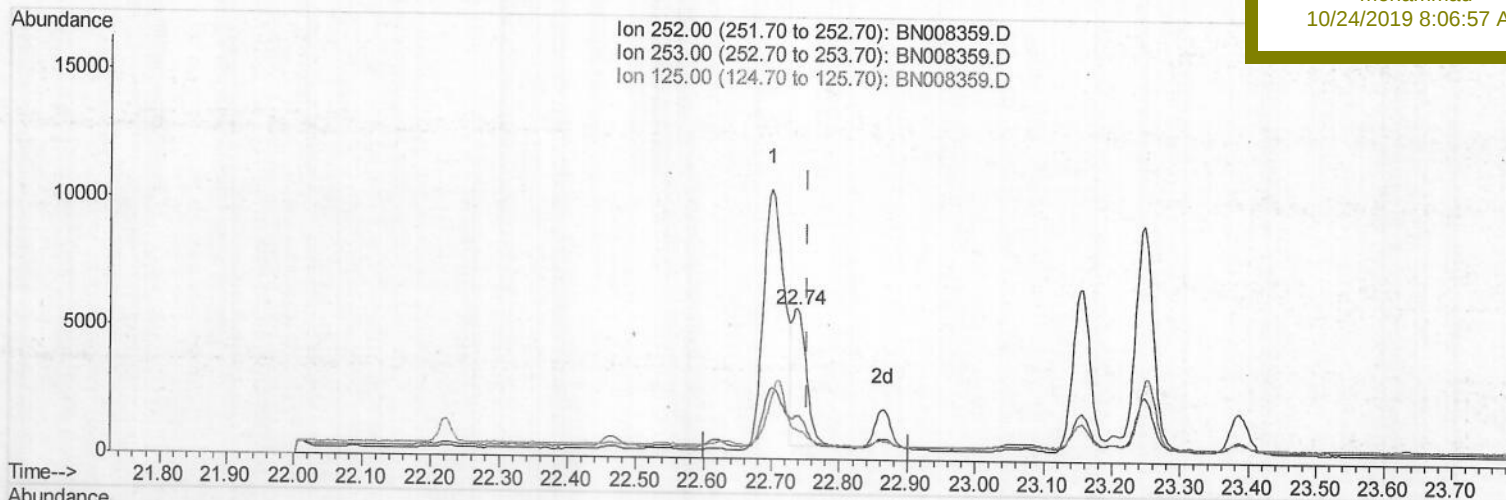
BEPA5DL

Manual Integrations

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10/24/2019 8:06:57 AM



(22) Benzo(k)fluoranthene

22.738min (-0.014) 0.22ng/ul m

response 8172

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	28.86
125.00	15.40	19.85#
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	2080	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.34	136	10028	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	6359	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	13781	0.40	ng/ul	-0.01
16) Chrysene-d12	21.17	240	11905	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	10063	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	1779	0.11	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	5239	0.11	ng/ul	-0.01
Target Compounds						
					Ovalue	
3) Naphthalene	10.39	128	640	0.022	ng/ul#	73
7) Acenaphthylene	13.93	152	936	0.031	ng/ul#	83
8) Acenaphthene	14.27	153	744	0.032	ng/ul	94
9) Fluorene	15.27	166	799	0.030	ng/ul#	95
12) Phenanthrene	17.00	178	15494	0.384	ng/ul	100
13) Anthracene	17.09	178	3001	0.085	ng/ul#	94
15) Fluoranthene	19.03	202	35733	0.671	ng/ul	96
17) Pyrene	19.40	202	41901	0.884	ng/ul	99
18) Benzo(a)anthracene	21.15	228	20591	0.476	ng/ul	97
19) Chrysene	21.21	228	23611	0.445	ng/ul	98
21) Benzo(b)fluoranthene	22.70	252	21588m	0.659	ng/ul	} JU 10/24/19
22) Benzo(k)fluoranthene	22.74	252	8172m	0.220	ng/ul	
23) Benzo(a)pyrene	23.25	252	16156	0.469	ng/ul#	
24) Indeno(1,2,3-cd)pyrene	25.49	276	8638	0.213	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.49	278	2100	0.065	ng/ul#	71
26) Benzo(a,h,i)perylene	26.14	276	6957	0.187	ng/ul#	94

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