

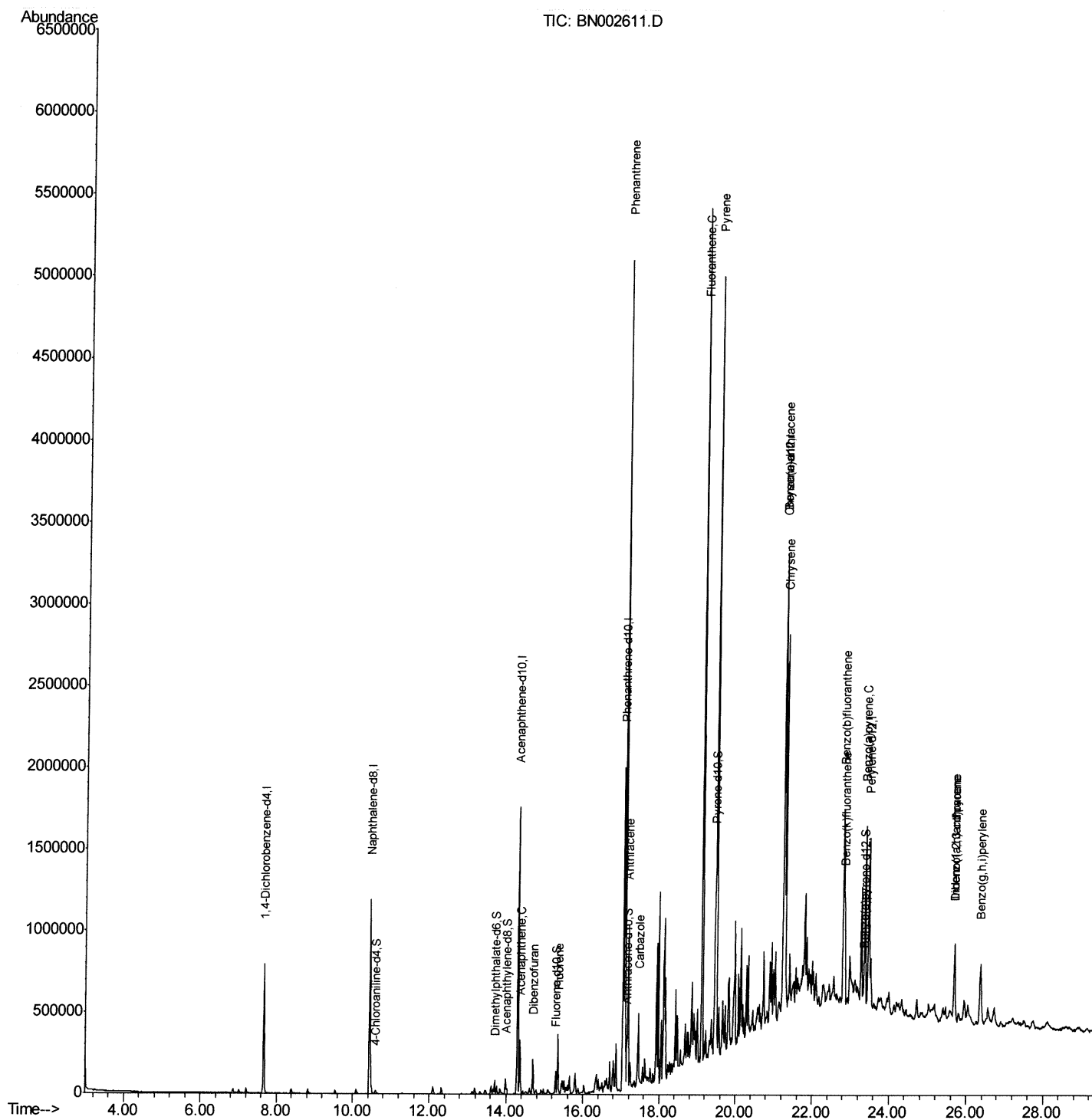
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090518\
Data File : BN002611.D
Acq On : 06 Sep 2018 11:25
Operator : JU/SJ
Sample : J4688-08DL 20X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A3YY6DL

Manual Integrations
APPROVED

Sohil
9/6/2018 4:30:05 PM

Quant Time: Sep 06 13:49:40 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 06 06:54:57 2018
Response via : Initial Calibration



Quantitation Report (Qedit)

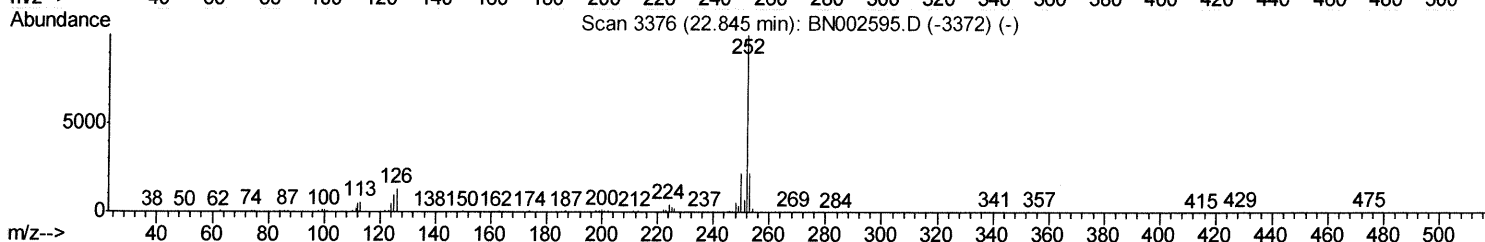
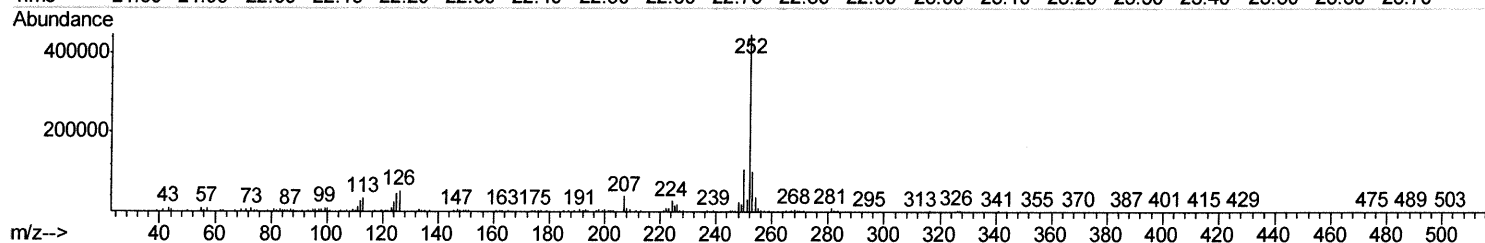
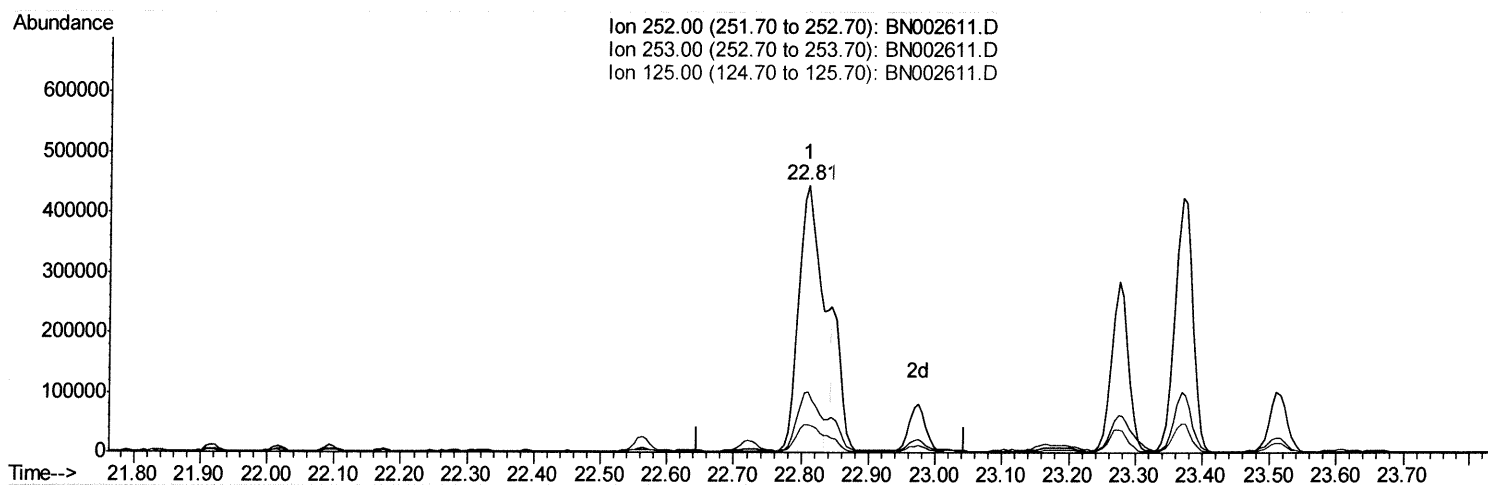
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Response via : Initial Calibration



TIC: BN002611.D

(88) Benzo(k)fluoranthene

22.810min (-0.035) 23.83ng/ui

response 967667

Ion	Exp%	Act%
252.00	100	100
253.00	21.60	22.52
125.00	10.20	10.68
0.00	0.00	0.00

Quantitation Report (Qedit)

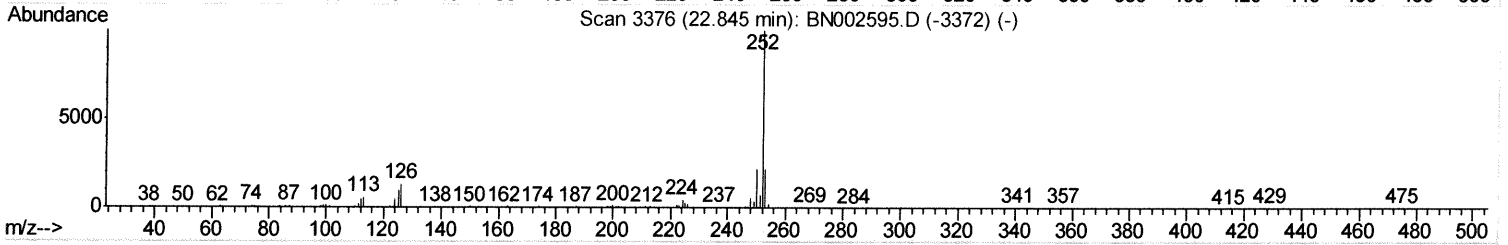
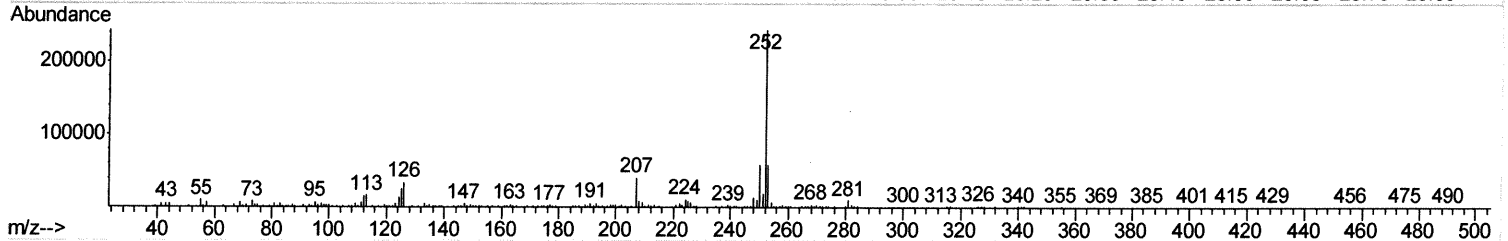
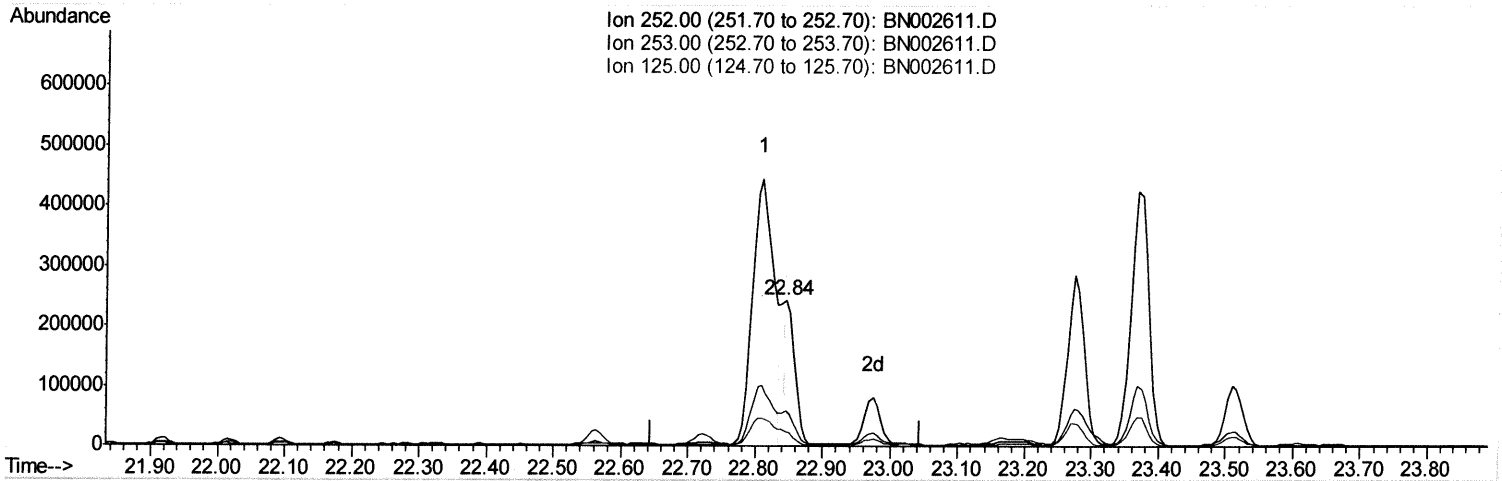
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Manual Integrations
APPROVED

Sohil
9/6/2018 4:30:05 PM

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Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 06 06:54:57 2018
Response via : Initial Calibration



TIC: BN002611.D

(88) Benzo(k)fluoranthene

22.845min (-0.000) 8.41ng/ul m

response 341605

Ion	Exp%	Act%
252.00	100	100
253.00	21.60	24.10
125.00	10.20	10.27
0.00	0.00	0.00

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 Operator : JU/SJ
 Sample : J4688-08DL 20X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
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Manual Integrations
 APPROVED

Sohil
 9/6/2018 4:30:05 PM

Quant Time: Sep 06 13:49:40 2018
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 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 06 06:54:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	215141	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	967179	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.30	164	555563	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.05	188	1188353	20.00	ng/ul	0.00
77) Chrysene-d12	21.25	240	824451	20.00	ng/ul	0.00
85) Perylene-d12	23.47	264	718954	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.24	96	597	0.13	ng/uL	0.00
5) Phenol-d5	6.86	99	17420	0.88	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.01	67	11838	0.96	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	14827	0.99	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	15126	0.96	ng/ul	0.01
19) Nitrobenzene-d5	8.82	128	7136	0.93	ng/ul	0.00
22) 2-Nitrophenol-d4	9.53	143	7640	0.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	14417	0.97	ng/ul	0.02
29) 4-Chloroaniline-d4	10.59	131	16617	1.03	ng/ul	0.02
43) Dimethylphthalate-d6	13.71	166	55910	1.21	ng/ul	0.00
46) Acenaphthylene-d8	13.99	160	65321	1.16	ng/ul	0.00
51) 4-Nitrophenol-d4	14.59	143	271	0.03	ng/ul	0.08
57) Fluorene-d10	15.29	176	52592	1.32	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.15	188	77701	1.37	ng/ul	0.00
78) Pyrene-d10	19.45	212	68256	1.73	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.33	264	50623	1.35	ng/ul	0.00

Target Compounds

					Qvalue
49) Acenaphthene	14.36	153	131292	3.460	ng/ul 98
53) Dibenzofuran	14.70	168	124523	2.326	ng/ul 93
58) Fluorene	15.35	166	149516	3.414	ng/ul 98
69) Phenanthrene	17.09	178	2951626	45.055	ng/ul 99
71) Anthracene	17.18	178	630992	9.497	ng/ul 99
74) Carbazole	17.45	167	259809	4.409	ng/ul 98
76) Fluoranthene	19.12	202	2996803	40.192	ng/ul 100
79) Pyrene	19.48	202	3004573	60.439	ng/ul 100
82) Benzo(a)anthracene	21.23	228	1207989	23.916	ng/ul 99
84) Chrysene	21.29	228	1125253	23.539	ng/ul 97
87) Benzo(b)fluoranthene	22.81	252	967667	22.454	ng/ul 99
88) Benzo(k)fluoranthene	22.84	252	341605m	8.414	ng/ul
90) Benzo(a)pyrene	23.37	252	778916	18.974	ng/ul 97
91) Indeno(1,2,3-cd)pyrene	25.68	276	428658	8.777	ng/ul# 92
92) Dibenzo(a,h)anthracene	25.69	278	122176	2.992	ng/ul# 80
93) Benzo(g,h,i)perylene	26.36	276	426782	10.473	ng/ul 96

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

S 9/10/18