

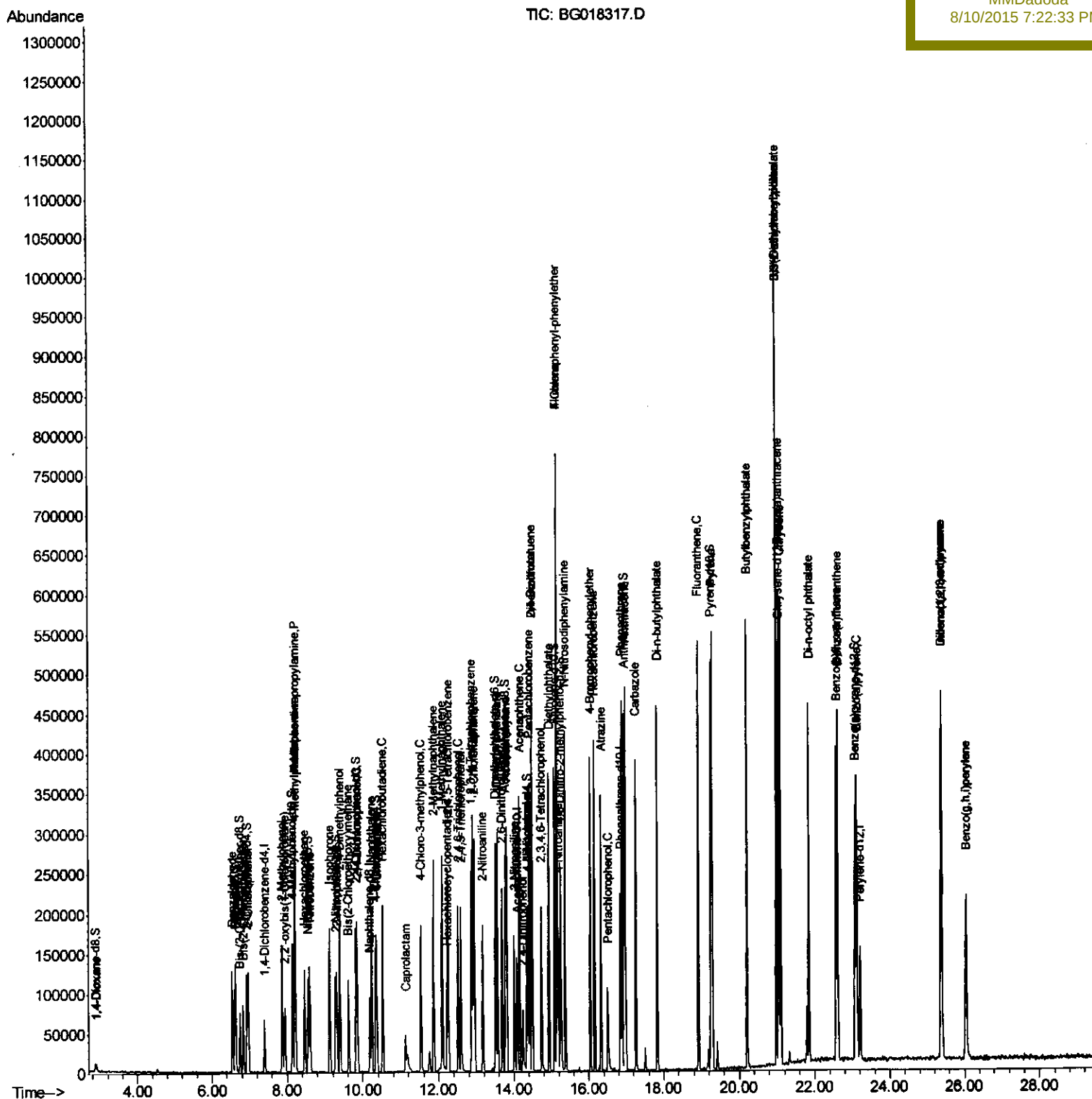
```
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\  
Data File : BG018317.D  
Acq On    : 8 Aug 2015    6:40  
Operator  : UM/TP  
Sample    : SSTD04075  
Misc      :  
ALS Vial  : 5    Sample Multiplier: 1
```

Quant Time: Aug 08 08:06:52 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations

MMDadoda
8/10/2015 7:22:33 PM

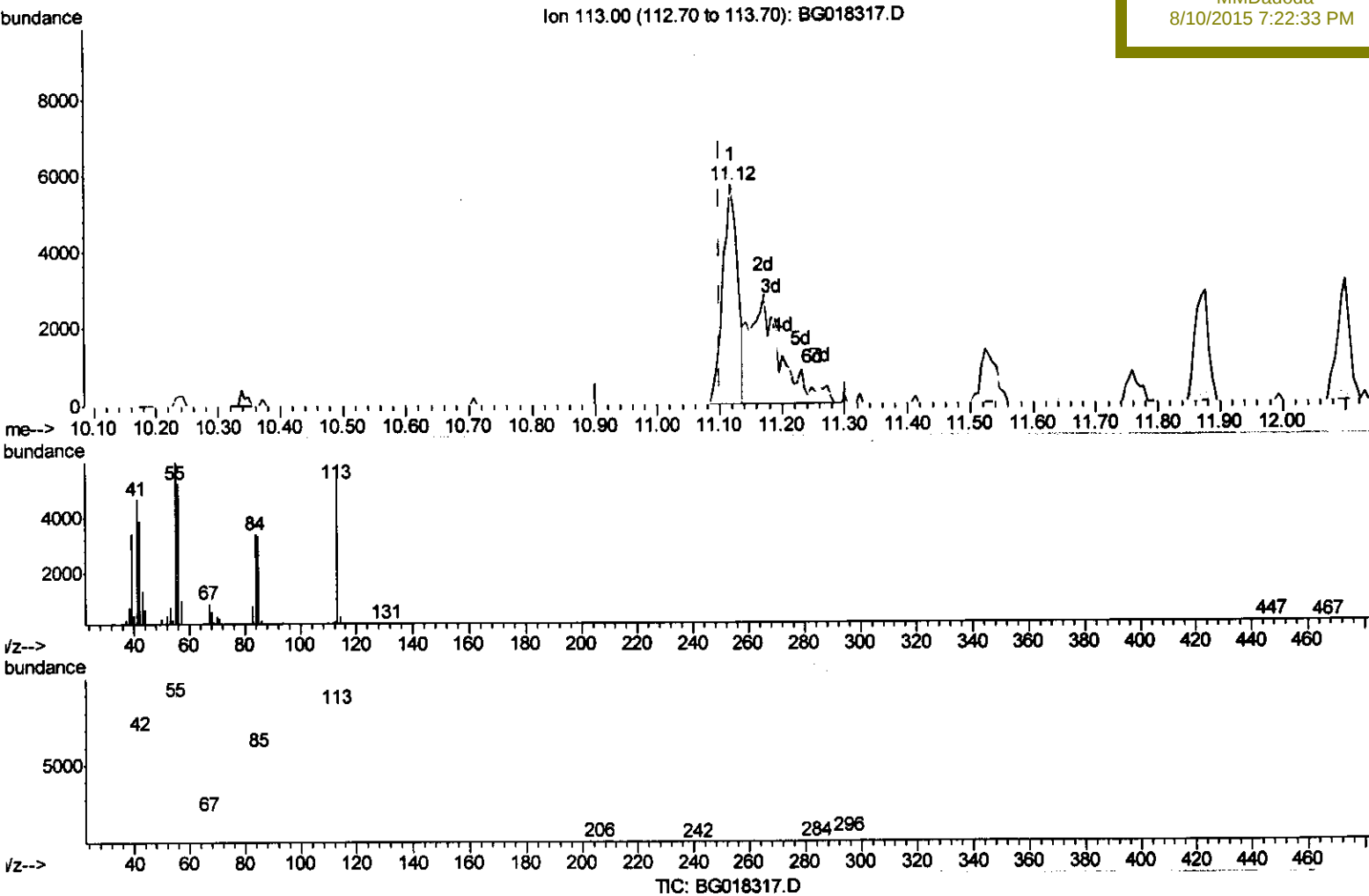


Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampled :
SSTD04075

Manual Integrations
APPROVED
MMDadoda
8/10/2015 7:22:33 PM



(32) Caprolactam
11.119min (+0.018) 18.88ng/ul
response 9447

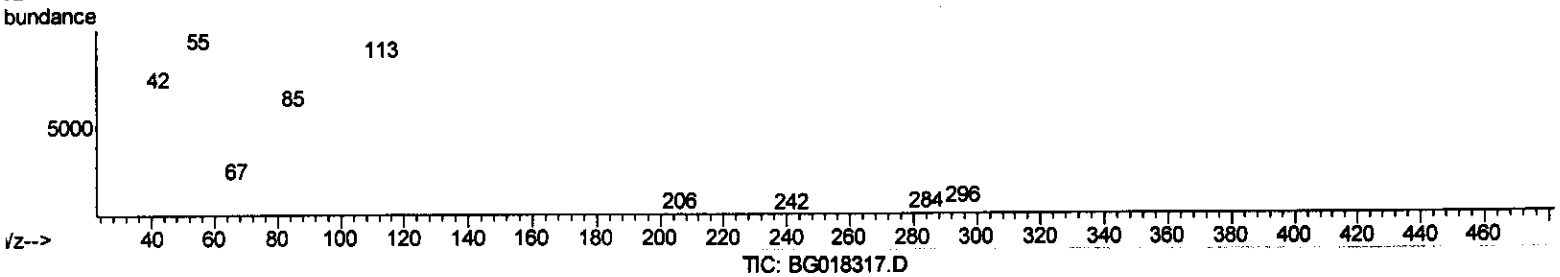
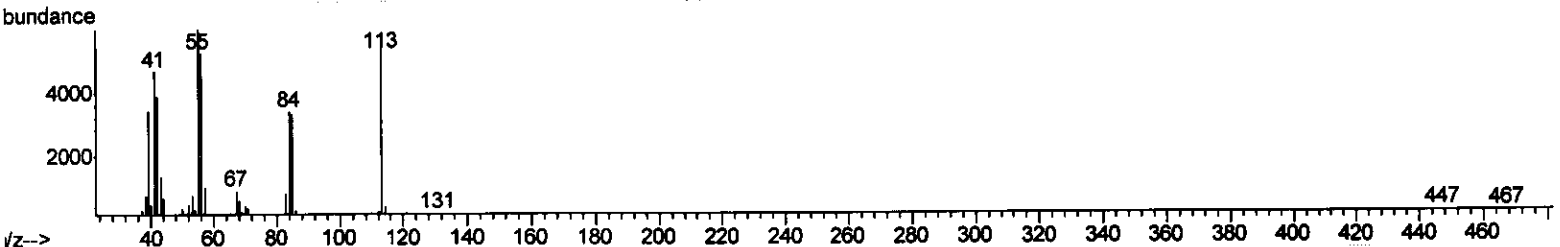
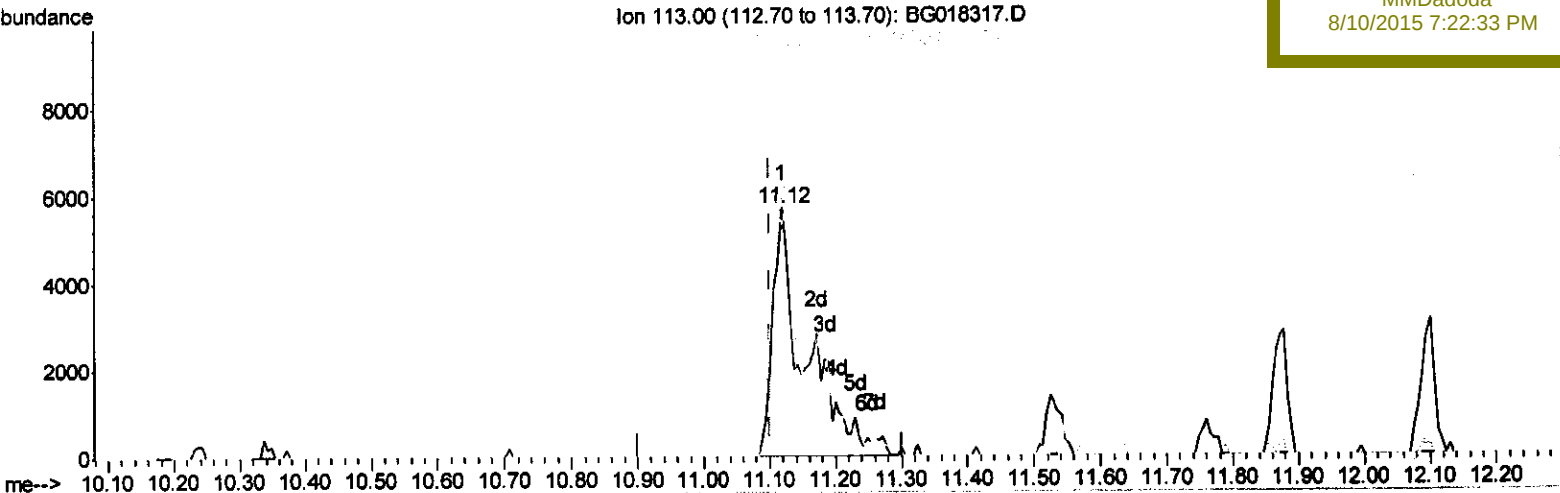
Ion	Exp%	Act%
113.00	100	100
55.00	120.80	104.56
56.00	112.30	91.02
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SST04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED
MMDadoda
8/10/2015 7:22:33 PM



(32) Caprolactam
11.119min (+0.018) 38.25ng/ul m > U.M
response 19136
0811115

Ion	Exp%	Act%
113.00	100	100
55.00	120.80	104.56
56.00	112.30	91.02
0.00	0.00	0.00

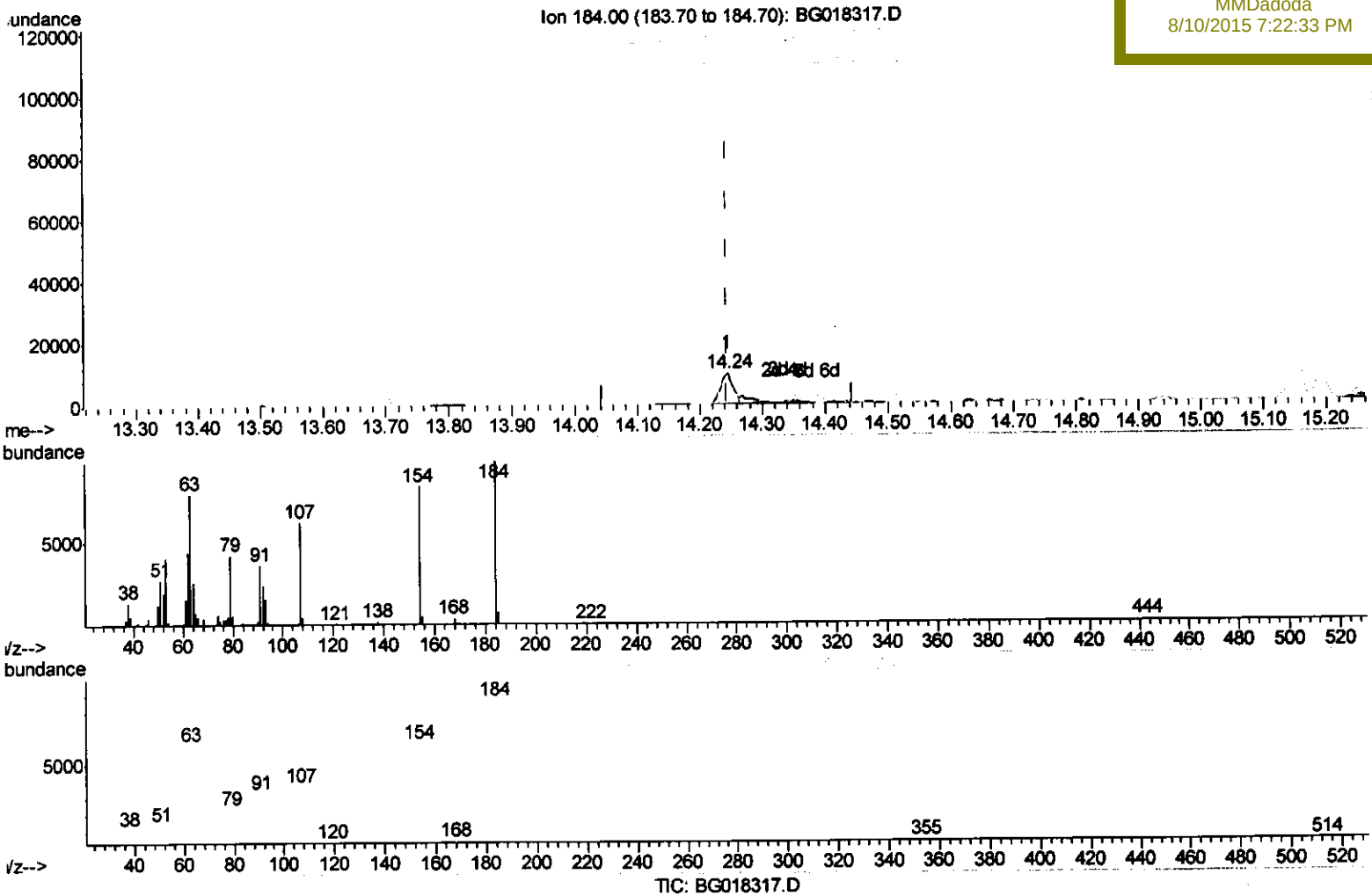
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED

MMDadoda
8/10/2015 7:22:33 PM



(51) 2,4-Dinitrophenol

14.244min (+0.000) 33.30ng/ui

response 13471

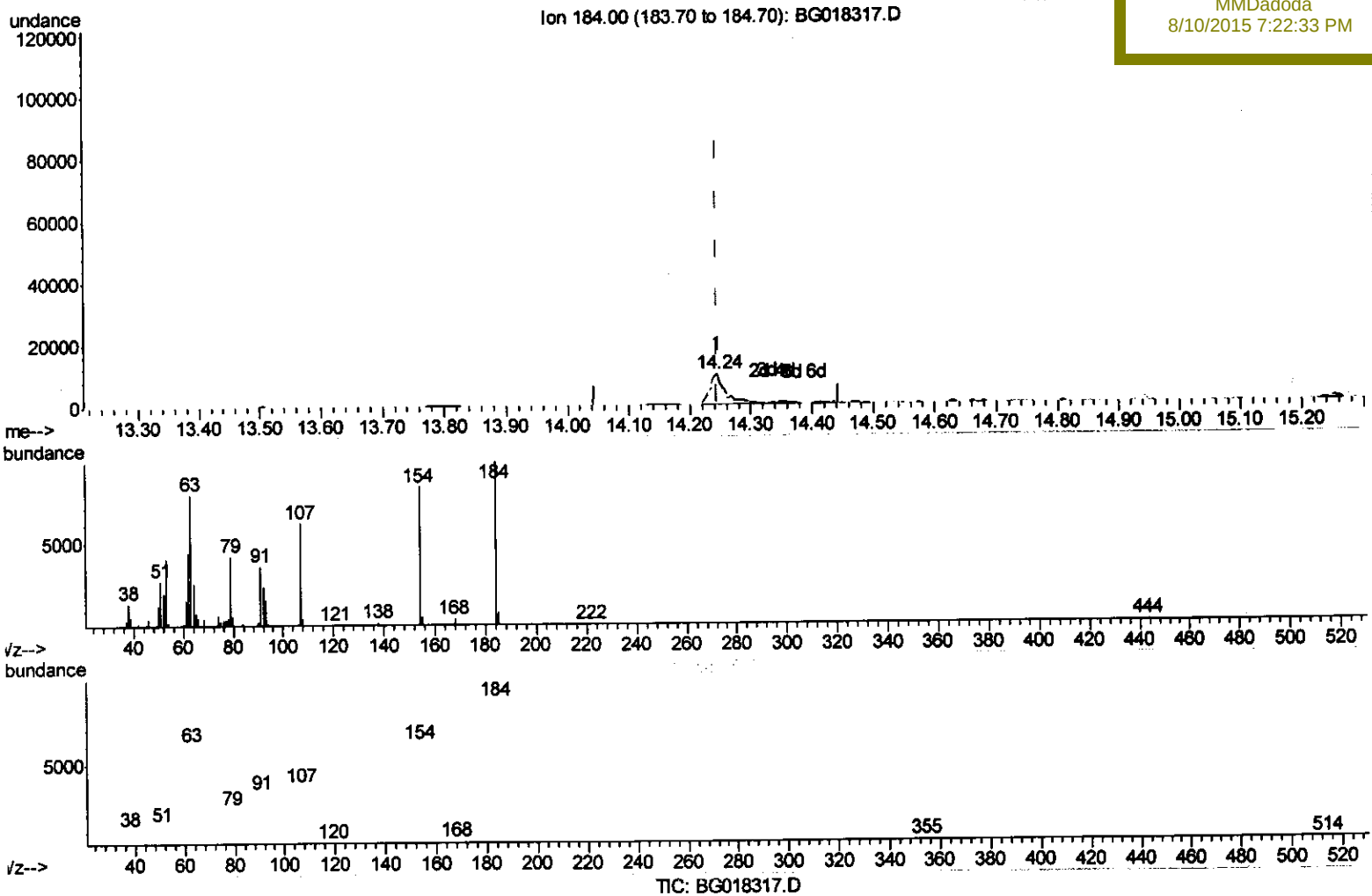
Ion	Exp%	Act%
184.00	100	100
63.00	61.40	80.80#
154.00	62.00	85.14#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED
MMDadoda
8/10/2015 7:22:33 PM



(51) 2,4-Dinitrophenol

14.244min (+0.000) 41.01ng/ul m } U.M
08111115

response 16588

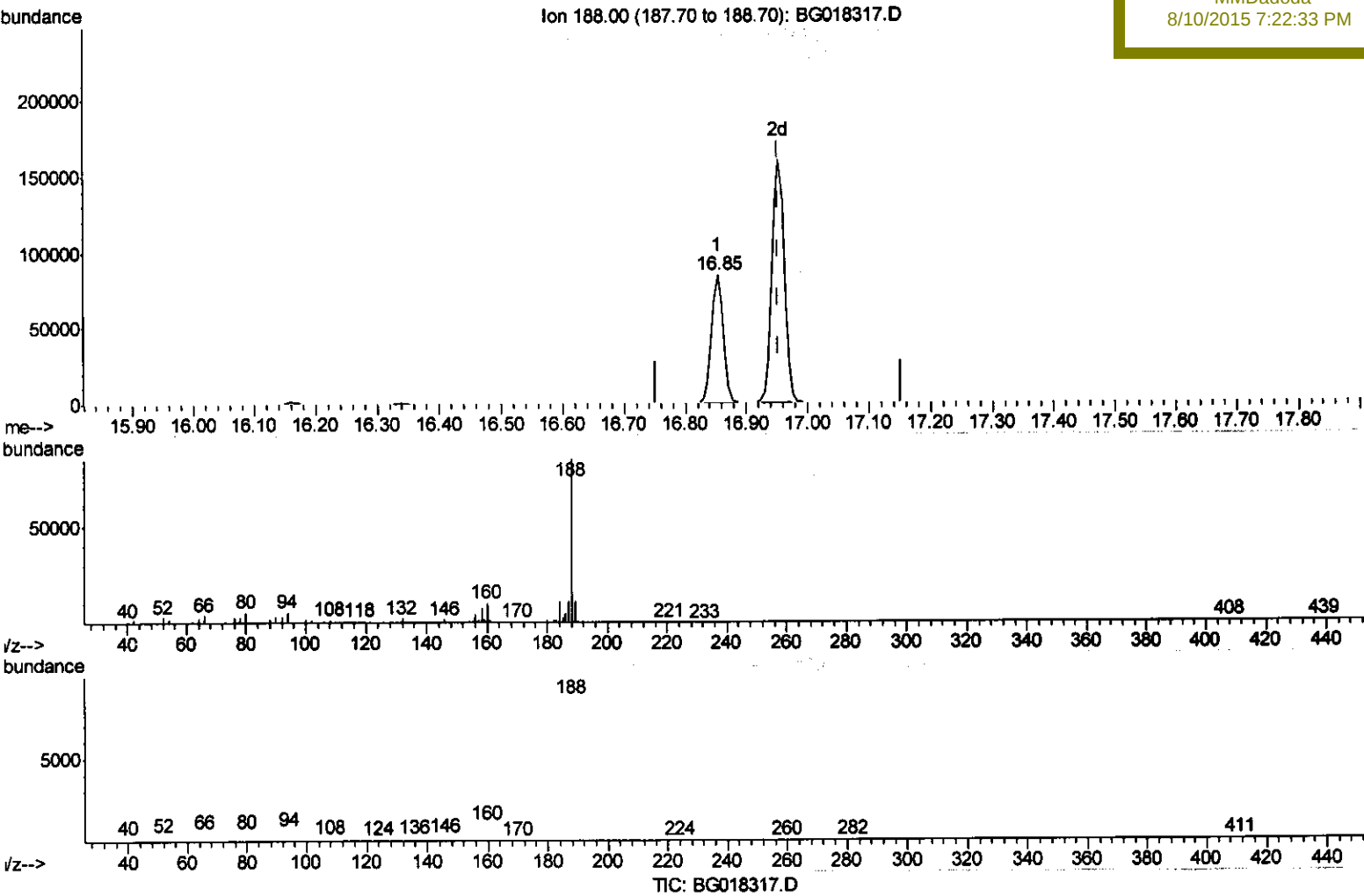
Ion	Exp%	Act%
184.00	100	100
63.00	61.40	80.80#
154.00	62.00	85.14#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED
MMDadoda
8/10/2015 7:22:33 PM



(71) Anthracene-d10 (S)

16.853min (-0.100) 18.53ng/ul

response 111025

Ion	Exp%	Act%
188.00	100	100
94.00	6.70	6.65
80.00	5.40	6.46
0.00	0.00	0.00

Instrument :
BNA_G
ClientSampleId :
SSTD04075

MMDadoda
8/10/2015 7:22:33 PM

Ion 188.00 (187.70 to 188.70): BG018317.D

abundance

200000

150000

100000

50000

0

me--> 16.00 16.10 16.20 16.30 16.40 16.50 16.60 16.70 16.80 16.90 17.00 17.10 17.20 17.30 17.40 17.50 17.60 17.70 17.80 17.90

abundance

100000

5000

42 52 66 80 94 104 114 132 146 160 170 188 222 333 358 409

188

42 52 66 80 94 108 124 136 146 160 170 224 260 282 411

TIC: BG018317.D

16.953min (+0.000) 37.31ng/ul m

Ion	Exp%	Act%
188.00	100	100
94.00	6.70	6.97
80.00	5.40	6.78#
0.00	0.00	0.00

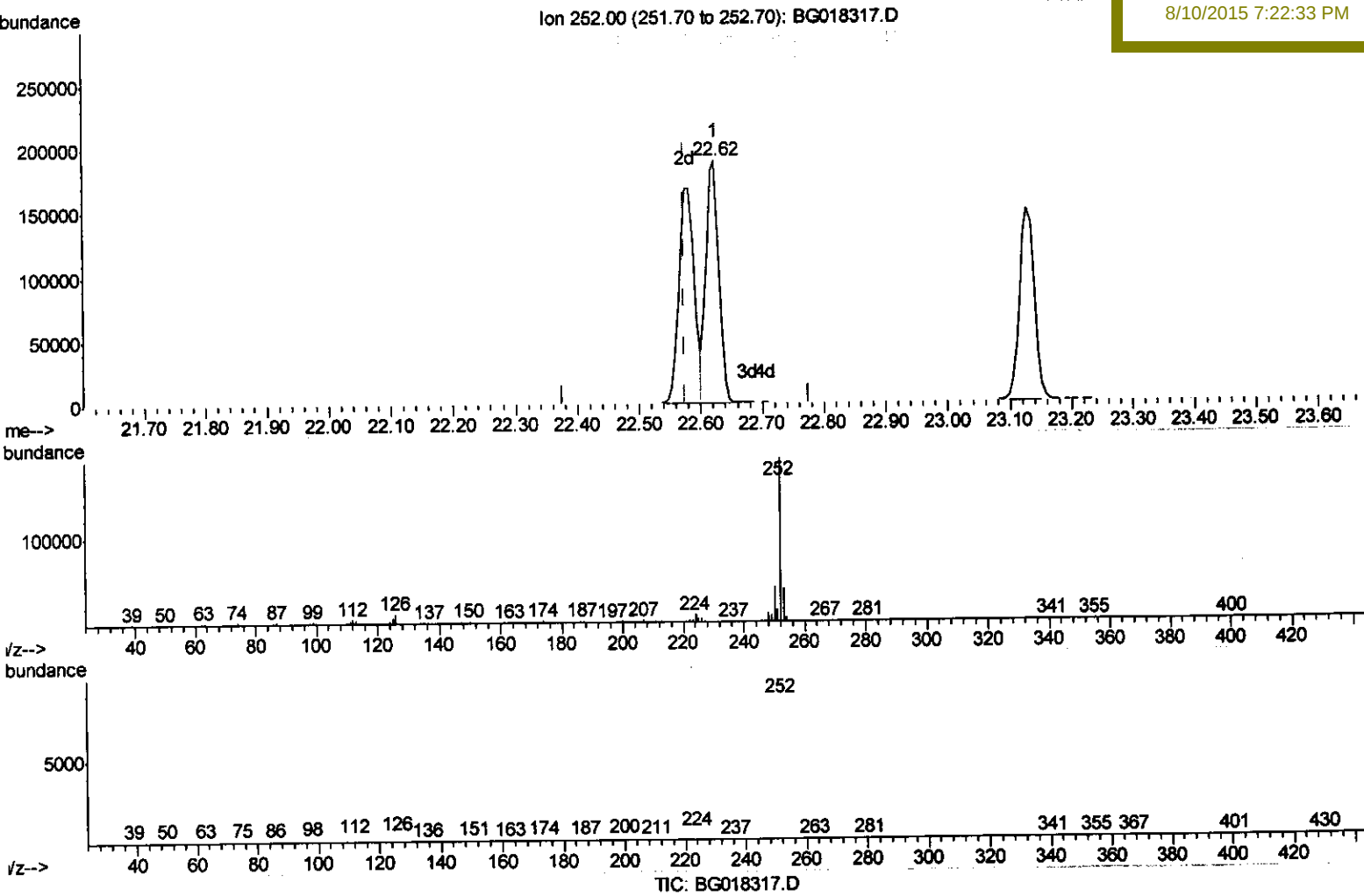
U.M
0811115

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SST04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED
MMDadoda
8/10/2015 7:22:33 PM



(88) Benzo(b)fluoranthene

22.623min (+0.047) 35.73ng/ul

response 272035

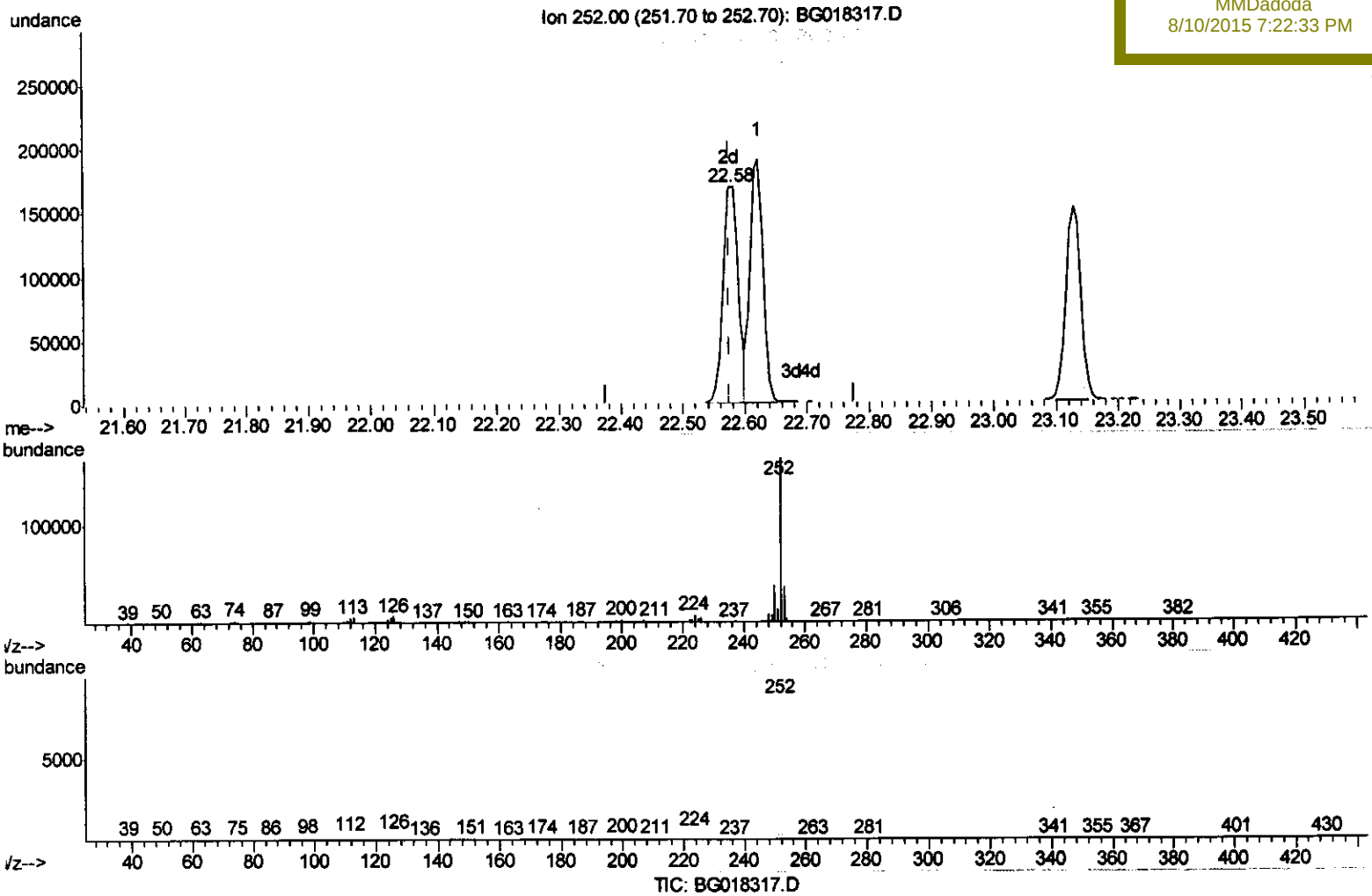
Ion	Exp%	Act%
252.00	100	100
253.00	20.30	20.85
125.00	3.00	3.21
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED
MMDadoda
8/10/2015 7:22:33 PM



(88) Benzo(b)fluoranthene

22.576min (+0.000) 38.08ng/ul m

response 289906

Ion Exp% Act%

252.00 100 100

253.00 20.30 21.21

125.00 3.00 3.61#

0.00 0.00 0.00

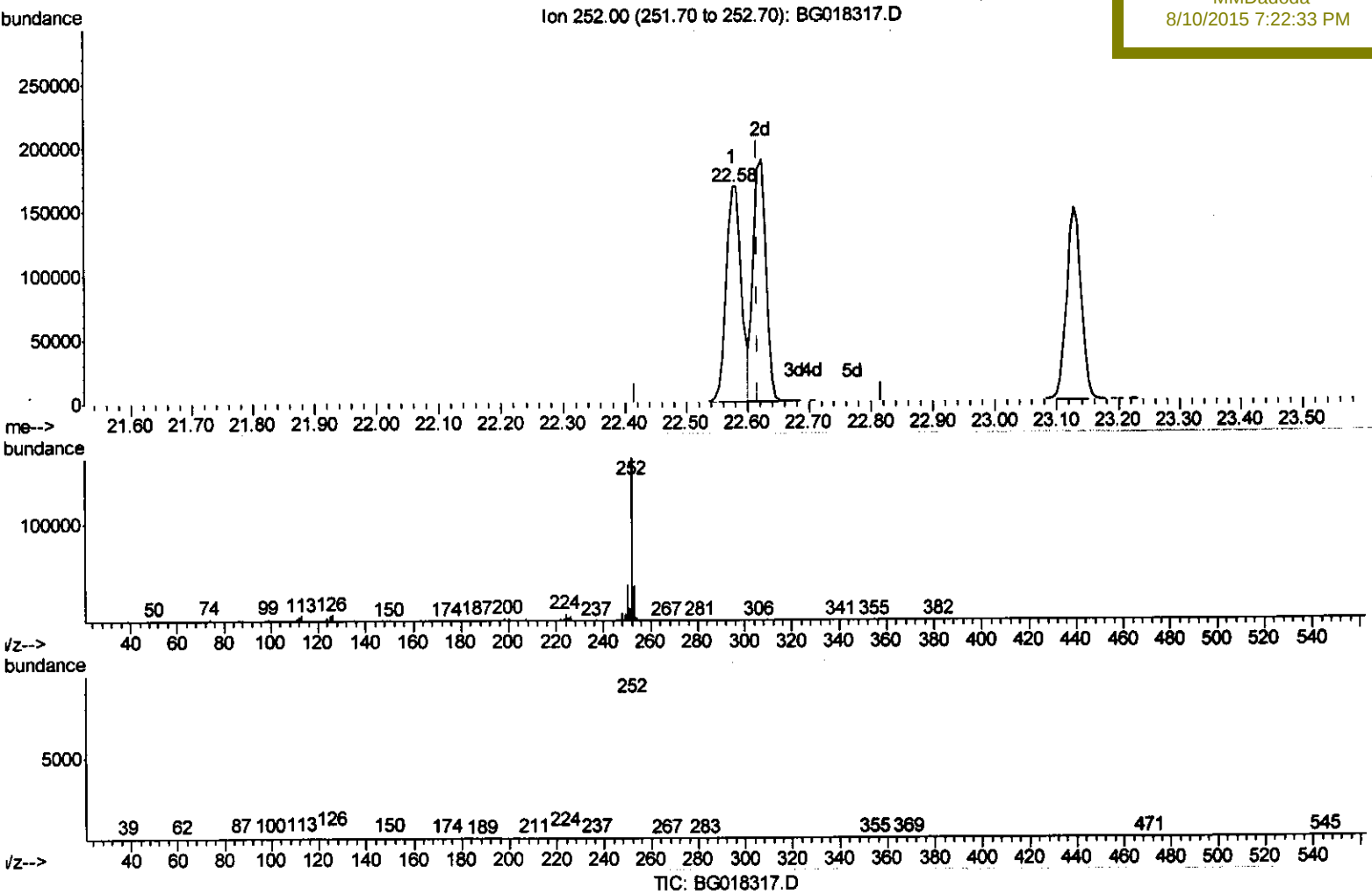
U.M
0811115

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED
MMDadoda
8/10/2015 7:22:33 PM



(89) Benzo(k)fluoranthene

22.576min (-0.041) 39.87ng/ul

response 289961

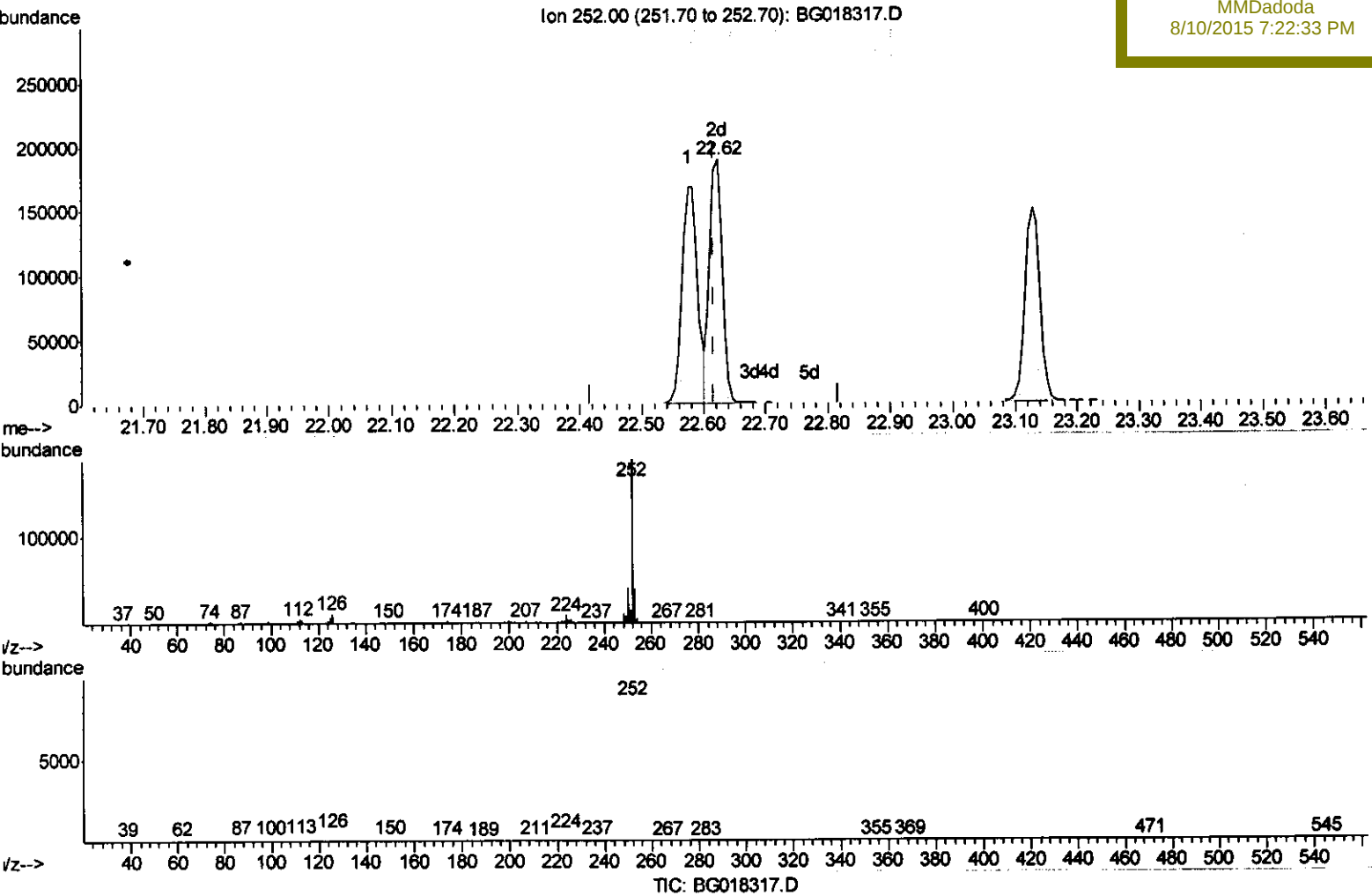
Ion	Exp%	Act%
252.00	100	100
253.00	21.70	21.21
125.00	4.10	3.61
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SST04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED
MMDadoda
8/10/2015 7:22:33 PM



(89) Benzo(k)fluoranthene

22.623min (+0.006) 37.38ng/ui m

response 271837

Ion	Exp%	Act%
252.00	100	100
253.00	21.70	20.85
125.00	4.10	3.21#
0.00	0.00	0.00

U.M
08/11/15

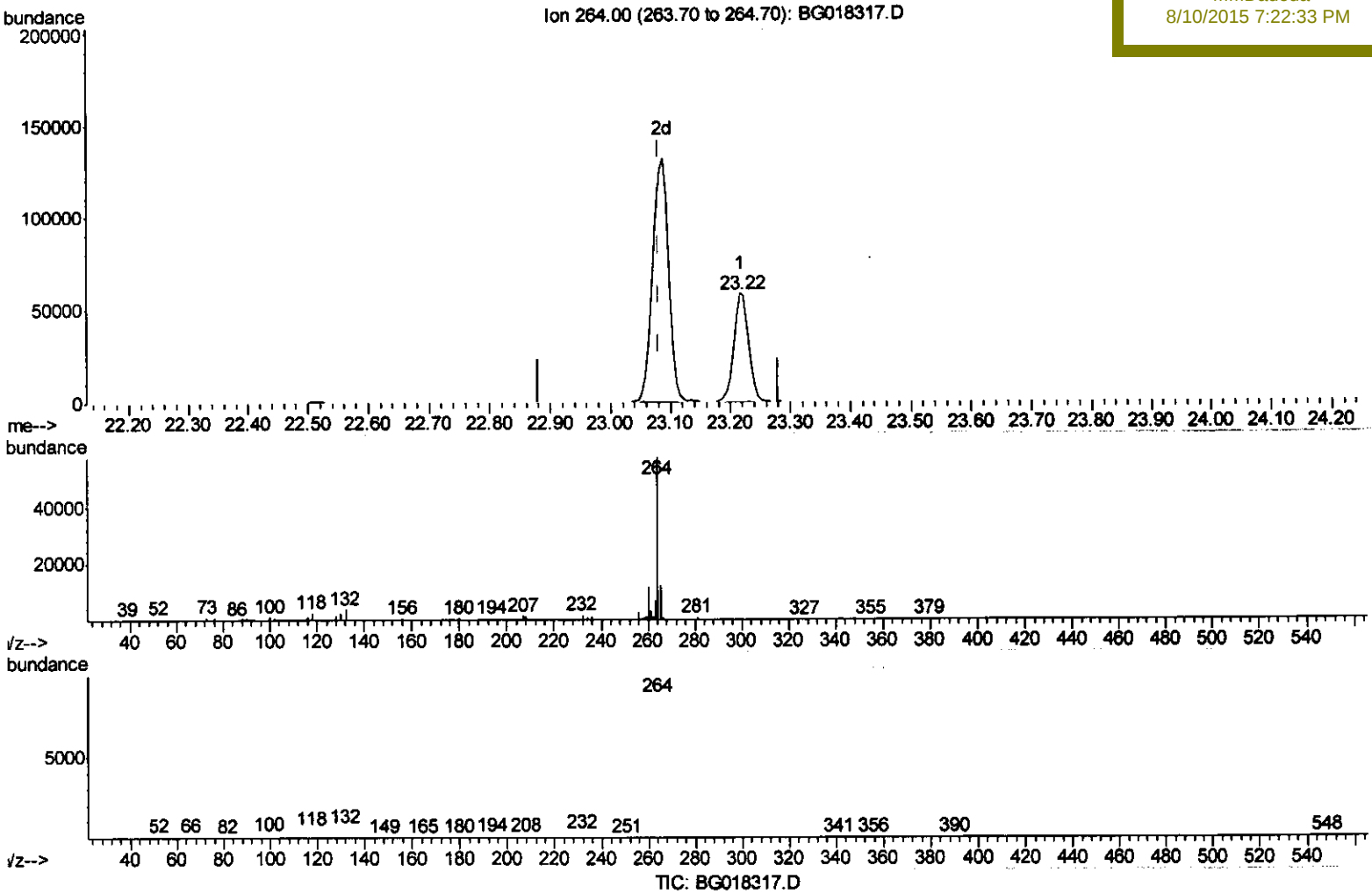
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED

MMDadoda
8/10/2015 7:22:33 PM



(90) Benzo(a)pyrene-d12 (S)
23.216min (+0.135) 15.28ng/ul

response 99327

Ion	Exp%	Act%
264.00	100	100
132.00	6.80	6.88
118.00	5.30	4.69
0.00	0.00	0.00

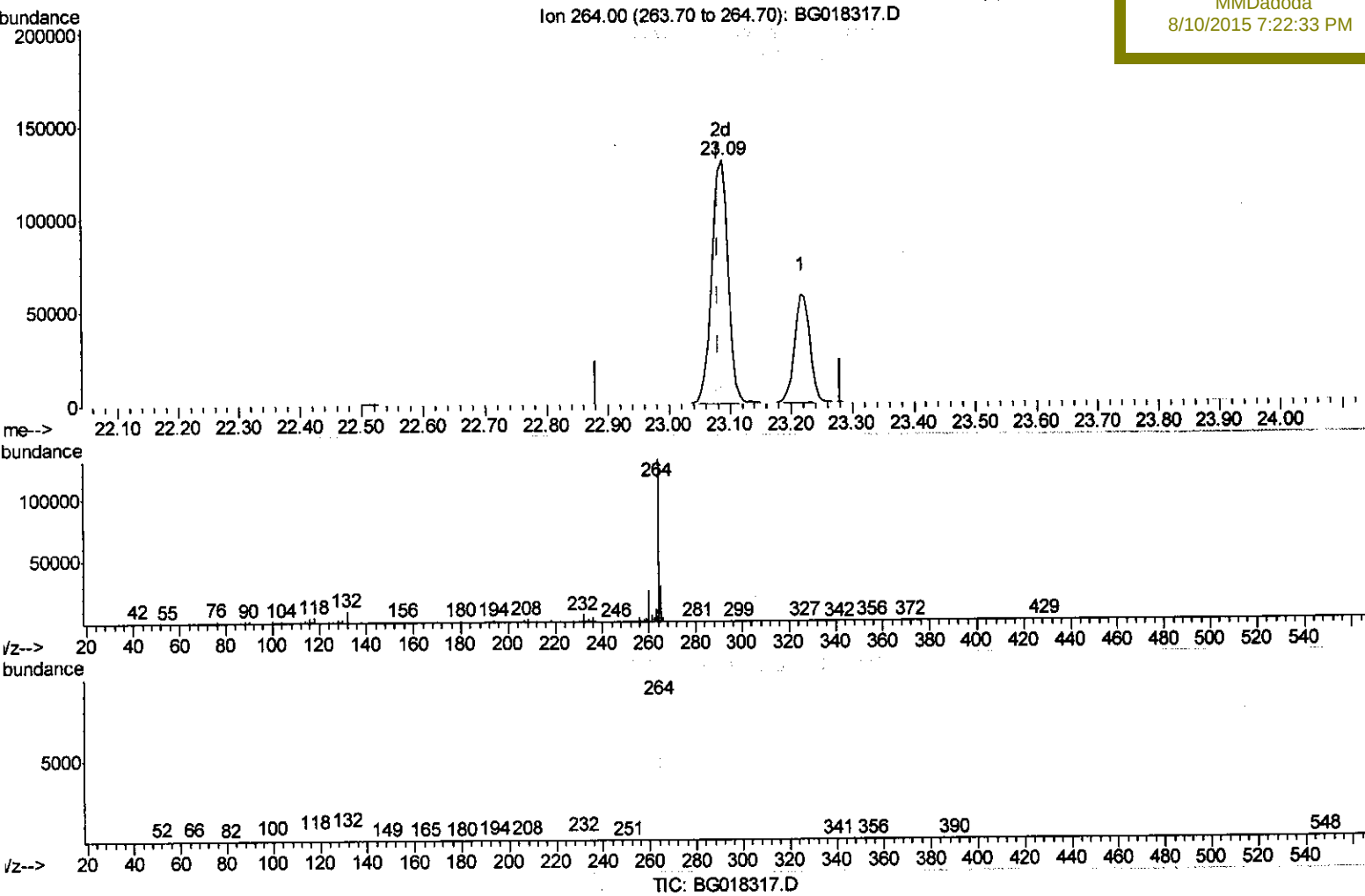
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 07:51:30 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED

MMDadoda
8/10/2015 7:22:33 PM



(90) Benzo(a)pyrene-d12 (S)

23.087min (+0.006) 37.03ng/ul m

response 240691

Ion	Exp%	Act%
264.00	100	100
132.00	6.80	7.08
118.00	5.30	3.34#
0.00	0.00	0.00

> U.M
08/11/15

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
 Data File : BG018317.D
 Acq On : 8 Aug 2015 6:40
 Operator : UM/TP
 Sample : SSTD04075
 Disc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 Client Sample Id :
 SSTD04075

Quant Time: Aug 08 08:06:52 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Aug 08 07:47:13 2015
 Response via : Initial Calibration

Manual Integrations
 APPROVED

MMDadoda
 8/10/2015 7:22:33 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	14218	20.00	ng/ul	0.00
18) Naphthalene-d8	10.18	136	62262	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.09	164	42928	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	111025	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	111594	20.00	ng/ul	0.00
86) Perylene-d12	23.22	264	100487	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.87	96	4065	15.27	ng/uL	0.00
5) Phenol-d5	6.61	99	49901	38.39	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.75	67	28524	38.29	ng/ul	0.00
9) 2-Chlorophenol-d4	6.95	132	39271	36.07	ng/ul	0.00
13) 4-Methylphenol-d8	8.15	113	42391	37.93	ng/ul	0.00
19) Nitrobenzene-d5	8.56	128	19758	34.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.29	143	23449	35.90	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.83	165	44347	36.80	ng/ul	0.00
29) 4-Chloroaniline-d4	10.34	131	56859	42.52	ng/ul	0.00
44) Dimethylphthalate-d6	13.52	166	155006	35.82	ng/ul	0.00
47) Acenaphthylene-d8	13.78	160	172399	36.04	ng/ul	0.00
52) 4-Nitrophenol-d4	14.36	143	21352	38.11	ng/ul	0.00
58) Fluorene-d10	15.10	176	139629	36.82	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.25	200	32774	39.82	ng/ul	0.00
71) Anthracene-d10	16.95	188	223482m	37.31	ng/ul	0.00
79) Pyrene-d10	19.27	212	233994	36.24	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.09	264	240691m	37.03	ng/ul	0.00

U.M
 08/11/15
 U.M
 08/11/15

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.90	88	4069	12.32	ng/uL#	79
4) Benzaldehyde	6.55	77	38517	45.26	ng/ul	91
6) Phenol	6.64	94	52349	39.55	ng/ul	95
8) Bis(2-Chloroethyl)ether	6.84	93	35160	36.09	ng/ul	97
10) 2-Chlorophenol	6.98	128	38105	35.60	ng/ul#	93
11) 2-Methylphenol	7.88	108	39646	37.68	ng/ul#	75
12) 2,2'-oxybis(1-Chloropropan	7.95	45	34638	39.73	ng/ul#	86
14) Acetophenone	8.23	105	69564	37.85	ng/ul	96
15) N-Nitroso-di-n-propylamine	8.22	70	41676	36.84	ng/ul#	81
16) 4-Methylphenol	8.20	108	46231	39.47	ng/ul	89
17) Hexachloroethane	8.46	117	20141	36.26	ng/ul	97
20) Nitrobenzene	8.60	77	63795	40.03	ng/ul	98
21) Isophorone	9.13	82	111459	37.58	ng/ul	97
23) 2-Nitrophenol	9.31	139	24530	36.17	ng/ul#	83
24) 2,4-Dimethylphenol	9.40	107	60987	37.36	ng/ul	98
25) Bis(2-Chloroethoxy)methane	9.62	93	53950	38.33	ng/ul	92
27) 2,4-Dichlorophenol	9.86	162	45050	39.40	ng/ul	93
28) Naphthalene	10.24	128	129663	36.29	ng/ul	97
30) 4-Chloroaniline	10.37	127	54065	42.18	ng/ul	99
31) Hexachlorobutadiene	10.52	225	35517	35.94	ng/ul#	76
32) Caprolactam	11.12	113	19136m	38.25	ng/ul	
33) 4-Chloro-3-methylphenol	11.52	107	56266	36.78	ng/ul#	83
34) 2-Methylnaphthalene	11.87	142	103966	36.85	ng/ul	93

U.M
 08/11/15

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
 Data File : BG018317.D
 Acq On : 8 Aug 2015 6:40
 Operator : UM/TP
 Sample : SSTD04075
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 08:06:52 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Aug 08 07:47:13 2015
 Response via : Initial Calibration

Instrument :
 BNA_G
 ClientSampleId :
 SSTD04075

Manual Integrations
 APPROVED

MMDadoda
 8/10/2015 7:22:33 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.09	142	102086	36.45	ng/ul#	88
37) 1,2,4,5-Tetrachlorobenzene	12.26	216	61004	38.40	ng/ul#	80
38) Hexachlorocyclopentadiene	12.22	237	21781	37.10	ng/ul#	93
39) 2,4,6-Trichlorophenol	12.52	196	39352	37.91	ng/ul	94
40) 2,4,5-Trichlorophenol	12.59	196	40716	37.35	ng/ul	95
41) 1,1'-Biphenyl	12.92	154	141858	37.16	ng/ul#	94
42) 2-Chloronaphthalene	12.95	162	109693	37.54	ng/ul	98
43) 2-Nitroaniline	13.18	65	46268	42.80	ng/ul	93
45) Dimethylphthalate	13.56	163	156748	37.02	ng/ul#	97
46) 2,6-Dinitrotoluene	13.69	165	34677	38.00	ng/ul	91
48) Acenaphthylene	13.81	152	175793	36.78	ng/ul	97
49) 3-Nitroaniline	14.02	138	30980	39.27	ng/ul#	92
50) Acenaphthene	14.16	153	115629	35.94	ng/ul	98
51) 2,4-Dinitrophenol	14.24	184	16588m}	41.01	ng/ul	
53) 4-Nitrophenol	14.37	109	35203	39.58	ng/ul	96
54) Dibenzofuran	14.50	168	171130	36.12	ng/ul	89
55) 2,4-Dinitrotoluene	14.49	165	50766	37.30	ng/ul	96
56) 2,3,4,6-Tetrachlorophenol	14.74	232	34144	34.57	ng/ul#	85
57) Diethylphthalate	14.94	149	171521	36.88	ng/ul	98
59) Fluorene	15.16	166	152841	37.15	ng/ul	99
60) 4-Chlorophenyl-phenylether	15.15	204	85166	36.17	ng/ul	96
61) 4-Nitroaniline	15.20	138	31708	39.40	ng/ul	92
64) 4,6-Dinitro-2-methylphenol	15.26	198	31501	37.75	ng/ul#	87
65) N-Nitrosodiphenylamine	15.37	169	130851	36.20	ng/ul	92
66) 4-Bromophenyl-phenylether	16.05	248	58400	36.13	ng/ul	98
67) Hexachlorobenzene	16.16	284	70418	35.85	ng/ul	92
68) Atrazine	16.34	200	60298	38.01	ng/ul	92
69) Pentachlorophenol	16.52	266	19351	34.45	ng/ul#	85
70) Phenanthrene	16.89	178	250348	36.05	ng/ul	98
72) Anthracene	16.99	178	247222	35.39	ng/ul	97
73) 1,2,3,4-Tetrachlorobenzene	12.88	216	63872	38.61	ng/uL	95
74) Pentachlorobenzene	14.42	250	72577	38.28	ng/uL	93
75) Carbazole	17.27	167	212672	39.53	ng/ul	97
76) Di-n-butylphthalate	17.83	149	293271	36.02	ng/ul	99
77) Fluoranthene	18.93	202	297890	39.63	ng/ul	99
80) Pyrene	19.30	202	294898	36.26	ng/ul#	96
81) Butylbenzylphthalate	20.21	149	131131	36.89	ng/ul	92
82) 3,3'-Dichlorobenzidine	21.00	252	115607	40.01	ng/ul	92
83) Benzo(a)anthracene	21.05	228	276494	36.40	ng/ul	99
84) Bis(2-ethylhexyl)phthalate	21.00	149	174582	36.28	ng/ul	99
85) Chrysene	21.11	228	256141	36.45	ng/ul	97
87) Di-n-octyl phthalate	21.85	149	267337	38.66	ng/ul	100
88) Benzo(b)fluoranthene	22.58	252	289906m}	38.08	ng/ul	
89) Benzo(k)fluoranthene	22.62	252	271837m}	37.38	ng/ul	
91) Benzo(a)pyrene	23.13	252	252780	36.87	ng/ul	98
92) Indeno(1,2,3-cd)pyrene	25.39	276	335092	40.01	ng/ul	93
93) Dibenzo(a,h)anthracene	25.40	278	286115	39.21	ng/ul#	99
94) Benzo(g,h,i)perylene	26.05	276	263375	38.61	ng/ul#	92

U.M
 08/11/15

U.M
 08/11/15

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080815\
Data File : BG018317.D
Acq On : 8 Aug 2015 6:40
Operator : UM/TP
Sample : SSTD04075
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 08 08:06:52 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080815.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 08 07:47:13 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
SSTD04075

Manual Integrations
APPROVED

MMDadoda
8/10/2015 7:22:33 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

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