

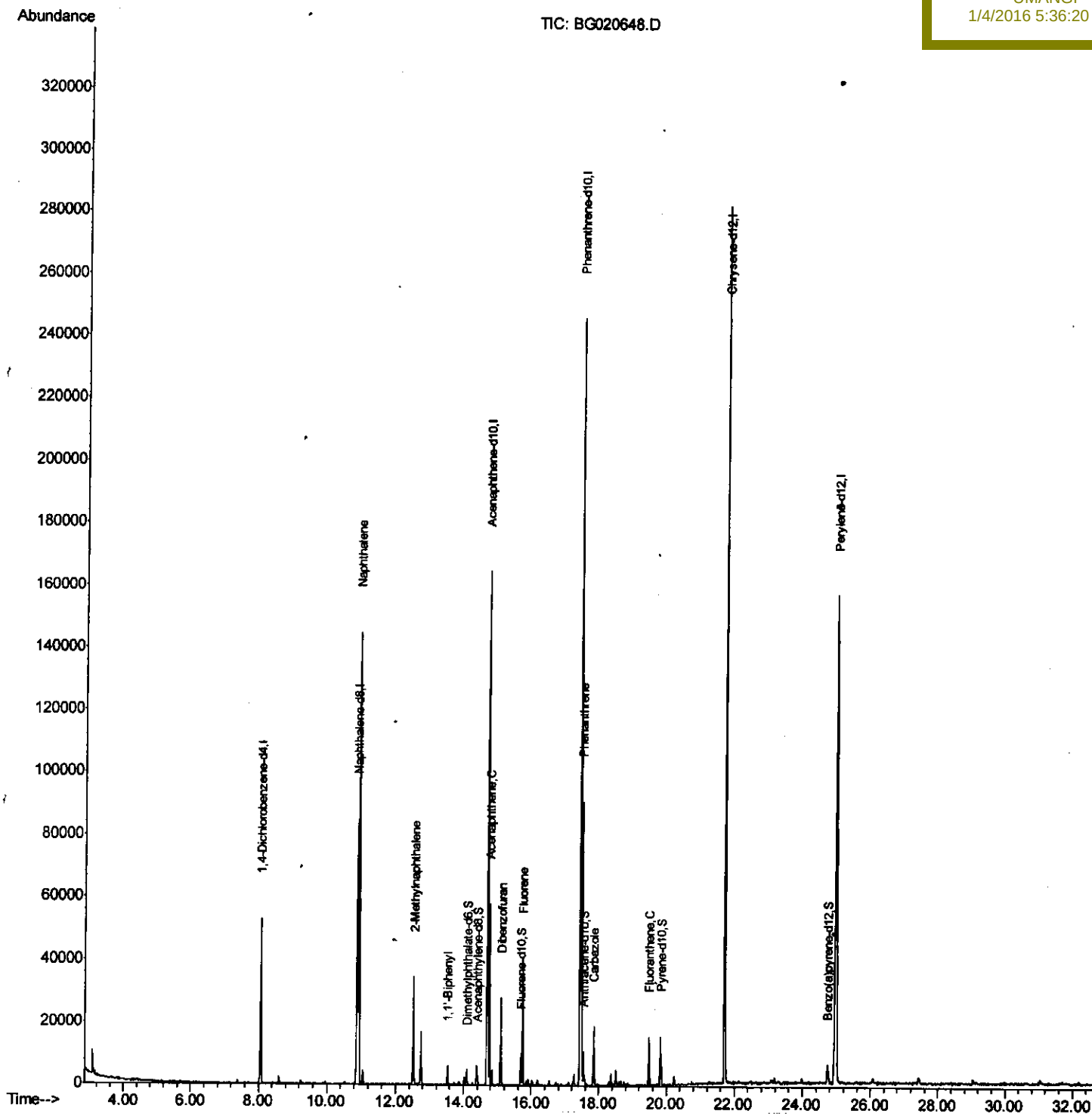
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG123115\
Data File : BG020648.D
Acq On : 31 Dec 2015 19:13
Operator : UM/SJ
Sample : G4847-07DL 30X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Quant Time: Jan 01 04:33:29 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG123015.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 01 03:15:46 2016
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
D9N71DL

Manual Integrations
APPROVED

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Quantitation Report (Qedit)

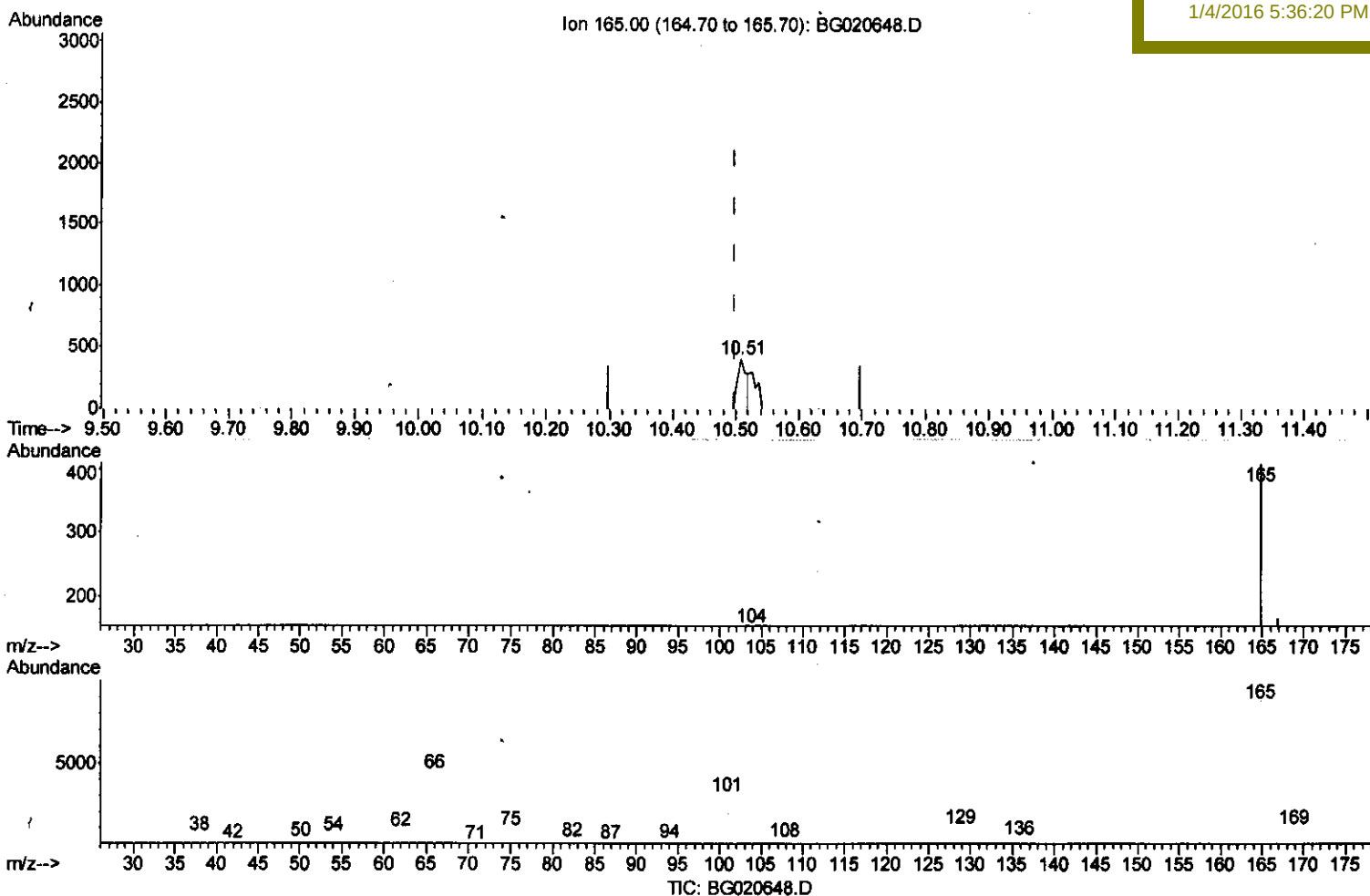
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Quant Time: Jan 01 03:21:04 2016
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(26) 2,4-Dichlorophenol-d3 (S)

10.508min (+0.009) 0.34ng/ul

response 429

Ion	Exp%	Act%
165.00	100	100
167.00	69.50	40.69#
101.00	31.20	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

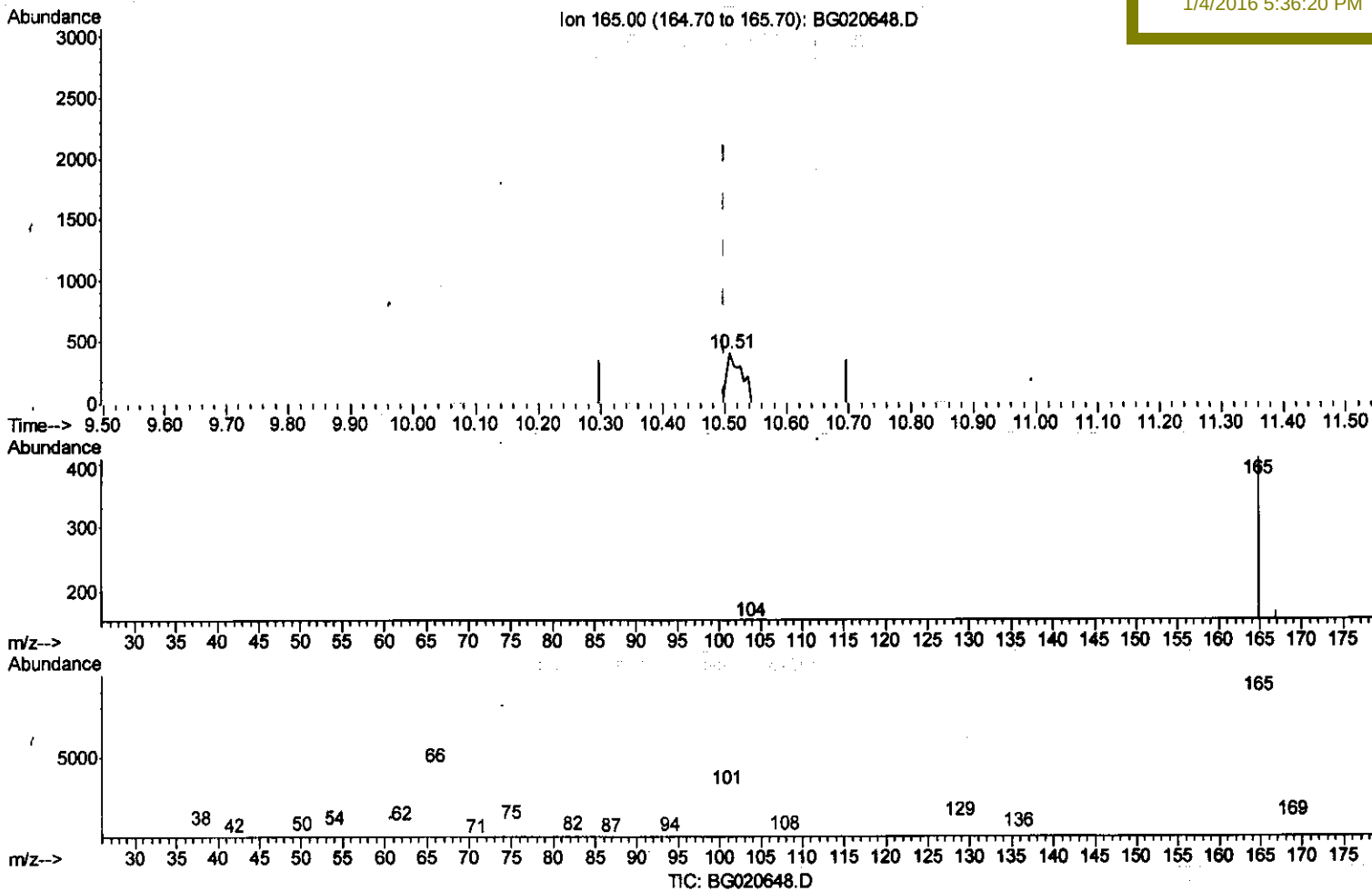
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(26) 2,4-Dichlorophenol-d3 (S)

10.508min (+0.009) 0.52ng/ul m U, M

response 671

1/2/16

Ion	Exp%	Act%
165.00	100	100
167.00	69.50	40.69#
101.00	31.20	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

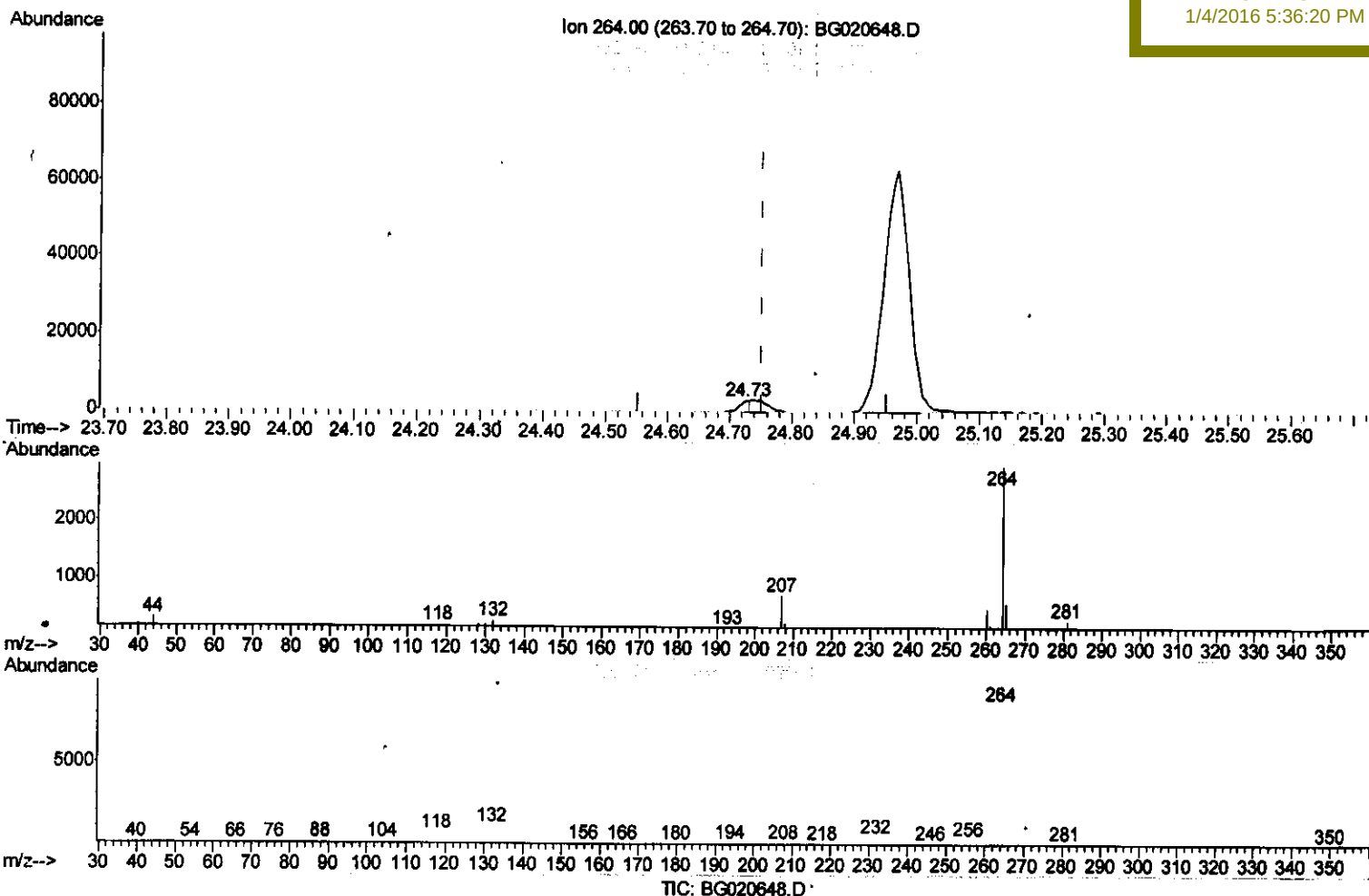
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Manual Integrations
 APPROVED

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

24.726min (-0.026) 0.43ng/ul

response 3774

Ion	Exp%	Act%
264.00	100	100
132.00	10.70	8.33%
118.00	6.10	6.03
0.00	0.00	0.00

Quantitation Report (Qedit)

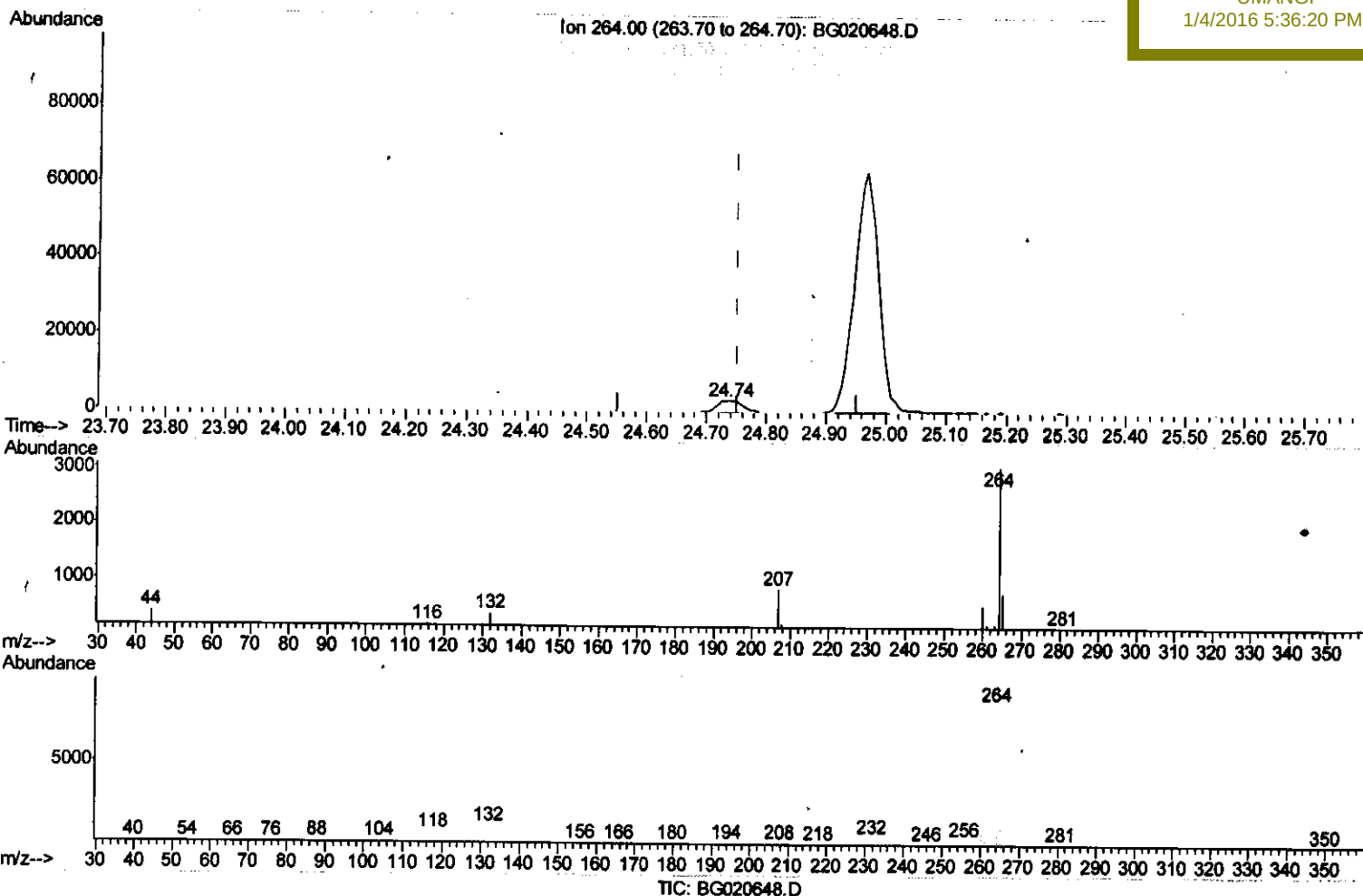
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG123115\
 Data File : BG020648.D
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 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
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Manual Integrations
 APPROVED

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

24.738min (-0.015) 1.09ng/ul m U.M

response 9563

1/2/16

Ion	Exp%	Act%
264.00	100	100
132.00	10.70	12.34
118.00	6.10	5.55
0.00	0.00	0.00

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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	8.06	152	16342	20.00	ng/ul	0.00
18) Naphthalene-d8	10.87	136	75077	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.70	164	58424	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.44	188	154876	20.00	ng/ul	0.00
78) Chrysene-d12	21.71	240	184848	20.00	ng/ul	0.00
86) Perylene-d12	24.97	264	176636	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	0.00	99	0d	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.39	67	561	0.67	ng/ul	0.00
9) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	9.24	128	361	0.62	ng/ul	0.00
22) 2-Nitrophenol-d4	9.97	143	365	0.54	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	10.51	165	671m	0.52	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	14.09	166	5199	1.09	ng/ul	0.00
47) Acenaphthylene-d8	14.39	160	5712	1.05	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.68	176	5141	1.19	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.53	188	8615	1.19	ng/ul	0.00
79) Pyrene-d10	19.82	212	9241	1.11	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.74	264	9563m	1.09	ng/ul	-0.01

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Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
28) Naphthalene	10.93	128	126888	31.91	ng/ul	98
34) 2-Methylnaphthalene	12.53	142	18914	6.36	ng/ul	87
41) 1,1'-Biphenyl	13.53	154	4899	1.11	ng/ul#	98
50) Acenaphthene	14.76	153	25960	6.63	ng/ul	97
54) Dibenzofuran	15.09	168	20063	3.39	ng/ul	93
59) Fluorene	15.74	166	19727	4.15	ng/ul	99
70) Phenanthrene	17.48	178	59394	7.05	ng/ul	96
75) Carbazole	17.85	167	16489	2.27	ng/ul	96
77) Fluoranthene	19.49	202	12037	1.17	ng/ul#	90

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