

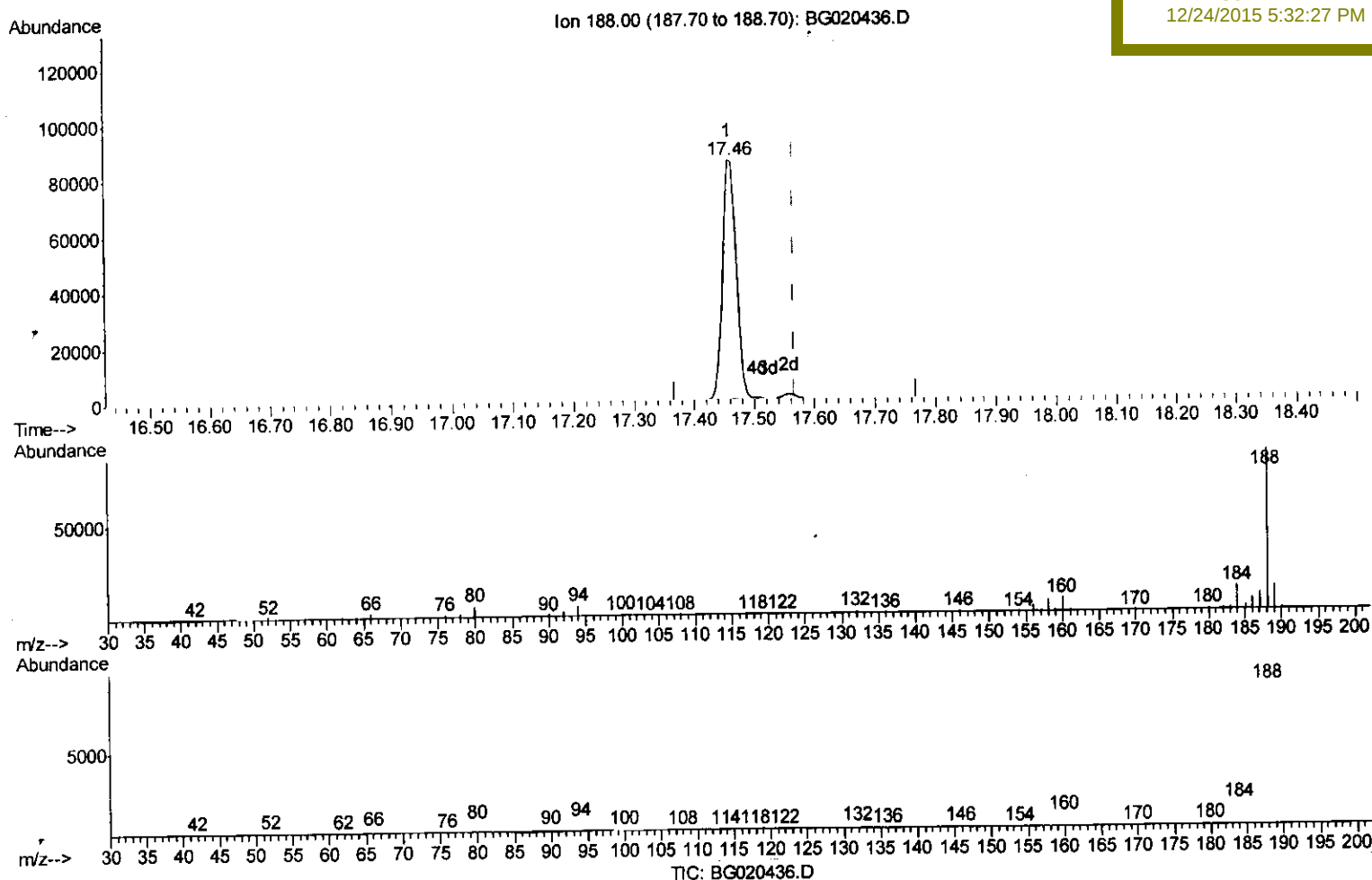
Data Path : Z:\HPCHEM1\BNA_G\Data\BG122215\
Data File : BG020436.D
Acq On : 23 Dec 2015 10:13
Operator : UM/SJ
Sample : G4768-13DL 125X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Dec 23 11:21:21 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 22 05:02:37 2015
Response via : Initial Calibration

Instrument :
BNA_G
Client Sampled :
D9N37DL

Manual Integrations
APPROVED

Sohil
12/24/2015 5:32:27 PM



(71) Anthracene-d10 (S)

17.458min (-0.110) 21.70ng/ul

response 132573

Ion	Exp%	Act%
188.00	100	100
94.00	7.70	6.42
80.00	7.50	7.29
0.00	0.00	0.00

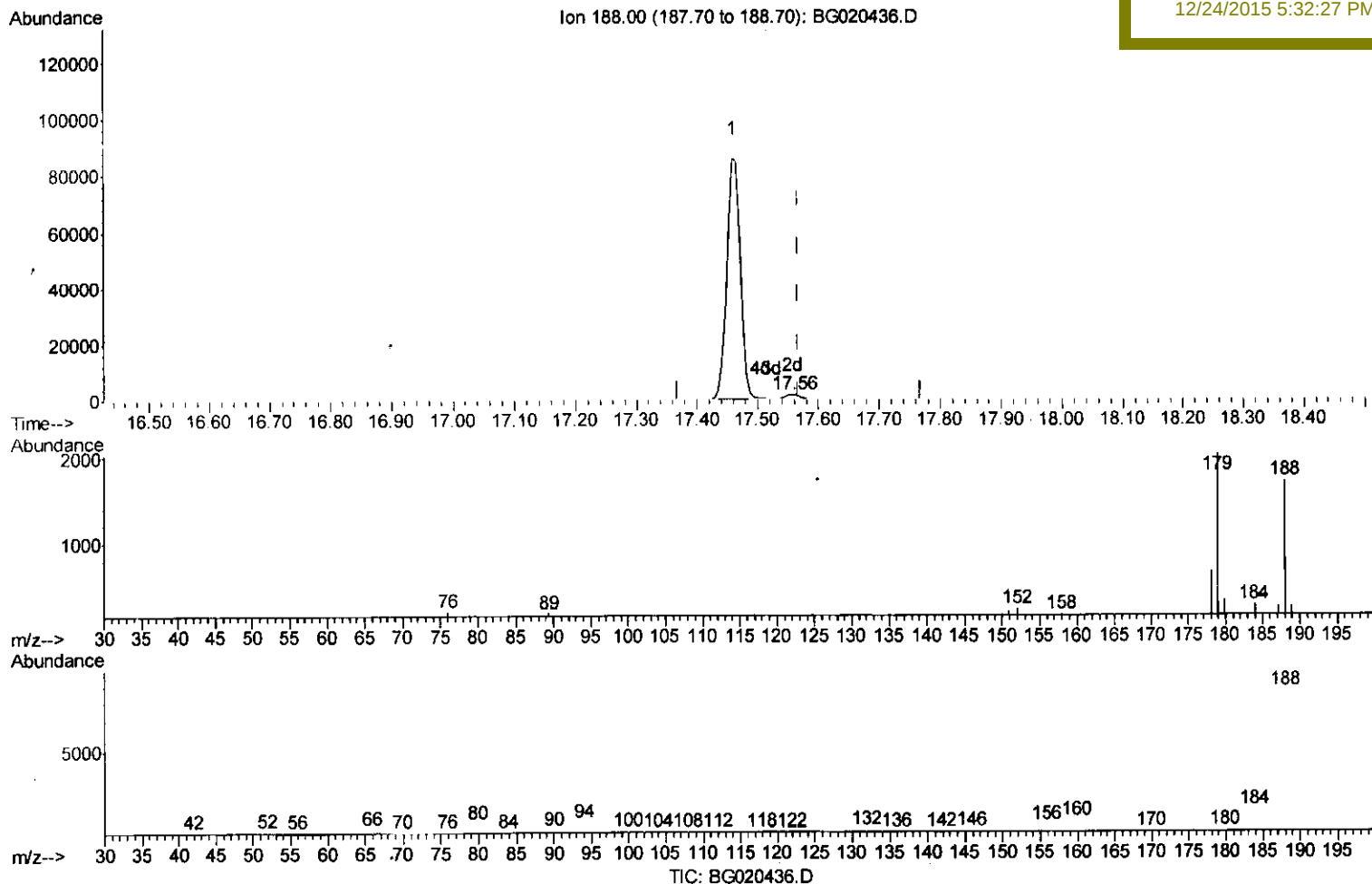
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(71) Anthracene-d10 (S)

17.558min (-0.010) 0.46ng/ul m

response 2839

Ion Exp% Act%

188.00 100 100

94.00 7.70 0.00#

80.00 7.50 0.00#

0.00 0.00 0.00

> U.M
28/12/15

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Quant Time: Dec 23 11:23:40 2015
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	12463	20.00	ng/ul	0.00
18) Naphthalene-d8	10.90	136	58157	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.71	164	46887	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.46	188	132573	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	171179	20.00	ng/ul	0.00
86) Perylene-d12	25.00	264	164929	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/ul	
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	8.62	113	178	0.19	ng/ul	-0.19
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	14.12	166	1473	0.39	ng/ul	0.00
47) Acenaphthylene-d8	14.41	160	1381	0.31	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.71	176	1442	0.42	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.56	188	2839m	0.46	ng/ul	0.00
79) Pyrene-d10	19.84	212	2577	0.32	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.77	264	3062	0.37	ng/ul	-0.03

Target Compounds

						Qvalue
28) Naphthalene	10.95	128	180862	59.55	ng/ul	99
34) 2-Methylnaphthalene	12.55	142	16316	7.03	ng/ul	100
41) 1,1'-Biphenyl	13.55	154	5508	1.56	ng/ul#	85
50) Acenaphthene	14.78	153	24588	7.80	ng/ul	97
54) Dibenzofuran	15.11	168	20275	4.32	ng/ul	97
59) Fluorene	15.76	166	13828	3.67	ng/ul	94
70) Phenanthrene	17.50	178	30575	4.22	ng/ul	98
75) Carbazole	17.86	167	35954	5.82	ng/ul	96

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

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