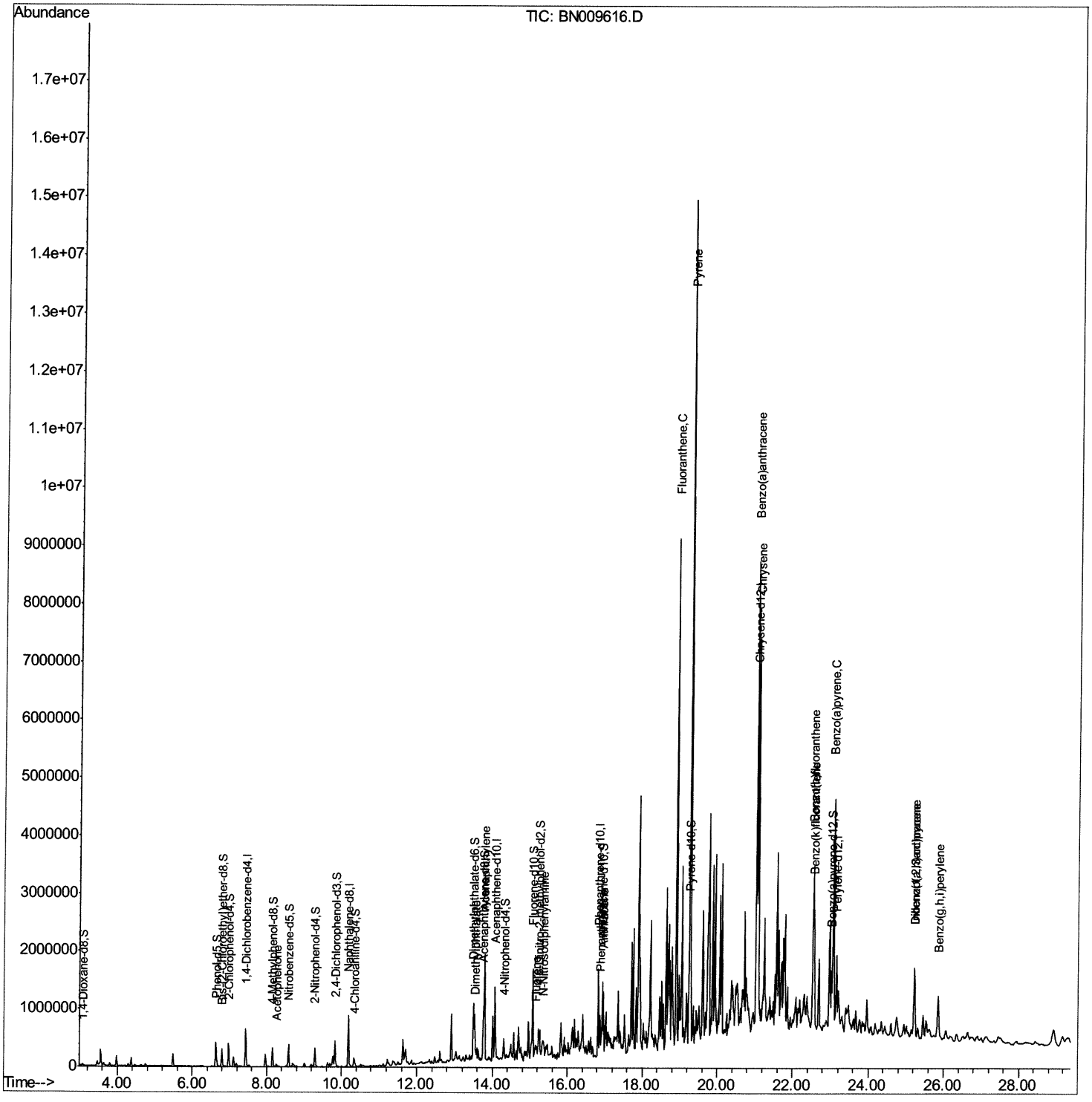
Quantitation Report (QT Reviewed)

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN020620\
Data File : BN009616.D
Acq On : 06 Feb 2020 19:11
Operator : CG/JU
Sample : L1323-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Feb 06 22:59:12 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN020320MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Feb 03 16:39:56 2020
Response via : Initial Calibration

Instrument :
BNA_N
ClientSampleId :
GAT42

Manual Integrations APPROVED
mohammad
2/10/2020 8:19:31 AM



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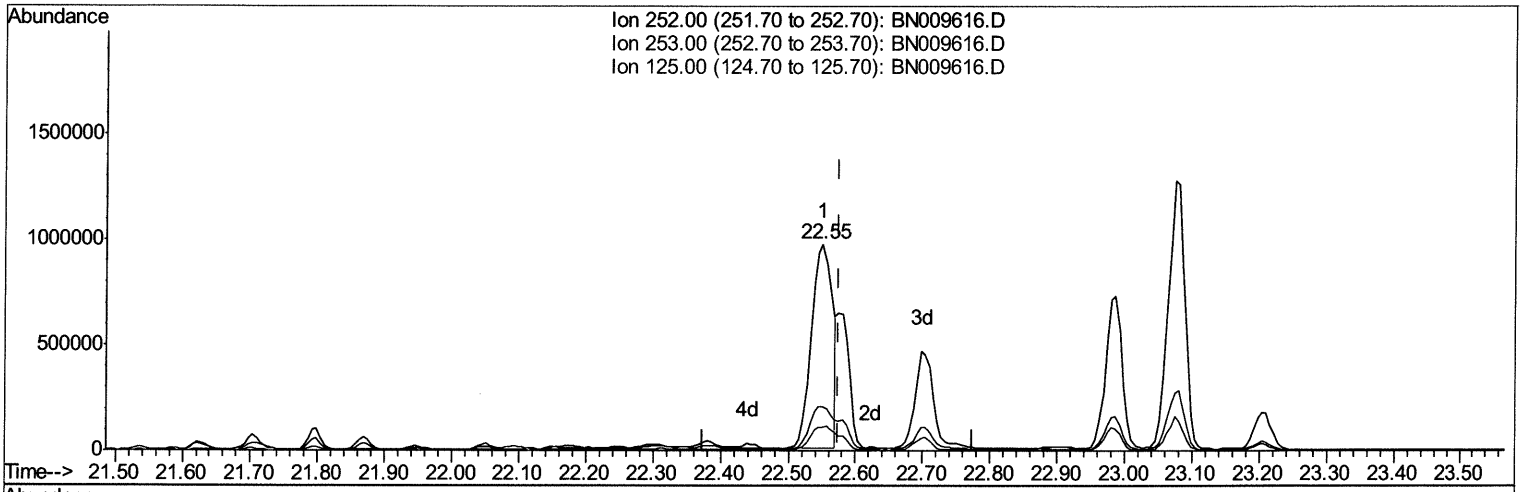

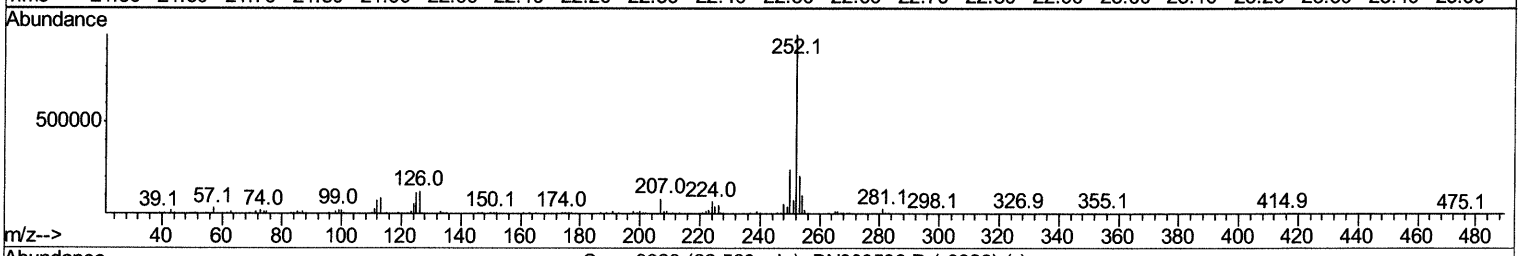

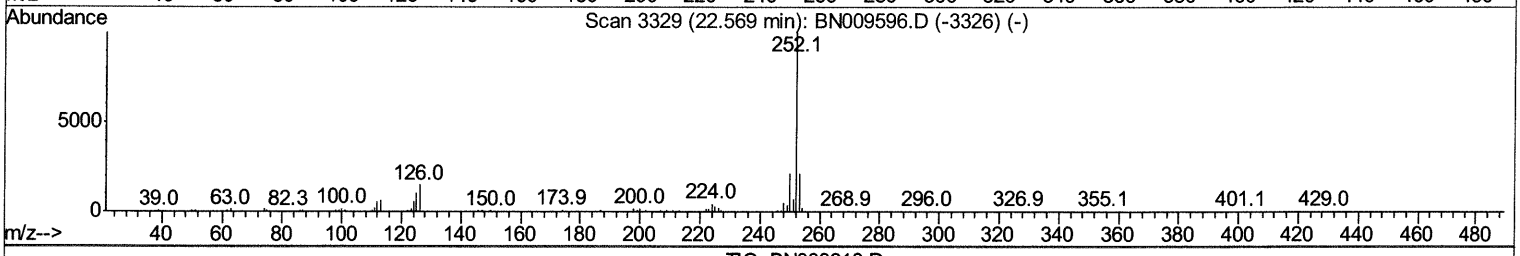

TIC: BN009616.D

(88) Benzo(k)fluoranthene

22.551min (-0.024) 45.02ng/ul

response 2114426

Ion	Exp%	Act%
252.00	100	100
253.00	21.20	21.61
125.00	10.00	11.53
0.00	0.00	0.00

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mohammad
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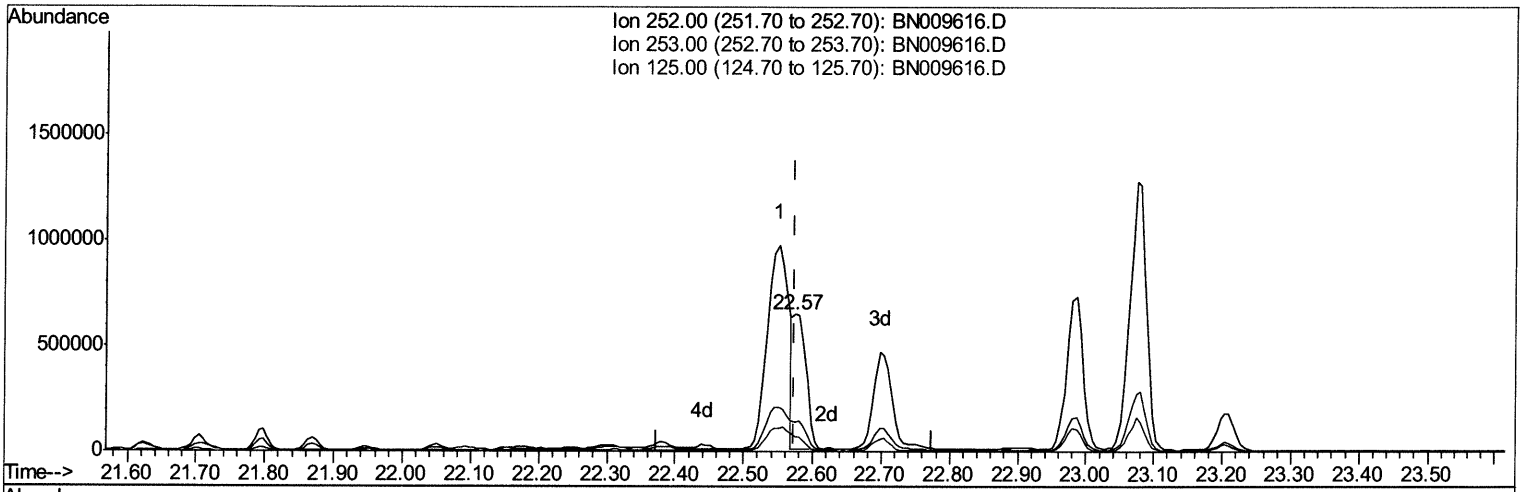

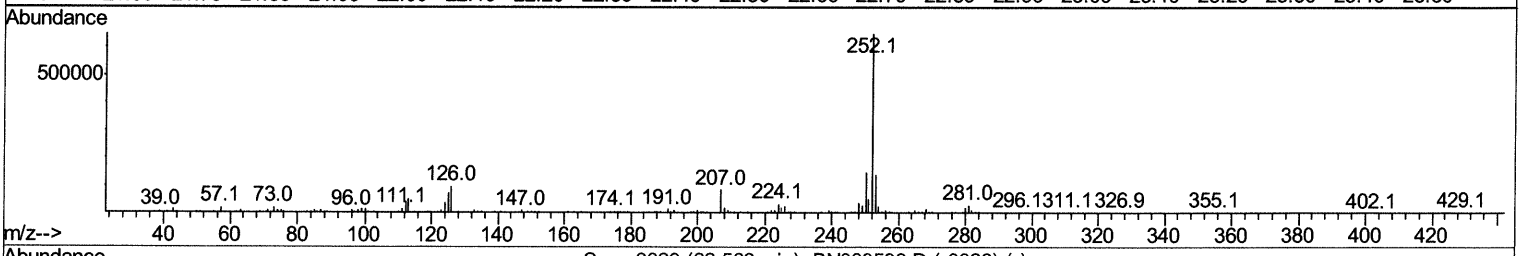

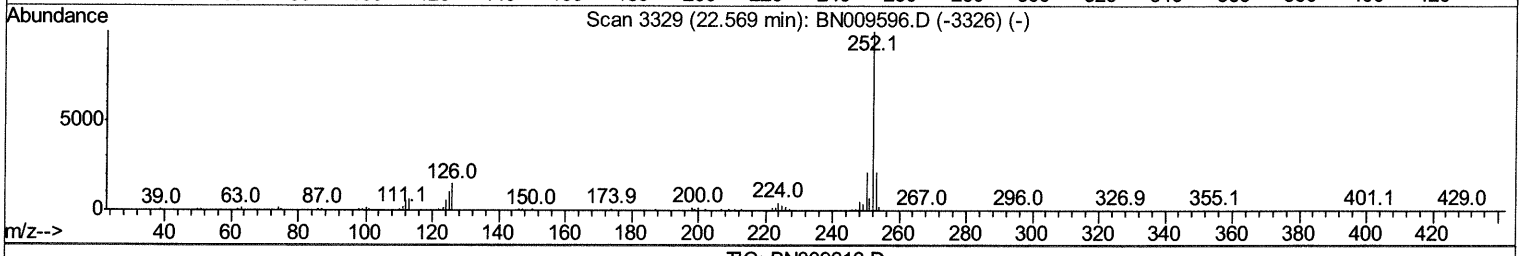

TIC: BN009616.D

(88) Benzo(k)fluoranthene

22.574min (0.000) 17.18ng/ul m JU 02/13/20

response 806622

Ion	Exp%	Act%
252.00	100	100
253.00	21.20	21.30
125.00	10.00	11.21
0.00	0.00	0.00

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mohammad
2/10/2020 8:19:31 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	152963	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.18	136	641078	20.00	ng/ul	-0.01
35) Acenaphthene-d10	14.07	164	391258	20.00	ng/ul	-0.01
61) Phenanthrene-d10	16.83	188	782852	20.00	ng/ul	0.00
77) Chrysene-d12	21.05	240	664061	20.00	ng/ul	0.00
85) Perylene-d12	23.16	264	762211	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.08	96	9205	2.47	ng/uL	0.00
5) Phenol-d5	6.63	99	192238	14.41	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.79	67	140219	15.65	ng/ul	0.00
9) 2-Chlorophenol-d4	6.98	132	152147	15.49	ng/ul	0.00
13) 4-Methylphenol-d8	8.15	113	107802	10.15	ng/ul	-0.01
19) Nitrobenzene-d5	8.58	128	74991	15.73	ng/ul	0.00
22) 2-Nitrophenol-d4	9.29	143	83498	15.69	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	9.82	165	150288	15.08	ng/ul	0.00
29) 4-Chloroaniline-d4	10.34	131	82762	7.72	ng/ul	0.00
43) Dimethylphthalate-d6	13.50	166	503456	17.63	ng/ul	0.00
46) Acenaphthylene-d8	13.76	160	594295	17.22	ng/ul	-0.01
51) 4-Nitrophenol-d4	14.31	143	72175	13.38	ng/ul	0.00
57) Fluorene-d10	15.07	176	437967	17.44	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.23	200	68903	14.09	ng/ul	0.00
70) Anthracene-d10	16.93	188	643638	18.09	ng/ul	0.00
78) Pyrene-d10	19.24	212	699814	20.05	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.03	264	671819	17.90	ng/ul	0.00
Target Compounds						Qvalue
14) Acetophenone	8.26	105	18552	1.091	ng/ul	95
44) Dimethylphthalate	13.55	163	89219	2.943	ng/ul	99
47) Acenaphthylene	13.79	152	1187833	31.452	ng/ul	99
58) Fluorene	15.15	166	98280	3.289	ng/ul#	37
64) N-Nitrosodiphenylamine	15.35	169	41556	1.776	ng/ul#	76
69) Phenanthrene	16.87	178	396108	8.872	ng/ul	98
71) Anthracene	16.96	178	638724	14.100	ng/ul	98
76) Fluoranthene	18.90	202	4308454	84.555	ng/ul	98
79) Pyrene	19.27	202	8388561	173.044	ng/ul	96
82) Benzo(a)anthracene	21.03	228	3466501	75.605	ng/ul#	78
84) Chrysene	21.09	228	3207568	70.272	ng/ul	94
87) Benzo(b)fluoranthene	22.55	252	2114426	44.339	ng/ul	98
88) Benzo(k)fluoranthene	22.57	252	806622m	17.175	ng/ul	98
90) Benzo(a)pyrene	23.07	252	2164932	48.999	ng/ul	96
91) Indeno(1,2,3-cd)pyrene	25.23	276	917566	17.468	ng/ul	95
92) Dibenzo(a,h)anthracene	25.23	278	349738	7.910	ng/ul#	95
93) Benzo(g,h,i)perylene	25.85	276	769513	17.479	ng/ul	96

JU 02/13/20

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