

Quantitation Report (Qedit)

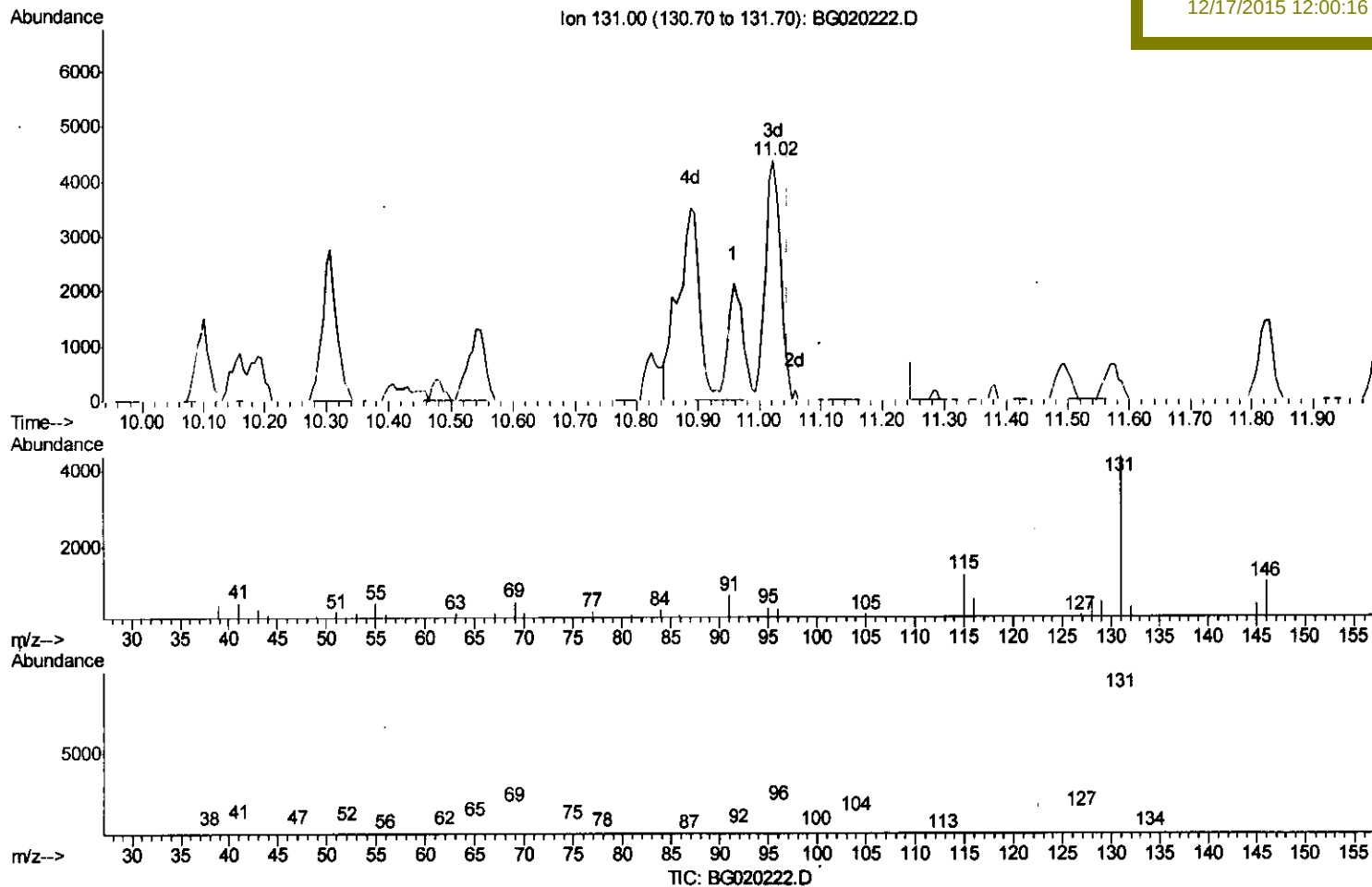
Data Path : Z:\HPCHEM1\BNA_G\Data\BG121715\
Data File : BG020222.D
Acq On : 16 Dec 2015 14:12
Operator : UM/SJ
Sample : G4786-17DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Dec 17 00:43:15 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 17 00:37:15 2015
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampled :
G9C90DL

Manual Integrations
APPROVED

MMdadoda
12/17/2015 12:00:16 PM



(29) 4-Chloroaniline-d4 (S)

11.023min (-0.023) 4.57ng/ul m

response 7426

Ion	Exp%	Act%
131.00	100	100
133.00	30.40	0.00#
69.00	17.80	12.93#
0.00	0.00	0.00

U.21
12/18/15

Quantitation Report (Qedit)

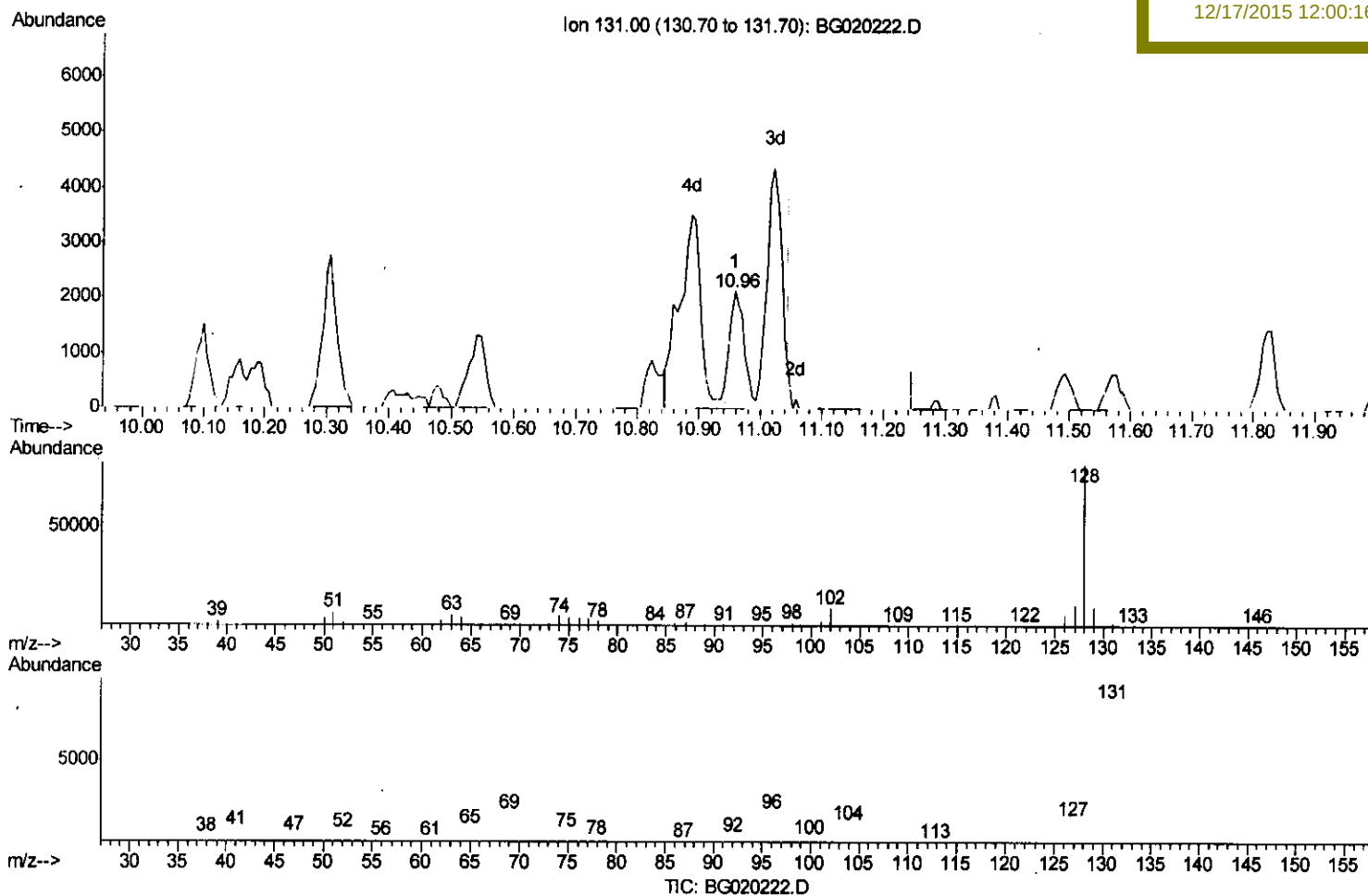
Data Path : Z:\HPCHEM1\BNA_G\Data\BG121715\
Data File : BG020222.D
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(29) 4-Chloroaniline-d4 (S)

10.958min (-0.088) 2.07ng/ul

response 3368

Ion	Exp%	Act%
131.00	100	100
133.00	30.40	27.68
69.00	17.80	16.43
0.00	0.00	0.00

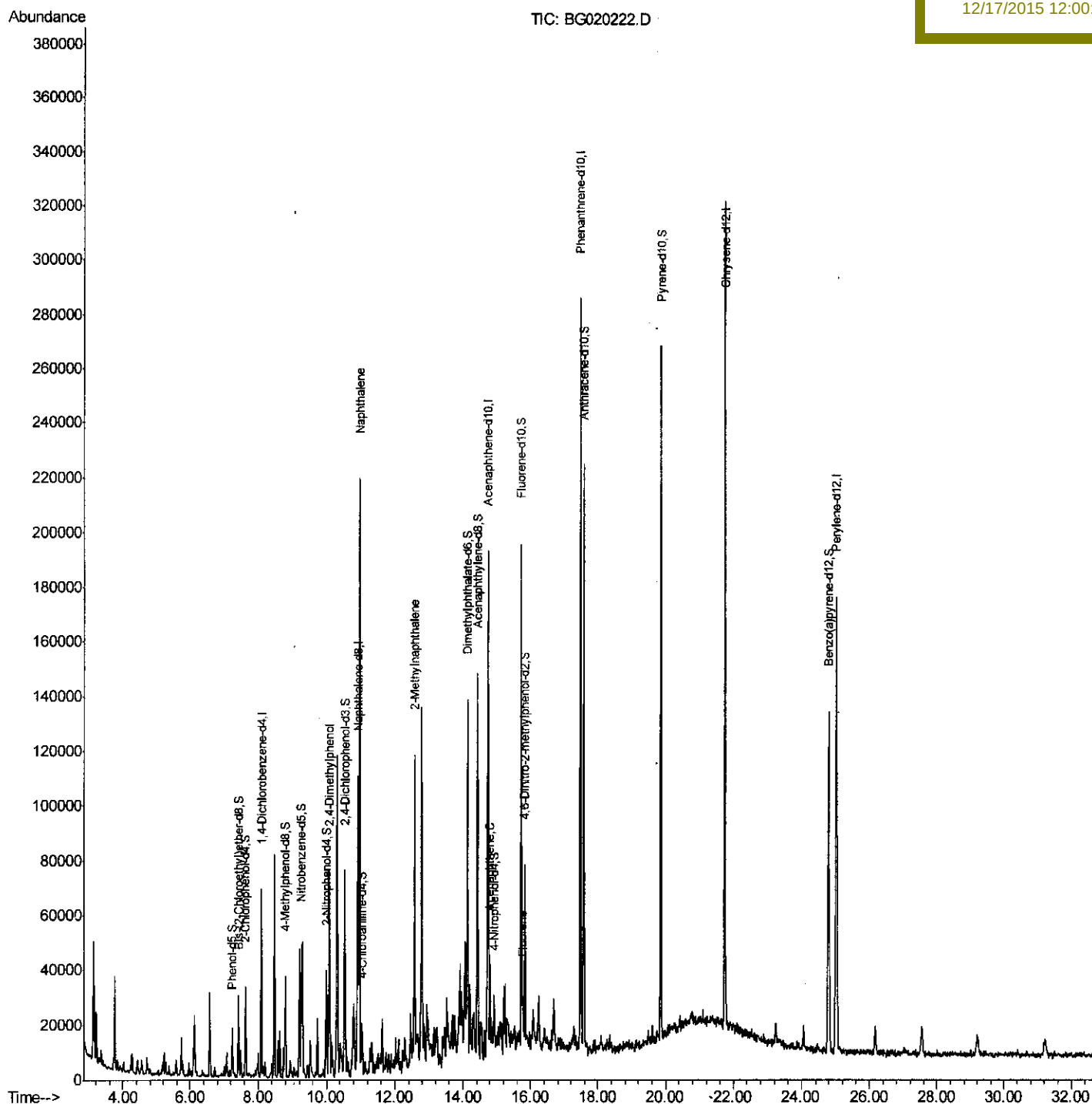
Data Path : Z:\HPCHEM1\BNA G\DATA\BG121715\
Data File : BG020222.D
Acq On : 16 Dec 2015 14:12
Operator : UM/SJ
Sample : G4786-17DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Dec 17 14:32:43 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 17 00:37:15 2015
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Instrument :
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ClientSampled :
G9C90DL

Manual Integrations
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MMdadoda
12/17/2015 12:00:16 PM



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 Operator : UM/SJ
 Sample : G4786-17DL 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA G
 ClientSampleId :
 G9C90DL

Quant Time: Dec 17 14:32:43 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 17 00:37:15 2015
 Response via : Initial Calibration

Manual Integrations
 APPROVED

MMdadoda
 12/17/2015 12:00:16 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	18765	20.00	ng/ul	0.00
18) Naphthalene-d8	10.91	136	82332	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.72	164	66442	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	165718	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	187625	20.00	ng/ul	0.00
86) Perylene-d12	25.01	264	174200	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	7.25	99	5830	3.52	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.42	67	14461	14.65	ng/ul	0.00
9) 2-Chlorophenol-d4	7.62	132	15172	12.61	ng/ul	0.00
13) 4-Methylphenol-d8	8.80	113	11818	8.35	ng/ul	0.00
19) Nitrobenzene-d5	9.27	128	10981	17.11	ng/ul	0.00
22) 2-Nitrophenol-d4	9.99	143	11644	16.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.53	165	23045	15.93	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	7426m	4.57	ng/ul	-0.02
44) Dimethylphthalate-d6	14.12	166	81233	15.12	ng/ul	0.00
47) Acenaphthylene-d8	14.42	160	95741	15.31	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	3266	3.39	ng/ul	0.00
58) Fluorene-d10	15.72	176	74908	15.37	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.82	200	14366	14.62	ng/ul	0.00
71) Anthracene-d10	17.57	188	123400	16.16	ng/ul	0.00
79) Pyrene-d10	19.85	212	137851	15.76	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.78	264	140383	16.16	ng/ul	0.00

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 12/18/15

Target Compounds

						Ovalue
24) 2,4-Dimethylphenol	10.07	107	27540	15.84	ng/ul	99
28) Naphthalene	10.96	128	171058	39.79	ng/ul	99
34) 2-Methylnaphthalene	12.56	142	50118	15.26	ng/ul	99
50) Acenaphthene	14.79	153	14131	3.16	ng/ul	96
59) Fluorene	15.77	166	5825	1.09	ng/ul#	97

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