

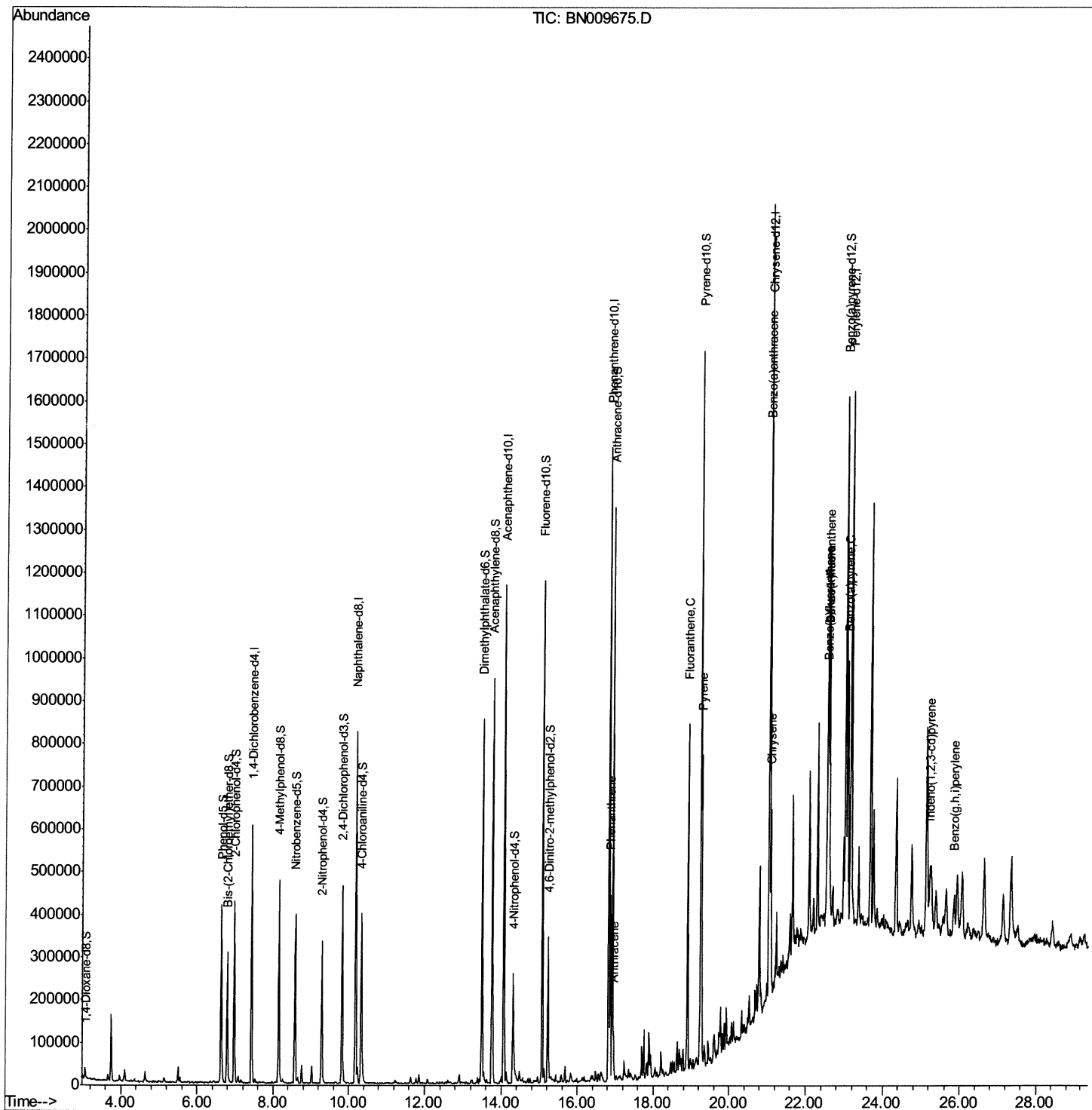
Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN020620\
Data File : BN009675.D
Acq On : 08 Feb 2020 08:58
Operator : CG/JU
Sample : L1401-19
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CB6W0

Manual Integrations
APPROVED

mohammad
2/11/2020 8:24:38 AM

Quant Time: Feb 10 03:39:22 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN020320MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 07 15:21:46 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

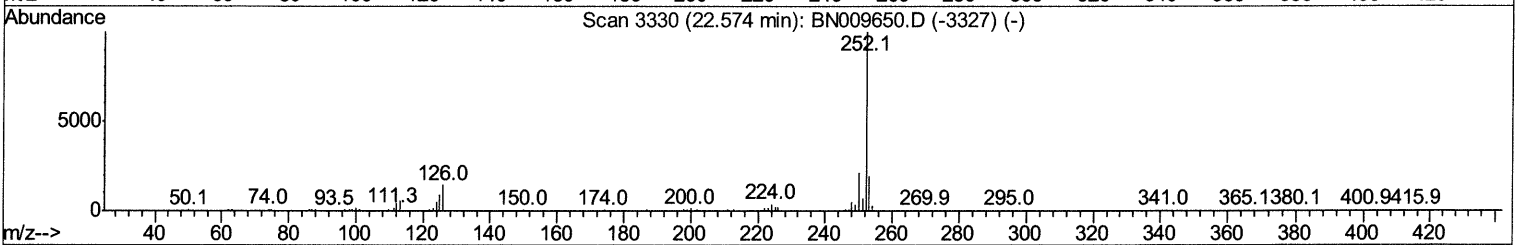
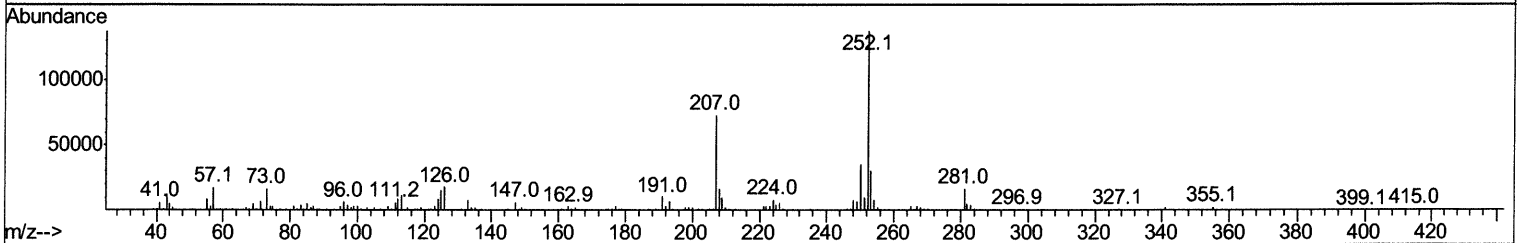
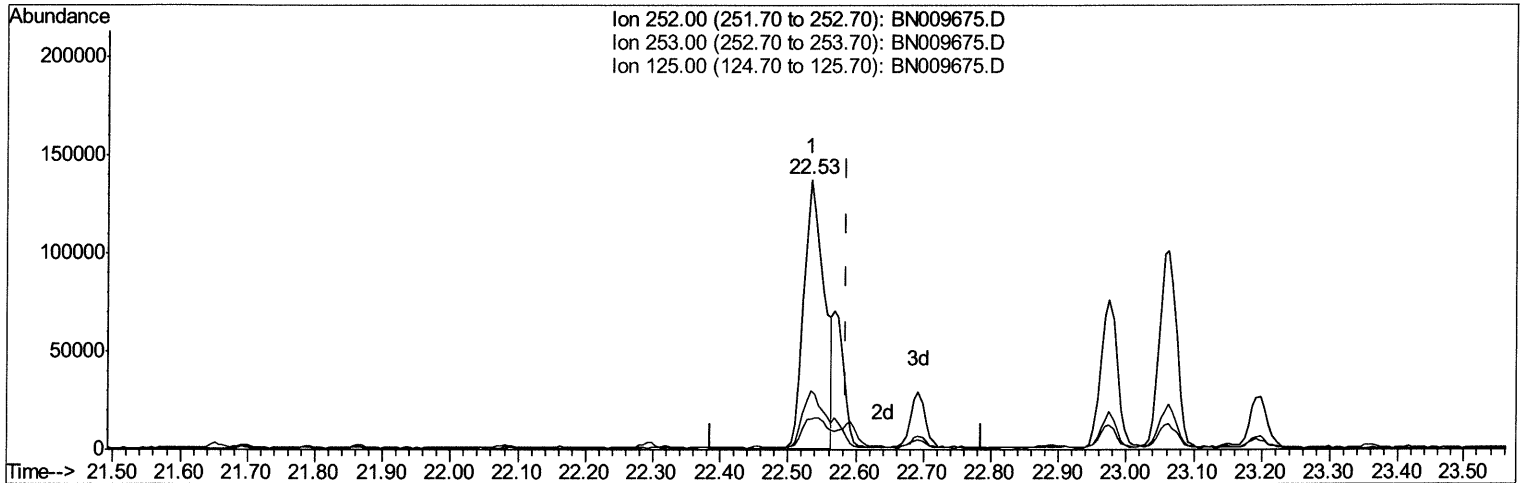
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TIC: BN009675.D

(88) Benzo(k)fluoranthene
 22.533min (-0.053) 6.02ng/ul
 response 286945

Ion	Exp%	Act%
252.00	100	100
253.00	21.20	22.00
125.00	10.00	11.51
0.00	0.00	0.00

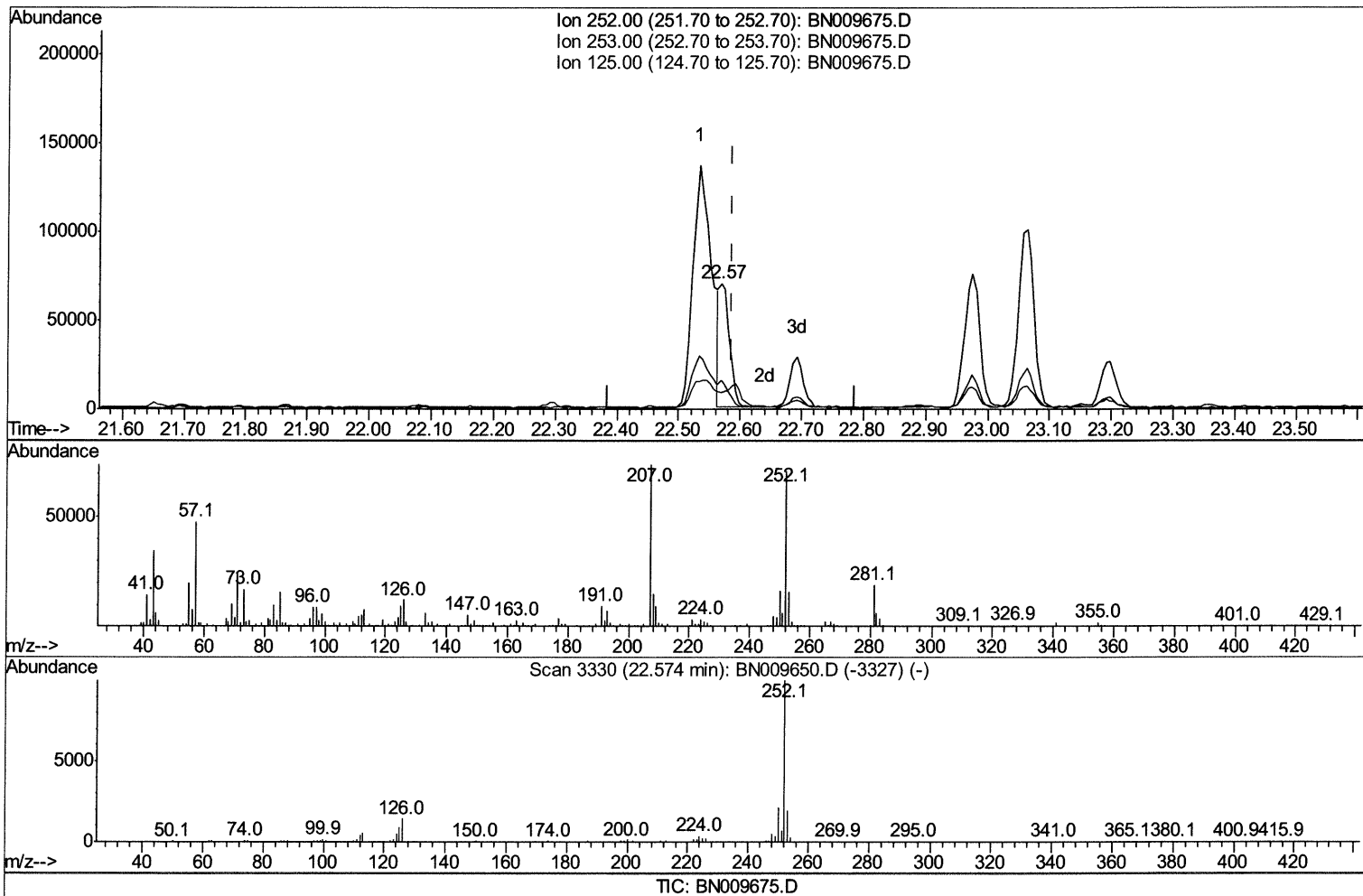
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(88) Benzo(k)fluoranthene

22.569min (-0.018) 1.59ng/ul m 2402/13/20

response 75727

Ion	Exp%	Act%
252.00	100	100
253.00	21.20	22.56
125.00	10.00	13.55#
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	7.43	152	142063	20.00	ng/ul	0.00
18) Naphthalene-d8	10.18	136	598832	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.07	164	375647	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.82	188	804395	20.00	ng/ul	0.00
77) Chrysene-d12	21.04	240	768685	20.00	ng/ul	0.00
85) Perylene-d12	23.15	264	773543	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.07	96	10422	3.01	ng/uL	0.00
5) Phenol-d5	6.63	99	202443	16.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.79	67	143089	17.19	ng/ul	0.00
9) 2-Chlorophenol-d4	6.97	132	160582	17.61	ng/ul	0.00
13) 4-Methylphenol-d8	8.14	113	160555	16.27	ng/ul	0.00
19) Nitrobenzene-d5	8.58	128	78686	17.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.29	143	86045	17.31	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.82	165	158813	17.06	ng/ul	0.00
29) 4-Chloroaniline-d4	10.33	131	210344	21.01	ng/ul	0.00
43) Dimethylphthalate-d6	13.49	166	503335	18.35	ng/ul	0.00
46) Acenaphthylene-d8	13.76	160	601045	18.14	ng/ul	0.00
51) 4-Nitrophenol-d4	14.31	143	68102	13.15	ng/ul	0.00
57) Fluorene-d10	15.07	176	436809	18.11	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.22	200	61594	12.26	ng/ul	0.00
70) Anthracene-d10	16.92	188	661871	18.11	ng/ul	0.00
78) Pyrene-d10	19.23	212	757003	18.73	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.02	264	758420	19.91	ng/ul	0.00

Target Compounds

					Ovalue	
69) Phenanthrene	16.86	178	237744	5.183	ng/ul	97
71) Anthracene	16.96	178	56021	1.204	ng/ul	97
76) Fluoranthene	18.89	202	449507	8.585	ng/ul	98
79) Pvrene	19.26	202	409176	7.292	ng/ul	98
82) Benzo(a)anthracene	21.02	228	234347	4.416	ng/ul	96
84) Chrysene	21.07	228	211974	4.012	ng/ul	99
87) Benzo(b)fluoranthene	22.53	252	286945	5.929	ng/ul	98
88) Benzo(k)fluoranthene	22.57	252	75727m >	1.589	ng/ul >	94
90) Benzo(a)pvrene	23.06	252	180623	4.028	ng/ul#	94
91) Indeno(1,2,3-cd)pyrene	25.22	276	104387	1.958	ng/ul	94
93) Benzo(a,h,i)perylene	25.84	276	101260	2.266	ng/ul#	92

JU 02/13/20

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