

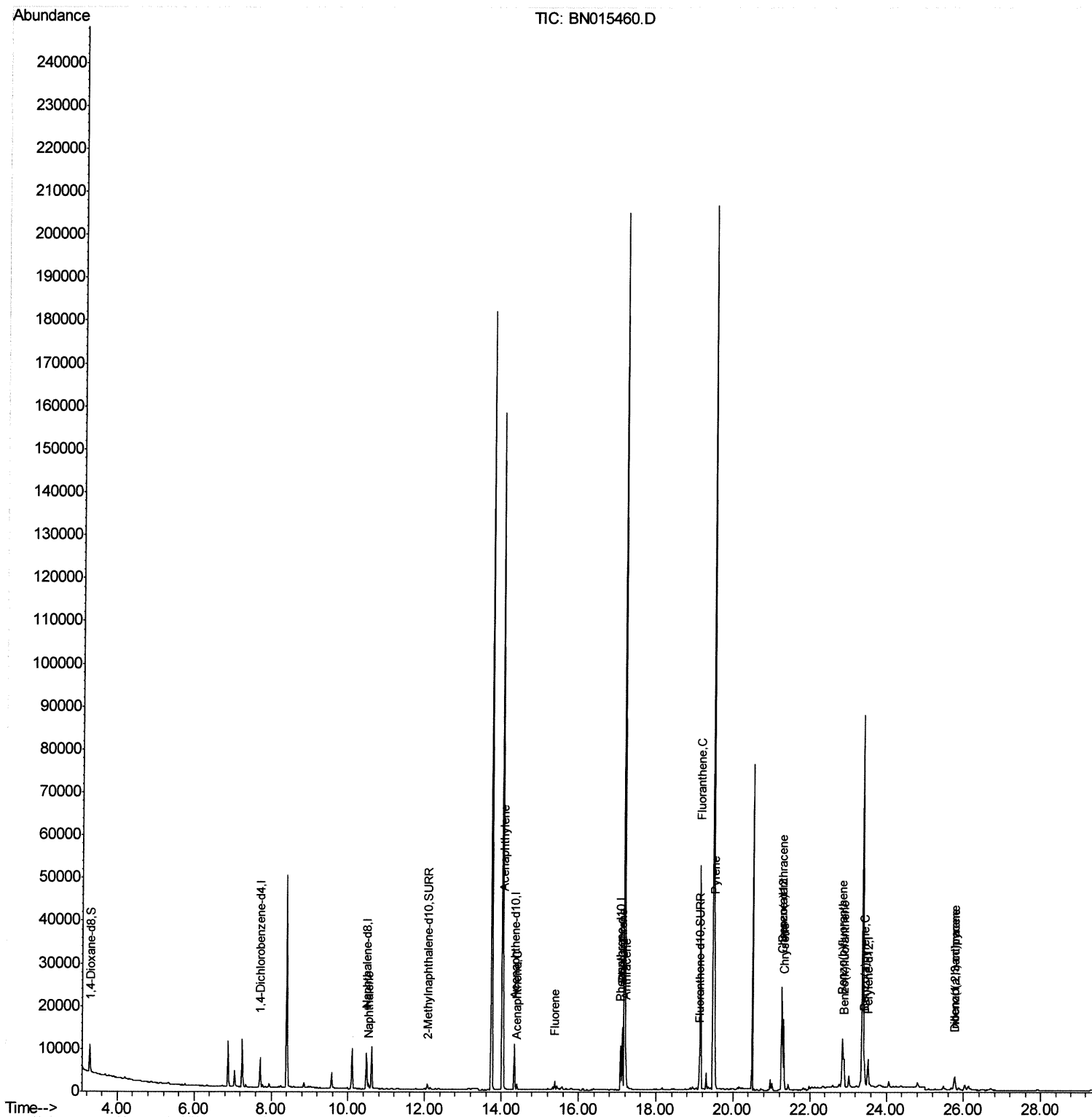
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN070821\
Data File : BN015460.D
Acq On : 09 Jul 2021 23:39
Operator : CG/JU
Sample : M2854-15DL 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDT9DL

Quant Time: Jul 10 02:44:57 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 08 11:17:31 2021
Response via : Initial Calibration

Manual Integrations
APPROVED

mohammad
7/12/2021 9:33:31 AM



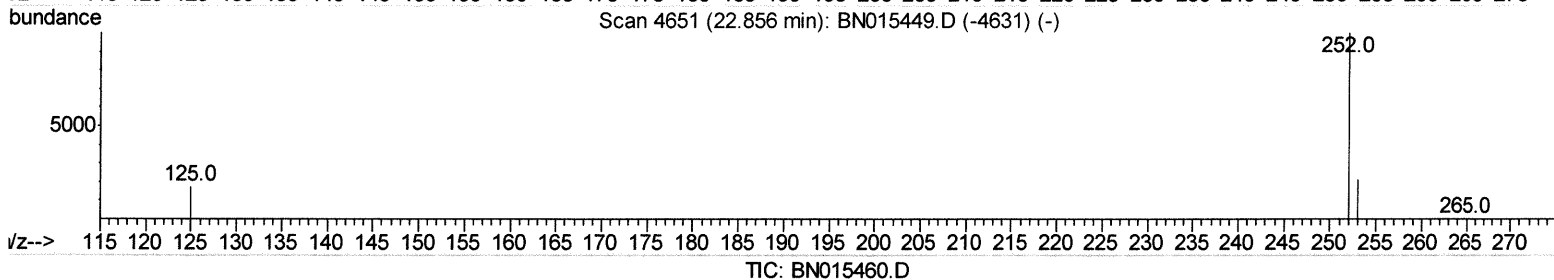
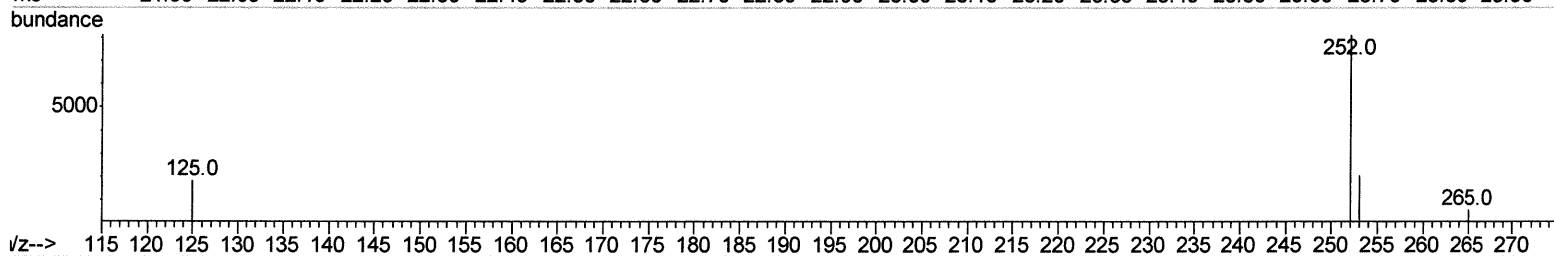
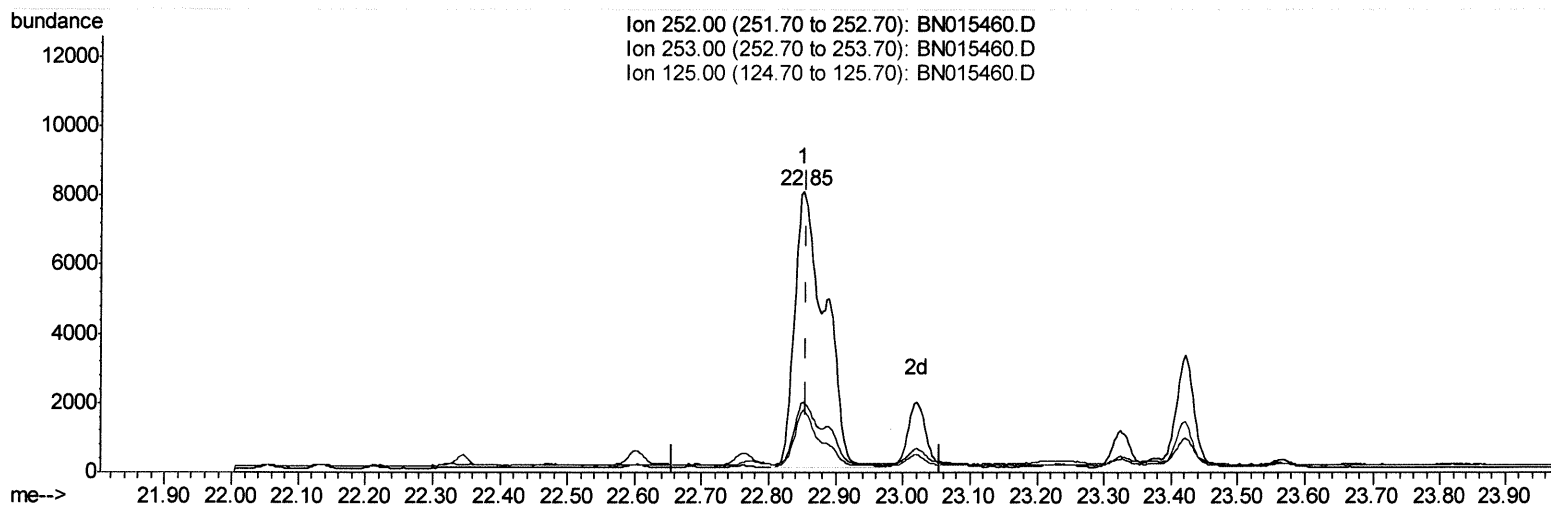
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(24) Benzo(b)fluoranthene
 22.851min (-0.005) 0.65ng/ul
 response 25181

Ion	Exp%	Act%
252.00	100	100
253.00	22.80	24.89
125.00	14.20	22.18
0.00	0.00	0.00

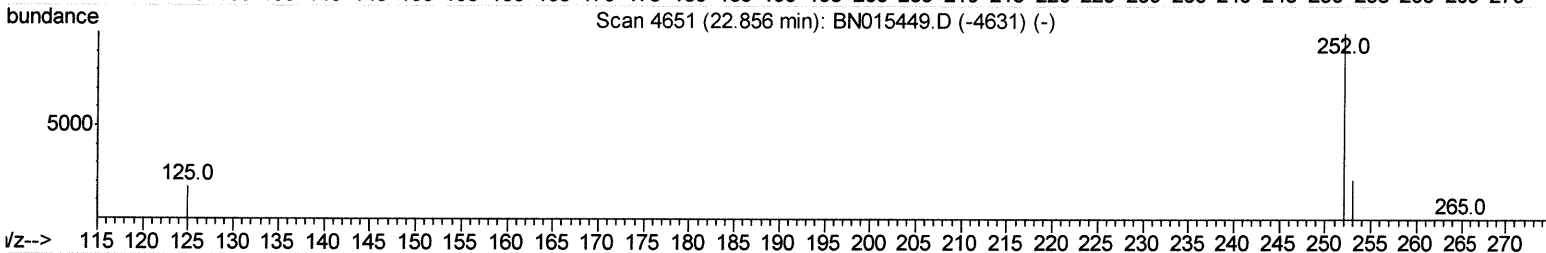
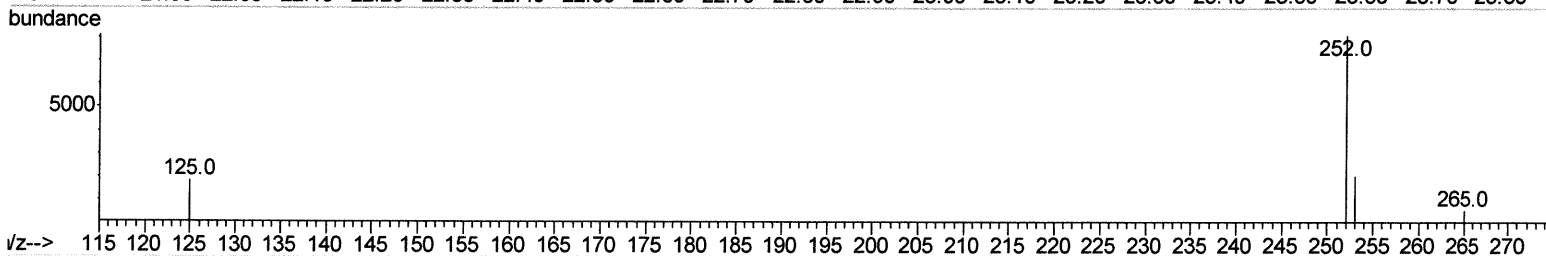
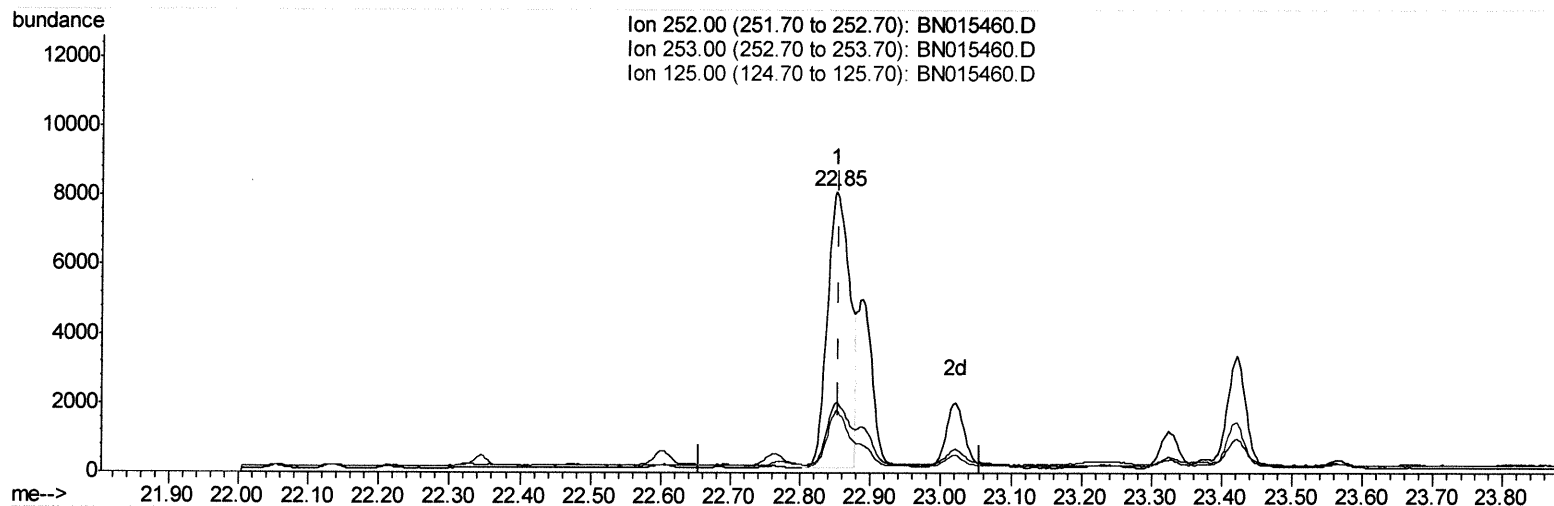
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(24) Benzo(b)fluoranthene

22.851min (-0.005) 0.45ng/ul m

response 17582

Ion	Exp%	Act%
252.00	100	100
253.00	22.80	24.89
125.00	14.20	22.18
0.00	0.00	0.00

547/14/21

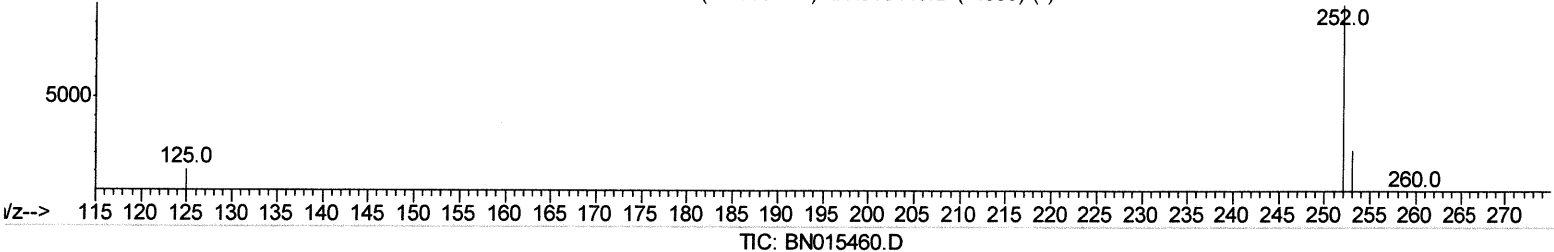
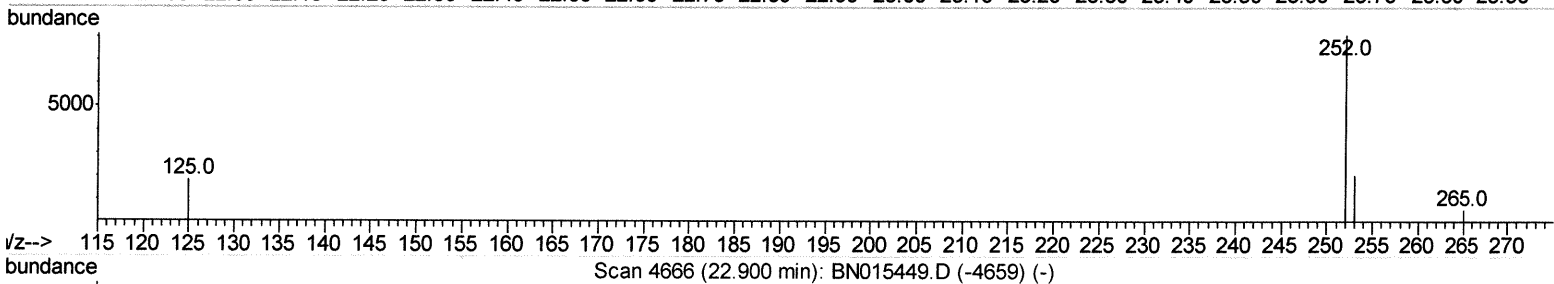
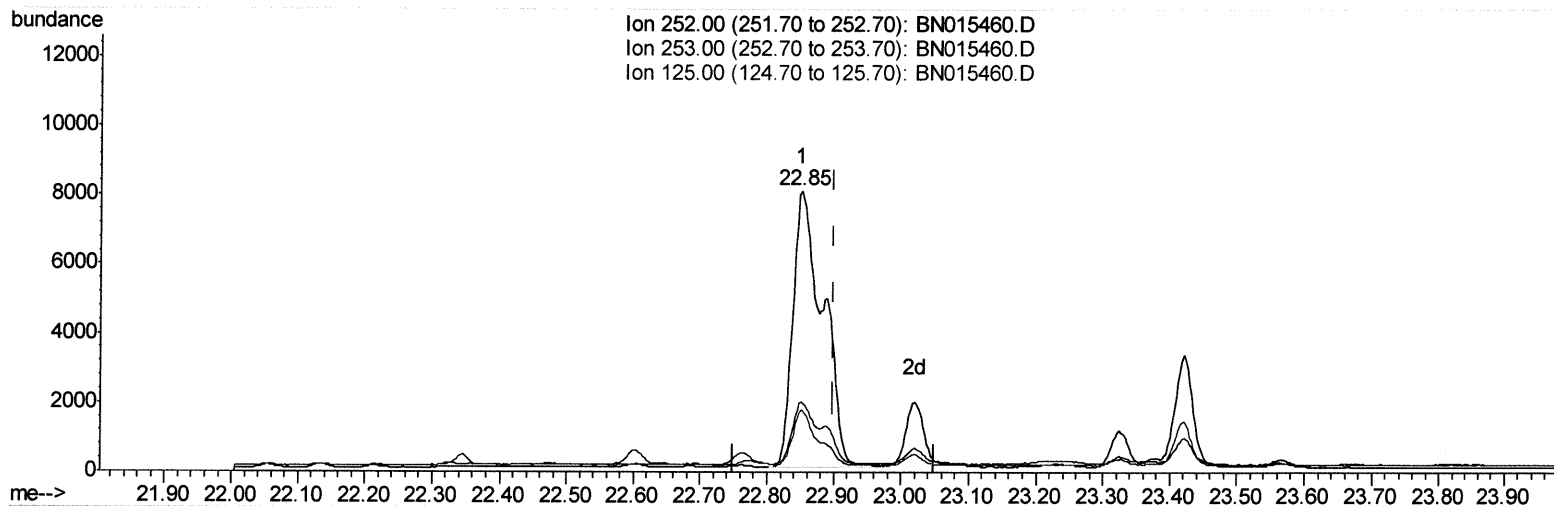
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Instrument :
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(25) Benzo(k)fluoranthene
 22.851min (-0.049) 0.57ng/ul
 response 25181

Ion	Exp%	Act%
252.00	100	100
253.00	22.60	24.89
125.00	14.40	22.18#
0.00	0.00	0.00

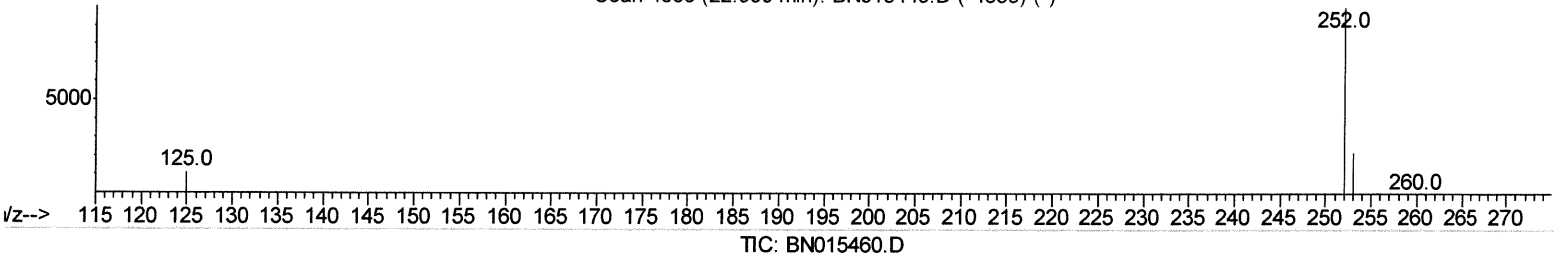
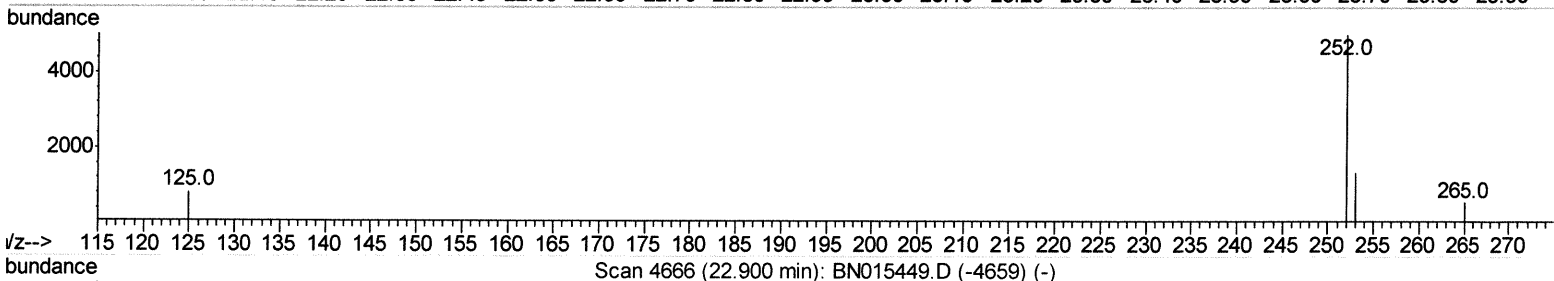
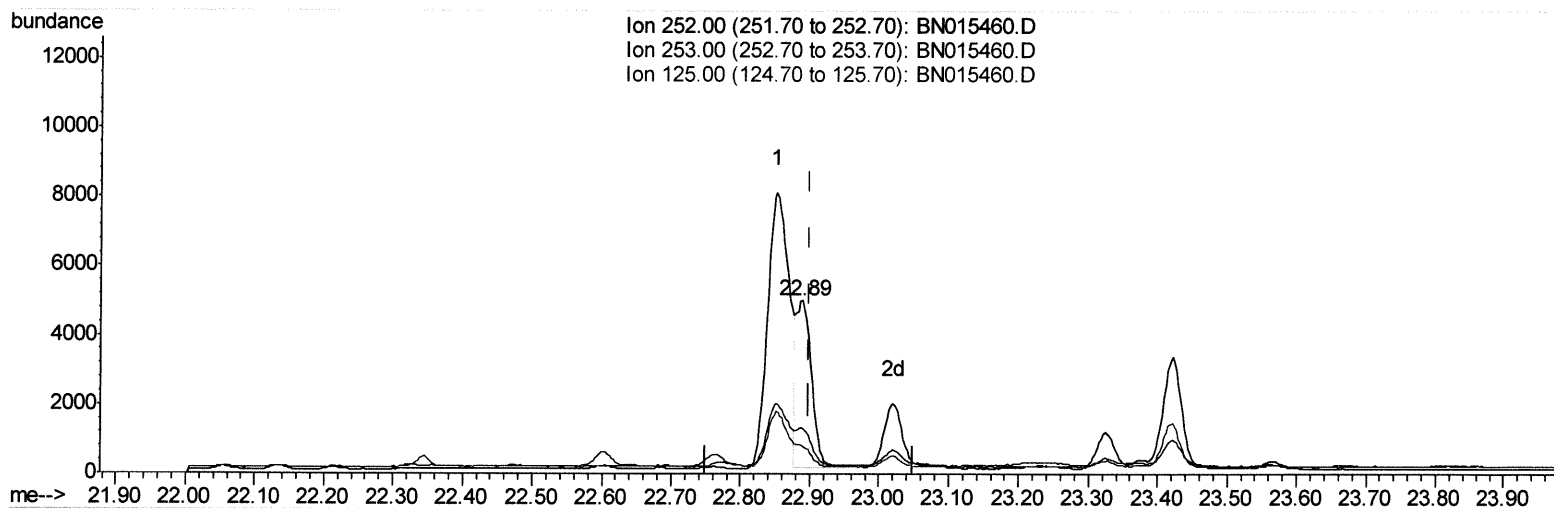
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(25) Benzo(k)fluoranthene

22.889min (-0.011) 0.17ng/ul m

response 7434

Ion	Exp%	Act%
252.00	100	100
253.00	22.60	26.81
125.00	14.40	16.30
0.00	0.00	0.00

547/14/21

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	3150	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	11022	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.34	164	5785	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.09	188	11537	0.40	ng/ul	0.00
17) Chrysene-d12	21.28	240	9190	0.40	ng/ul	0.00
23) Perylene-d12	23.52	264	8320	0.40	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.28	96	4478	1.26	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.07	152	1702	0.10	ng/ul	0.00
18) Fluoranthene-d10	19.12	212	3548	0.12	ng/ul	0.00
Target Compounds						
					Ovalue	
5) Naphthalene	10.53	128	1830	0.052	ng/ul#	91
10) Acenaphthylene	14.05	152	1942	0.065	ng/ul#	90
11) Acenaphthene	14.40	153	783	0.034	ng/ul	96
12) Fluorene	15.39	166	1371	0.054	ng/ul#	94
15) Phenanthrene	17.12	178	15604	0.373	ng/ul	99
16) Anthracene	17.22	178	6714	0.182	ng/ul	99
19) Fluoranthene	19.15	202	43696	1.029	ng/ul	99
20) Pyrene	19.51	202	25966	0.610	ng/ul	100
21) Benzo(a)anthracene	21.26	228	19591	0.575	ng/ul	97
22) Chrysene	21.32	228	15393	0.378	ng/ul	98
24) Benzo(b)fluoranthene	22.85	252	17582m	0.452	ng/ul	} 7/14/21
25) Benzo(k)fluoranthene	22.89	252	7434m	0.169	ng/ul	
26) Benzo(a)pyrene	23.42	252	6032	0.185	ng/ul#	
27) Indeno(1,2,3-cd)pyrene	25.76	276	4286	0.109	ng/ul	95
28) Dibenzo(a,h)anthracene	25.77	278	1910	0.062	ng/ul#	77

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