

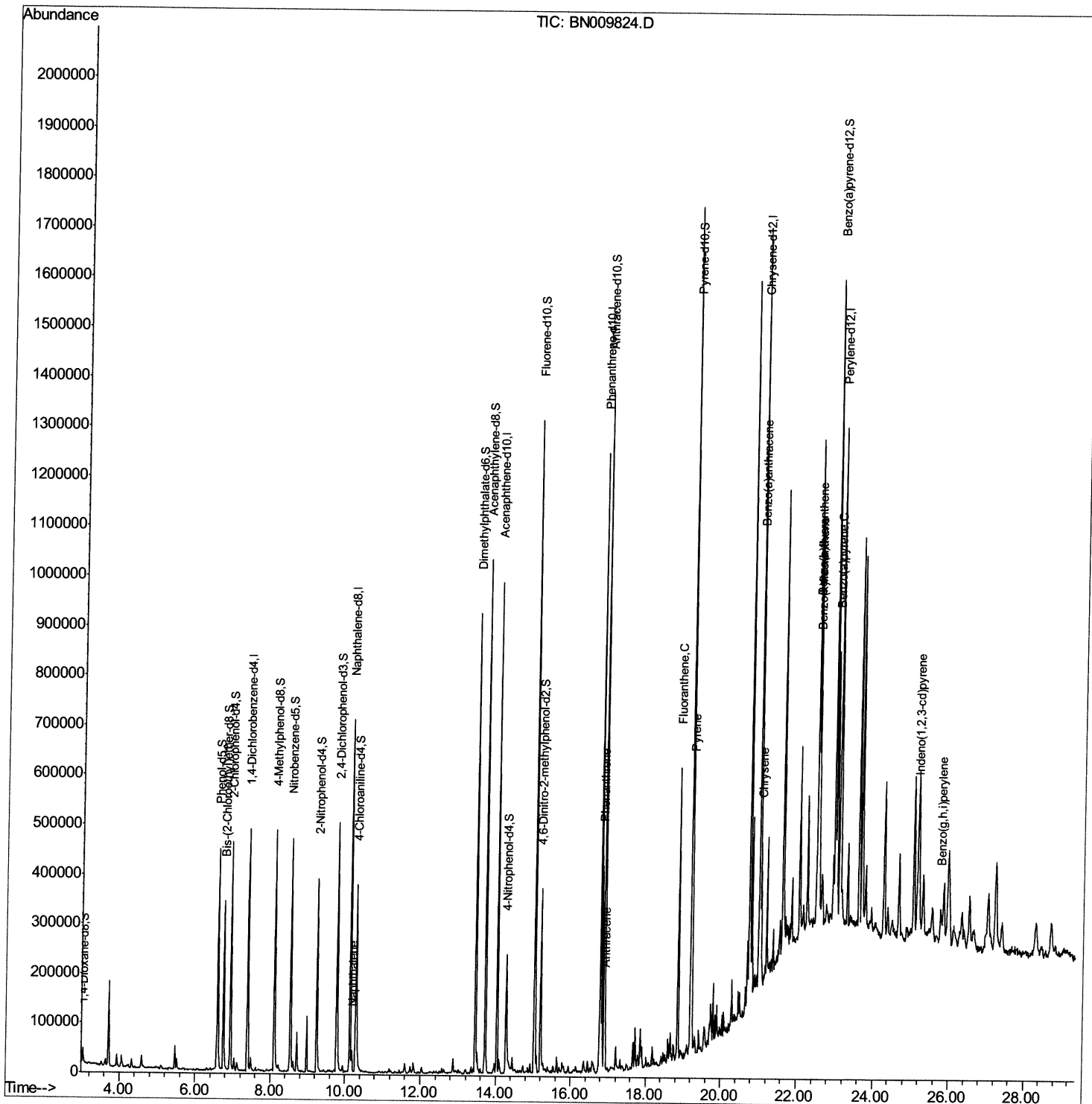
Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022220\
Data File : BN009824.D
Acq On : 22 Feb 2020 01:05
Operator : CG/JU
Sample : L1535-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CB6Q8

Manual Integrations
APPROVED

mohammad
2/24/2020 4:48:17 PM

Quant Time: Feb 22 03:04:10 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 21 22:14:17 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

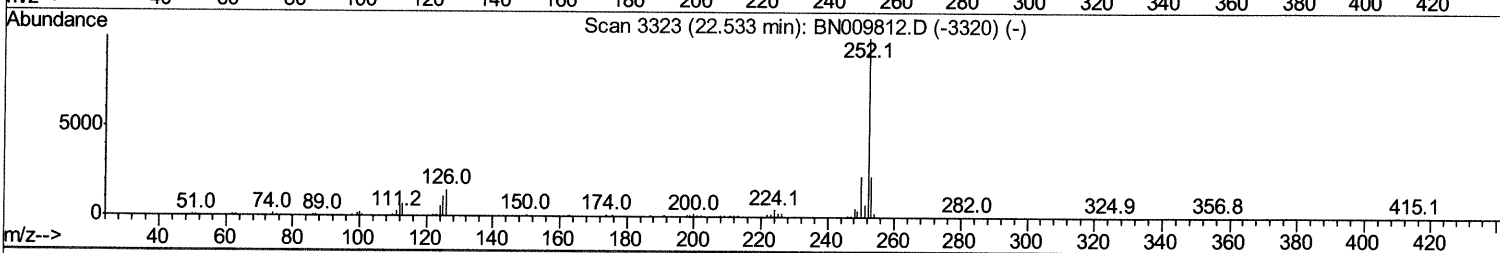
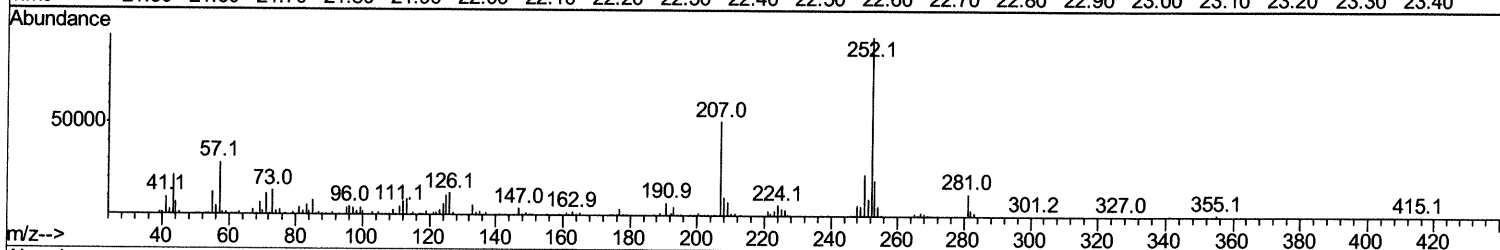
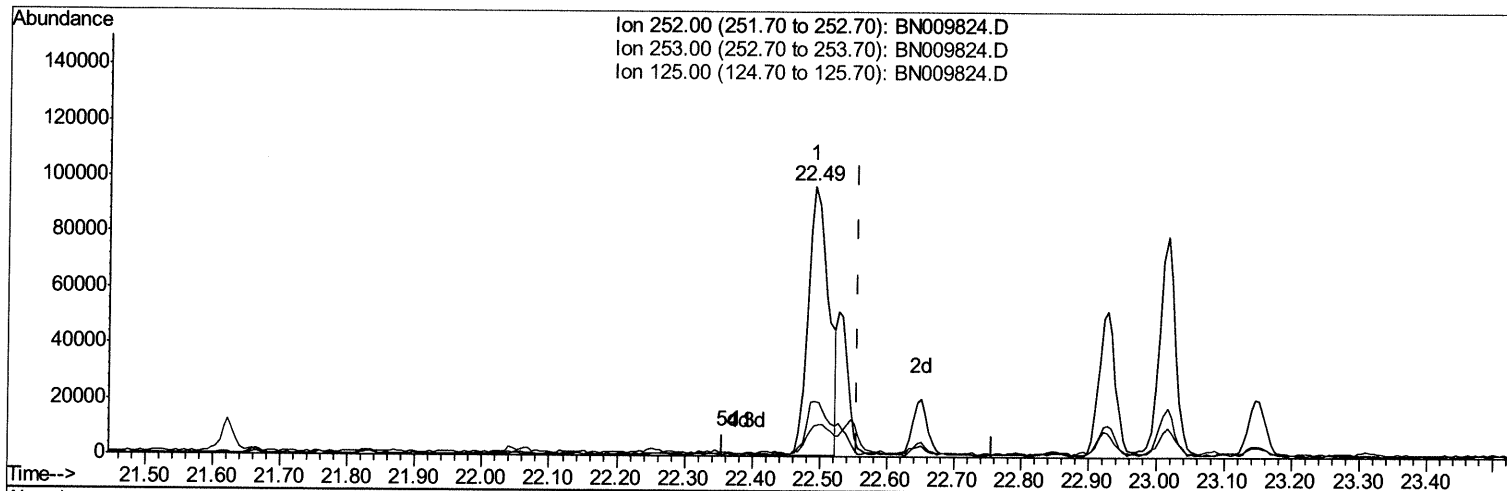
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TIC: BN009824.D

(89) Benzo(k)fluoranthene

22.492min (-0.065) 5.21ng/ul

response 203074

Ion	Exp%	Act%
252.00	100	100
253.00	22.50	20.46
125.00	10.20	11.52
0.00	0.00	0.00

Quantitation Report (Qedit)

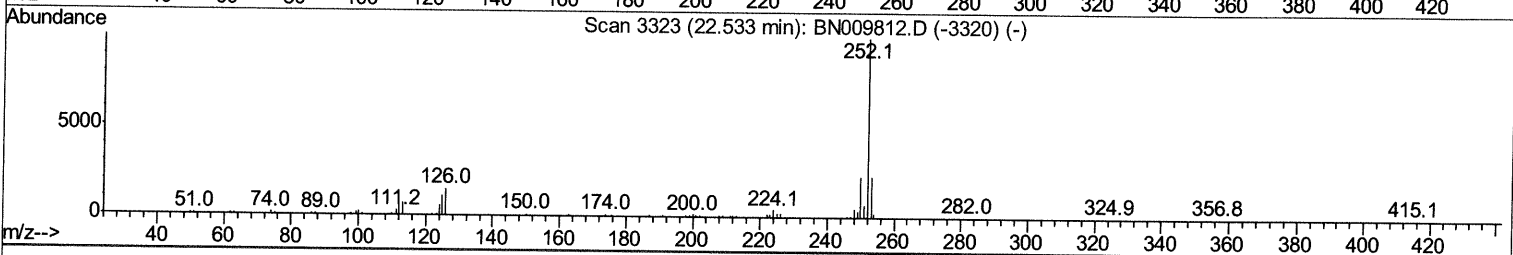
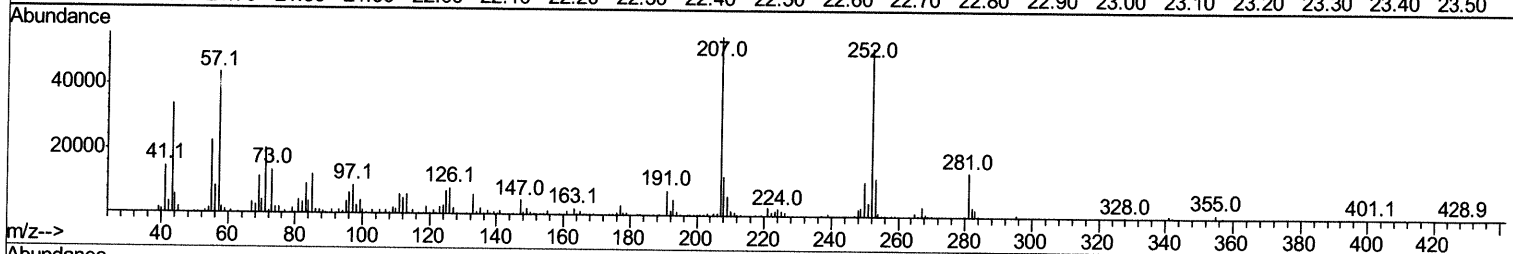
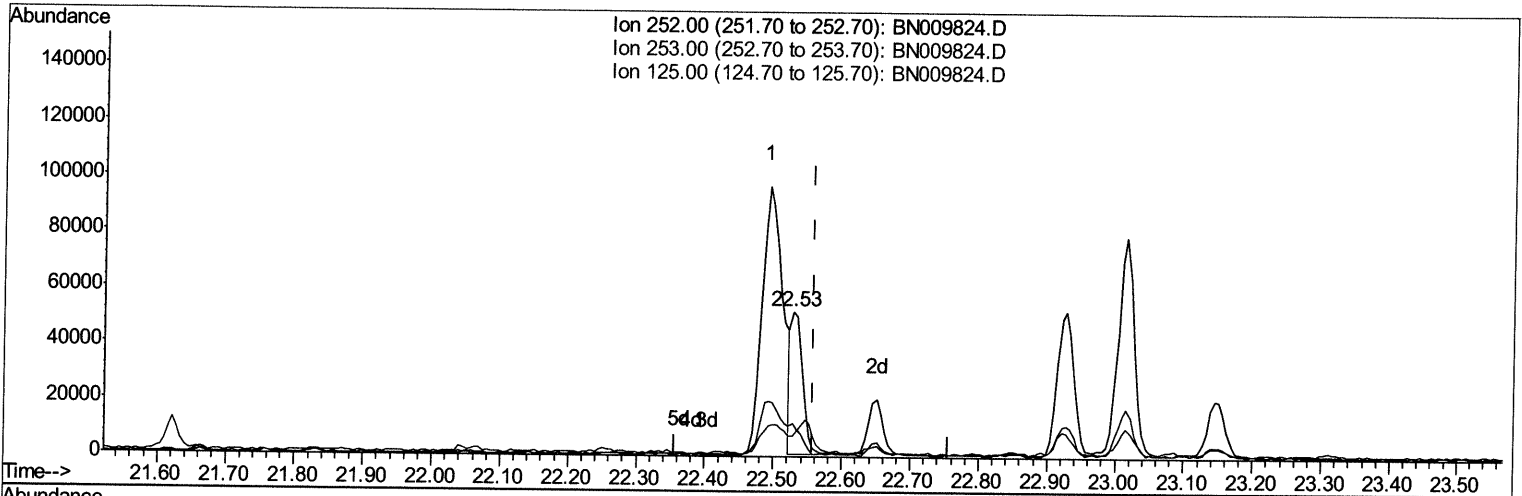
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Manual Integrations
 APPROVED

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 2/24/2020 4:48:17 PM

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 Response via : Initial Calibration



TIC: BN009824.D

(89) Benzo(k)fluoranthene

22.527min (-0.029) 1.48ng/ul m Ju 02/25/20

response 57835

Ion	Exp%	Act%
252.00	100	100
253.00	22.50	22.67
125.00	10.20	14.15#
0.00	0.00	0.00

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 Operator : CG/JU
 Sample : L1535-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
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Manual Integrations
 APPROVED

mohammad
 2/24/2020 4:48:17 PM

Quant Time: Feb 22 03:04:10 2020
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 21 22:14:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	117679	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.14	136	498316	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.03	164	313702	20.00	ng/ul	-0.02
62) Phenanthrene-d10	16.79	188	655235	20.00	ng/ul	-0.02
78) Chrysene-d12	21.00	240	582869	20.00	ng/ul	-0.02
86) Perylene-d12	23.10	264	635100	20.00	ng/ul	-0.03

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.03	96	11584	4.05	ng/uL	-0.02
5) Phenol-d5	6.59	99	212783	21.26	ng/ul	-0.02
7) Bis-(2-Chloroethyl) ether-d	6.75	67	161857	23.43	ng/ul	-0.02
9) 2-Chlorophenol-d4	6.93	132	173524	23.03	ng/ul	-0.02
13) 4-Methylphenol-d8	8.10	113	158767	19.94	ng/ul	-0.02
19) Nitrobenzene-d5	8.53	128	86399	23.70	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.25	143	94355	23.66	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	9.78	165	170618	22.25	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.29	131	197434	22.92	ng/ul	-0.02
44) Dimethylphthalate-d6	13.46	166	536062	22.88	ng/ul	-0.02
47) Acenaphthylene-d8	13.72	160	646430	23.42	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.27	143	63890	14.93	ng/ul	-0.02
58) Fluorene-d10	15.03	176	469553	23.21	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.19	200	64351	15.81	ng/ul	-0.02
71) Anthracene-d10	16.89	188	687192	22.88	ng/ul	-0.02
79) Pyrene-d10	19.19	212	770127	25.28	ng/ul	-0.02
90) Benzo(a)pyrene-d12	22.97	264	749162	23.62	ng/ul	-0.03

Target Compounds

					Ovalue	
28) Naphthalene	10.19	128	30879	1.149	ng/ul	98
70) Phenanthrene	16.83	178	217208	5.809	ng/ul	98
72) Anthracene	16.92	178	47399	1.240	ng/ul	95
77) Fluoranthene	18.86	202	321499	7.340	ng/ul	99
80) Pvrene	19.22	202	291489	6.937	ng/ul	98
83) Benzo(a)anthracene	20.99	228	164939	4.118	ng/ul	98
85) Chrysene	21.04	228	156524	3.969	ng/ul	98
88) Benzo(b)fluoranthene	22.49	252	203074	4.932	ng/ul	98
89) Benzo(k)fluoranthene	22.53	252	57835m >	1.483	ng/ul >	JU 02/25/20
91) Benzo(a)pyrene	23.02	252	127945	3.455	ng/ul	98
92) Indeno(1,2,3-cd)pyrene	25.15	276	77572	1.764	ng/ul#	93
94) Benzo(a,h,i)perylene	25.77	276	73942	2.006	ng/ul	96

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