

(LSC Reviewed)

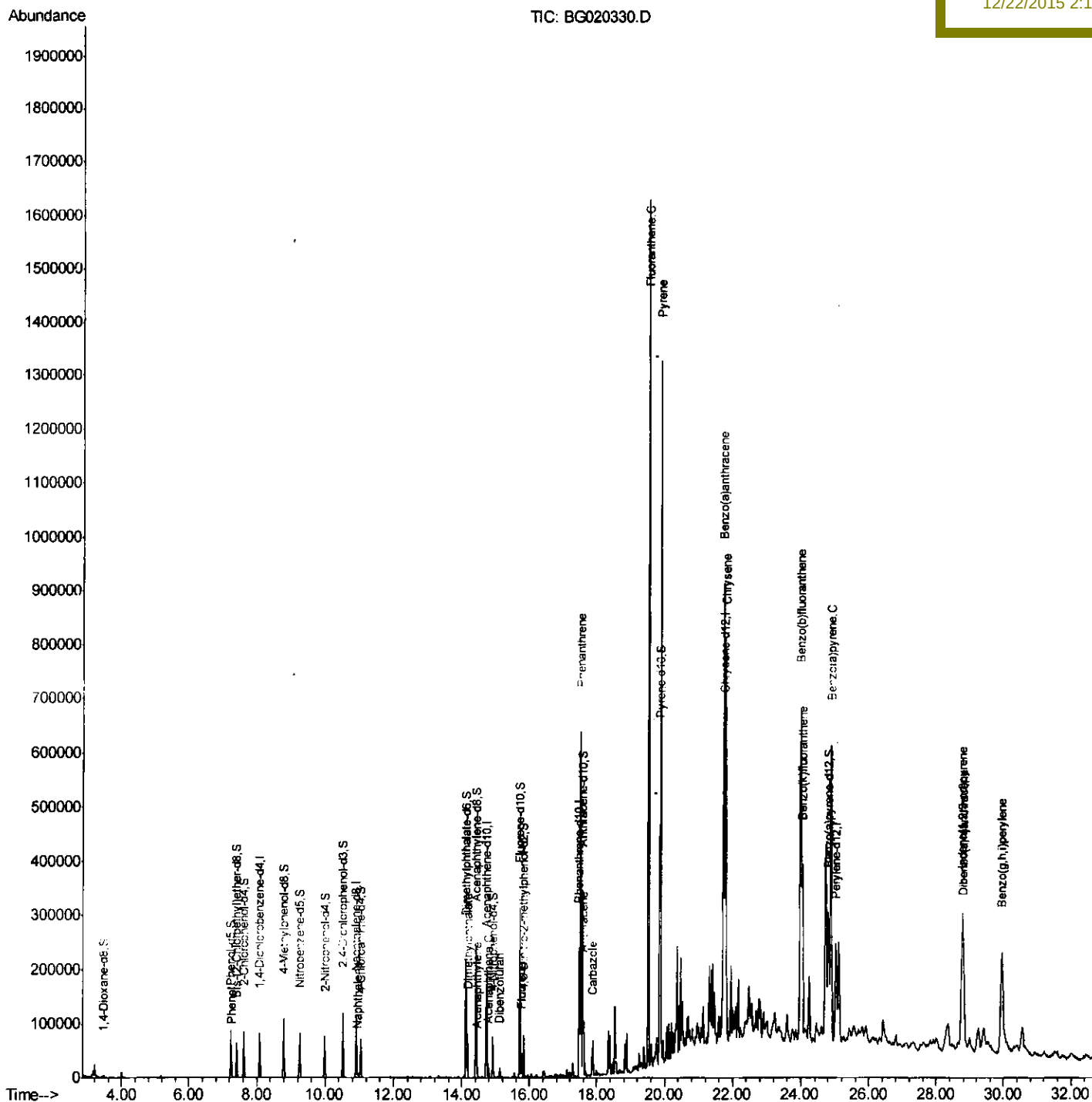
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Data Path : Z:\EPCHEM1\BNA G\DATA\BG121915\
Data File : BG020330.D
Acq On    : 20 Dec 2015    2:54
Operator  : UM/SJ
Sample    : G4866-15
Misc      :
ALS Vial  : 61    Sample Multiplier: 1
```

Instrument :
BNA_G
ClientSampleId :
E43M5

Quant Time: Dec 20 06:37:43 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION
Qlast Update : Sun Dec 20 05:53:46 2015
Response via : Initial Calibration

Manual Integrations APPROVED

apatel
12/22/2015 2:14:34 PM



Quantitation Report (Qedit)

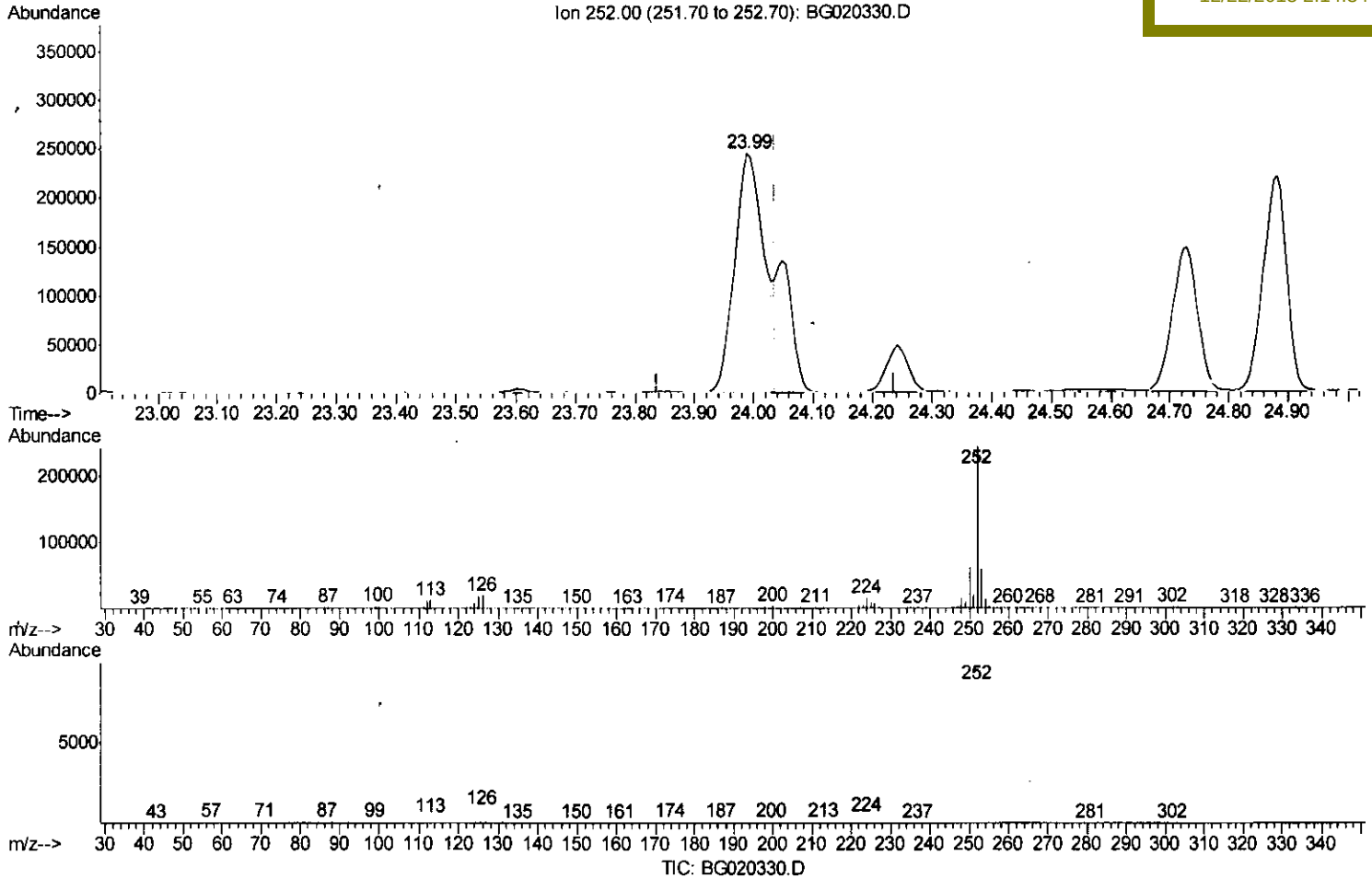
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(89) Benzo(k)fluoranthene
 23.989min (-0.048) 74.78ng/ul
 response 826953

Ion	Exp%	Act%
252.00	100	100
253.00	22.40	23.93
125.00	7.20	7.43
0.00	0.00	0.00

Quantitation Report (Qedit)

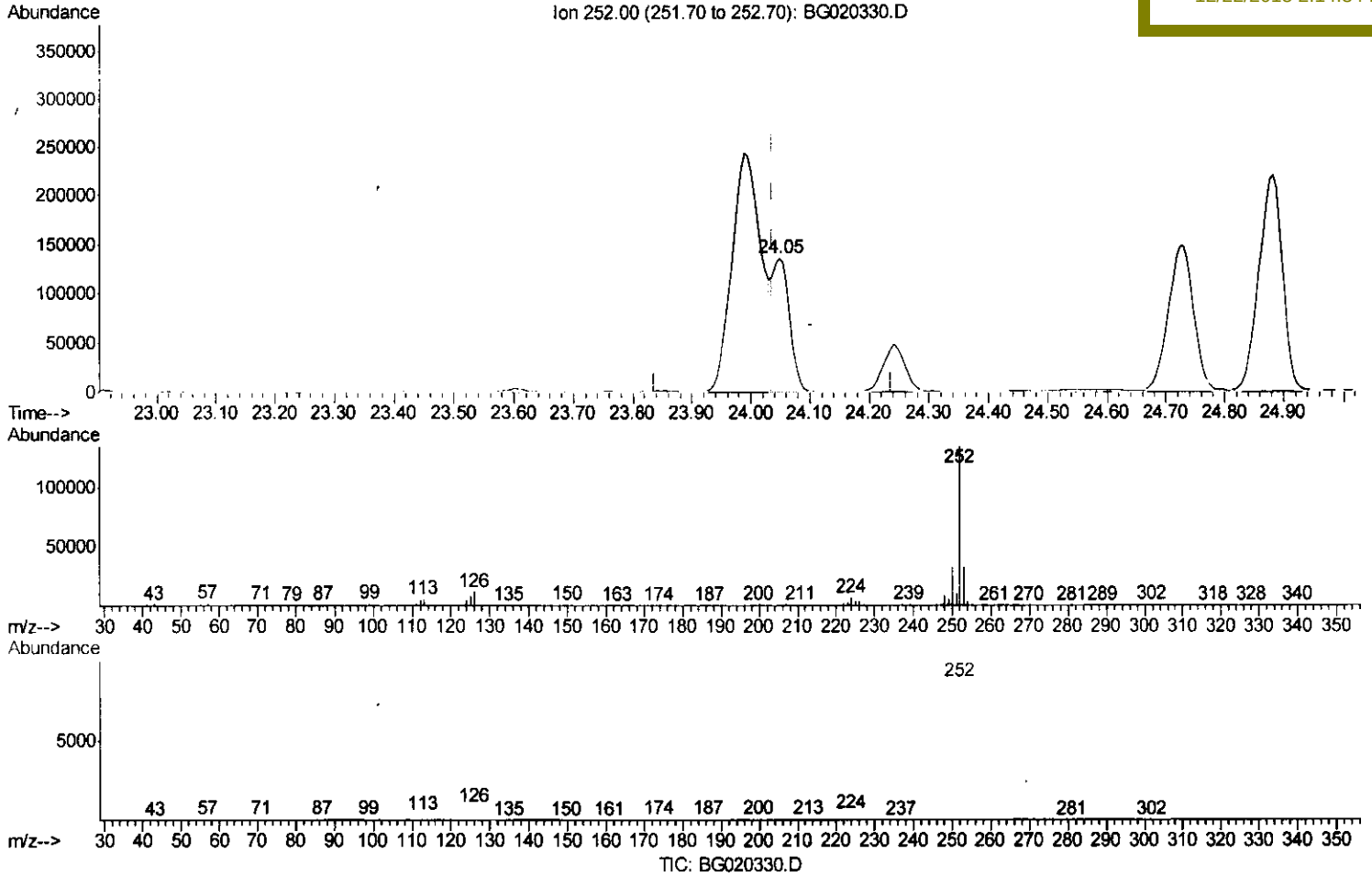
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 Data File : BG020330.D
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 APPROVED

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(89) Benzo(k)fluoranthene

24.048min (+0.011) 25.58ng/ul m U.M
 response 282858 12/22/15

Ion	Exp%	Act%
252.00	100	100
253.00	22.40	23.94
125.00	7.20	6.52
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121915\
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 Sample : G4866-15
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	23769	20.00	ng/ul	0.00
18) Naphthalene-d8	10.90	136	100499	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.72	164	68722	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	156453	20.00	ng/ul	0.00
78) Chrysene-d12	21.75	240	195165	20.00	ng/ul	0.01
86) Perylene-d12	25.02	264	191467	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	2003	4.65	ng/uL	0.00
5) Phenol-d5	7.24	99	52611	25.10	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.41	67	31454	25.15	ng/ul	0.00
9) 2-Chlorophenol-d4	7.62	132	42005	27.55	ng/ul	0.00
13) 4-Methylphenol-d8	8.79	113	43556	24.30	ng/ul	0.00
19) Nitrobenzene-d5	9.26	128	21264	27.14	ng/ul	0.00
22) 2-Nitrophenol-d4	9.98	143	26413	30.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.52	165	47665	26.99	ng/ul	0.00
29) 4-Chloroaniline-d4	11.04	131	42319	21.33	ng/ul	0.00
44) Dimethylphthalate-d6	14.12	166	151537	27.26	ng/ul	0.00
47) Acenaphthylene-d8	14.42	160	179145	27.70	ng/ul	0.00
52) 4-Nitrophenol-d4	14.91	143	20936	21.00	ng/ul	0.00
58) Fluorene-d10	15.71	176	134343	26.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	20538	22.14	ng/ul	0.00
71) Anthracene-d10	17.57	188	202985	28.16	ng/ul	0.00
79) Pyrene-d10	19.85	212	239112	26.29	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.80	264	273055	28.59	ng/ul	0.02

Target Compounds

					Ovalue	
6) Phenol	7.27	94	2670	1.19	ng/ul	96
28) Naphthalene	10.96	128	8247	1.57	ng/ul	99
45) Dimethylphthalate	14.17	163	53739	9.47	ng/ul	98
48) Acenaphthylene	14.45	152	10783	1.57	ng/ul	96
50) Acenaphthene	14.79	153	9237	2.00	ng/ul	98
54) Dibenzofuran	15.12	168	12522	1.82	ng/ul	92
59) Fluorene	15.77	166	21783	3.94	ng/ul	91
70) Phenanthrene	17.51	178	426983	49.89	ng/ul	99
72) Anthracene	17.60	178	101073	11.64	ng/ul	99
75) Carbazole	17.87	167	45432	6.23	ng/ul	97
77) Fluoranthene	19.52	202	1067557	106.74	ng/ul	99
80) Pyrene	19.88	202	852244	72.15	ng/ul	99
83) Benzo(a)anthracene	21.73	228	551681	48.45	ng/ul	97
85) Chrysene	21.80	228	555393	51.78	ng/ul	95
88) Benzo(b)fluoranthene	23.99	252	826953	71.76	ng/ul	95
89) Benzo(k)fluoranthene	24.05	252	282858m	25.58	ng/ul	96
91) Benzo(a)pyrene	24.88	252	598613	54.80	ng/ul	96
92) Indeno(1,2,3-cd)pyrene	28.78	276	480344	36.92	ng/ul	95
93) Dibenzo(a,h)anthracene	28.81	278	100573	9.31	ng/ul	93
94) Benzo(g,h,i)perylene	29.95	276	440809	41.33	ng/ul	100

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