

Quantitation Report (Qedit)

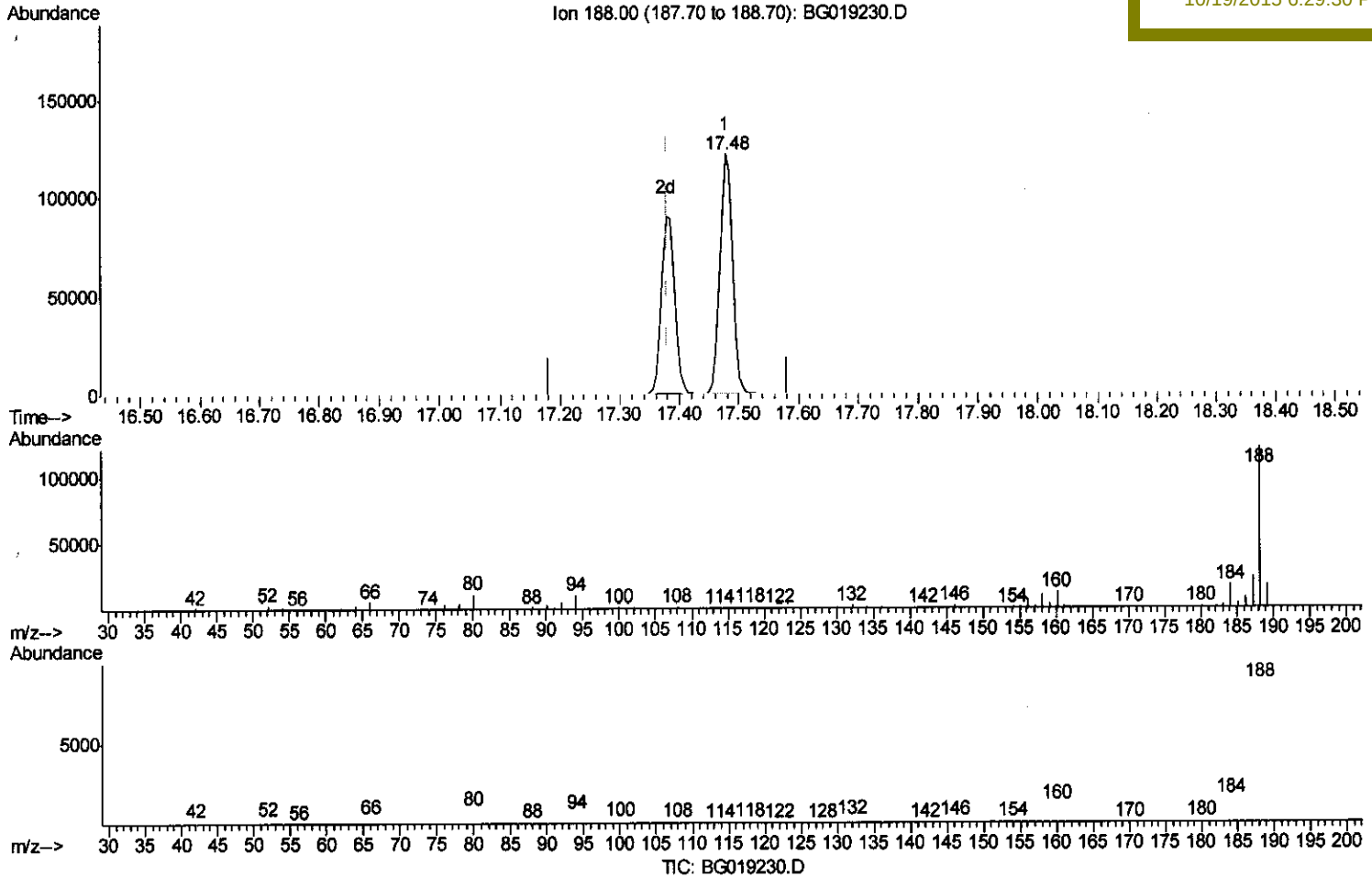
Data Path : Z:\HPCHEM1\BNA G\DATA\BG101615\
 Data File : BG019230.D
 Acq On : 16 Oct 2015 19:27
 Operator : UM/NP
 Sample : G3996-23MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Oct 17 06:34:18 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 17 06:30:39 2015
 Response via : Initial Calibration

Instrument :
 BNA_G
 ClientSampled :
 E5LL0MSD

Manual Integrations
 APPROVED

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(62) Phenanthrene-d10 (I)

17.479min (+0.097) 20.00ng/ul

response 182094

Ion	Exp%	Act%
188.00	100	100
94.00	9.20	8.47
80.00	10.90	9.32
0.00	0.00	0.00

Quantitation Report (Qedit)

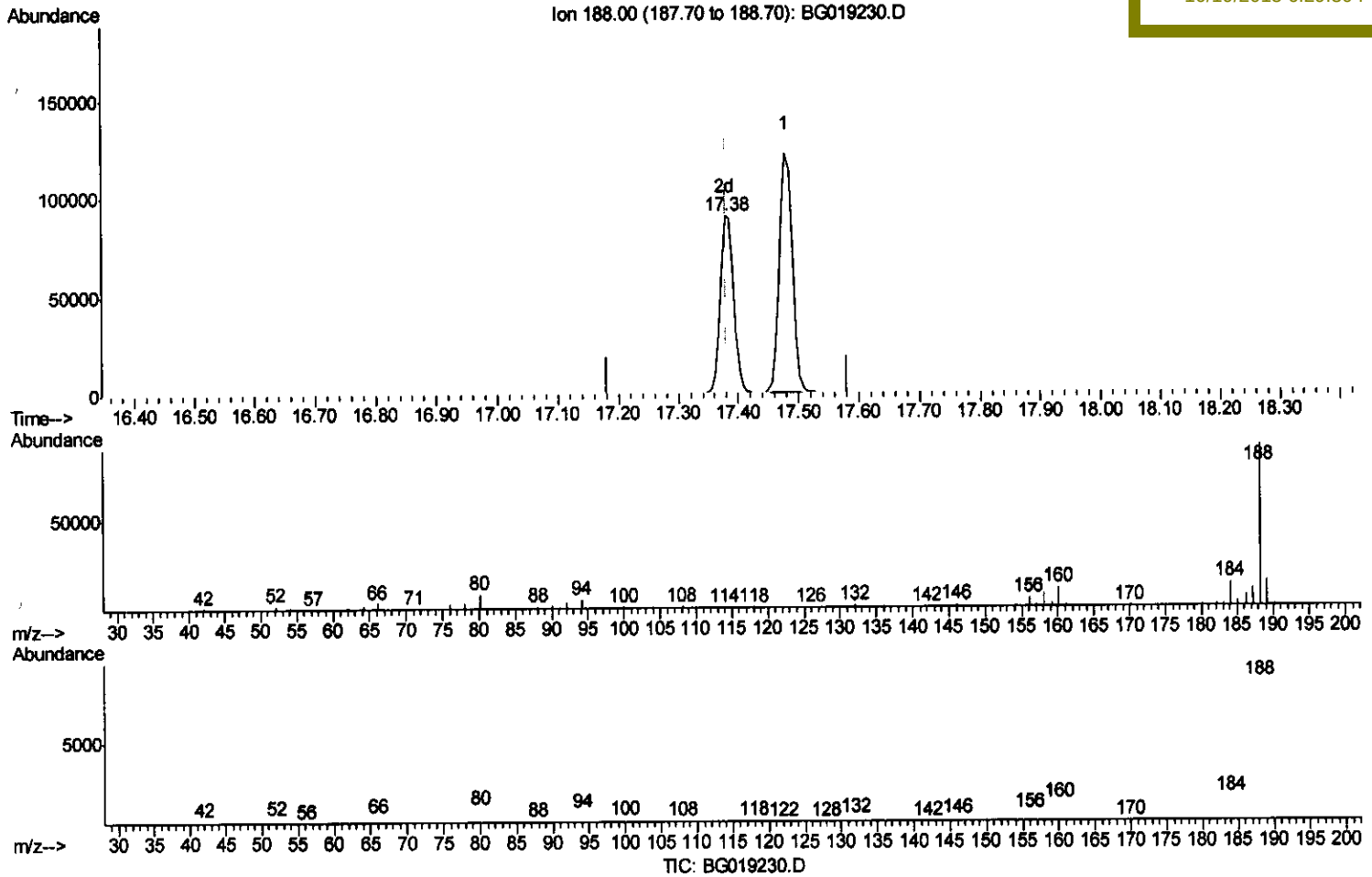
Data Path : Z:\HPCHEM1\BNA G\DATA\BG101615\
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(62) Phenanthrene-d10 (I)

17.379min (-0.002) 20.00ng/ul m

response 138110

Ion	Exp%	Act%
188.00	100	100
94.00	9.20	5.78#
80.00	10.90	9.41
0.00	0.00	0.00

Quantitation Report (Qedit)

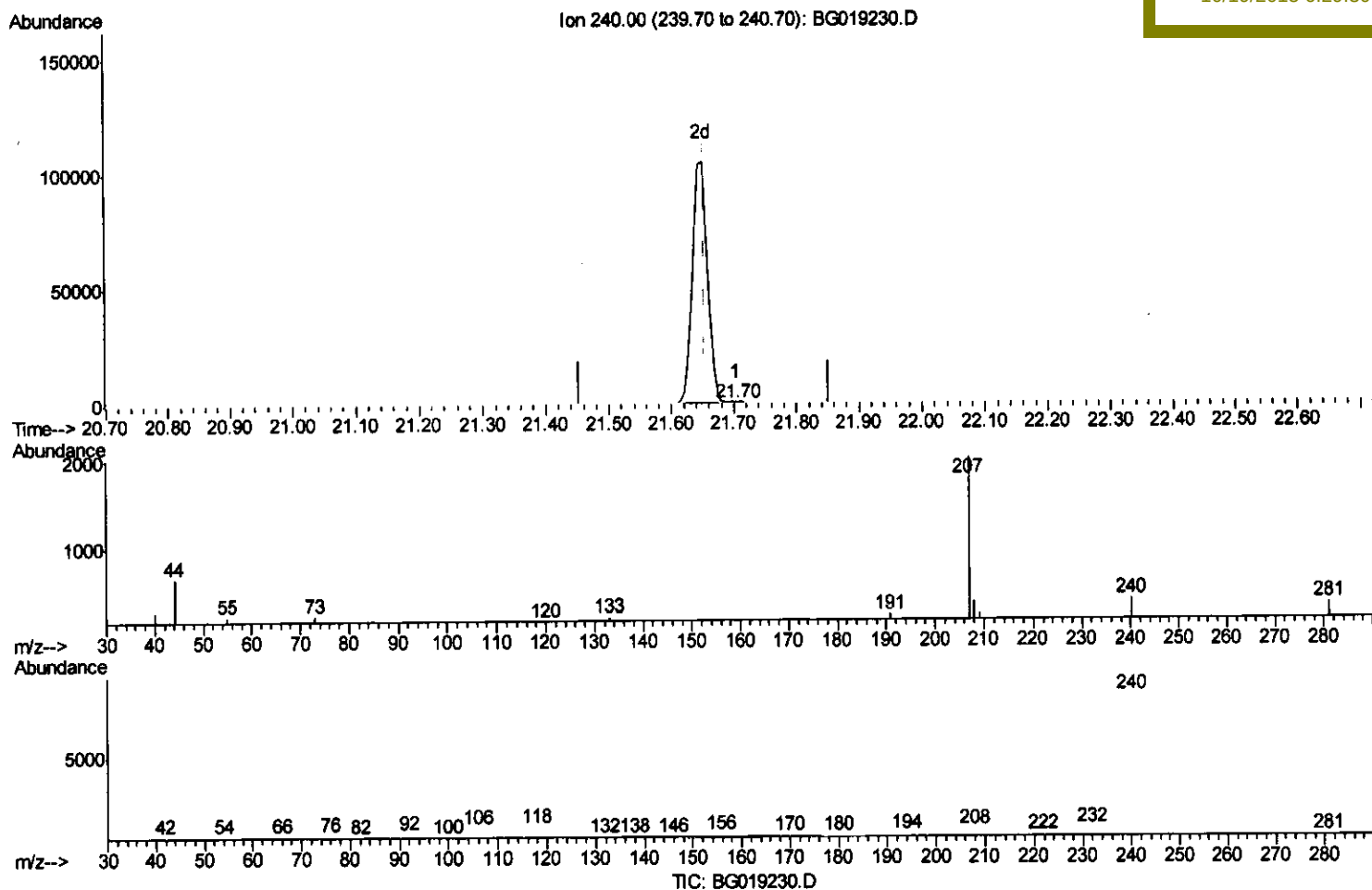
Data Path : Z:\HPCHEM1\BNA G\DATA\BG101615\
Data File : BG019230.D
Acq On : 16 Oct 2015 19:27
Operator : UM/NP
Sample : G3996-23MSD
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E5LL0MSD

Quant Time: Oct 17 07:18:39 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
Quant Title : SVOA CALIBRATION
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(78) Chrysene-d12 (I)

21.703min (+0.050) 20.00ng/ul

response 476

Ion	Exp%	Act%
240.00	100	100
120.00	10.30	39.63#
236.00	26.00	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

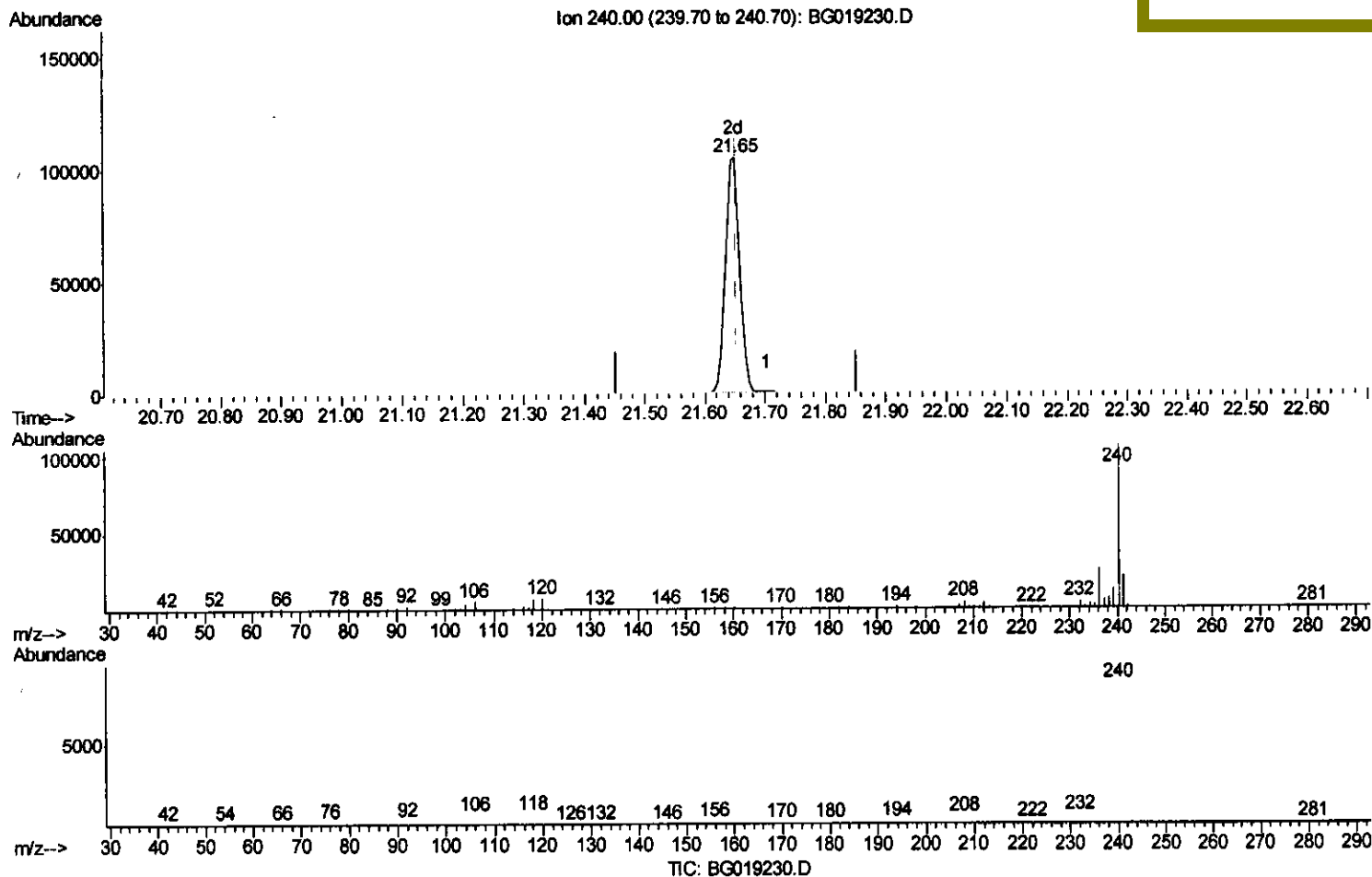
Data Path : Z:\HPCHEM1\BNA G\DATA\BG101615\
 Data File : BG019230.D
 Acq On : 16 Oct 2015 19:27
 Operator : UM/NP
 Sample : G3996-23MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E5LL0MSD

Quant Time: Oct 17 07:18:39 2015
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(78) Chrysene-d12 (I)

21.650min (-0.002) 20.00ng/ul m

response 173607

Ion	Exp%	Act%
240.00	100	100
120.00	10.30	7.15#
236.00	26.00	24.30
0.00	0.00	0.00

Quantitation Report (Qedit)

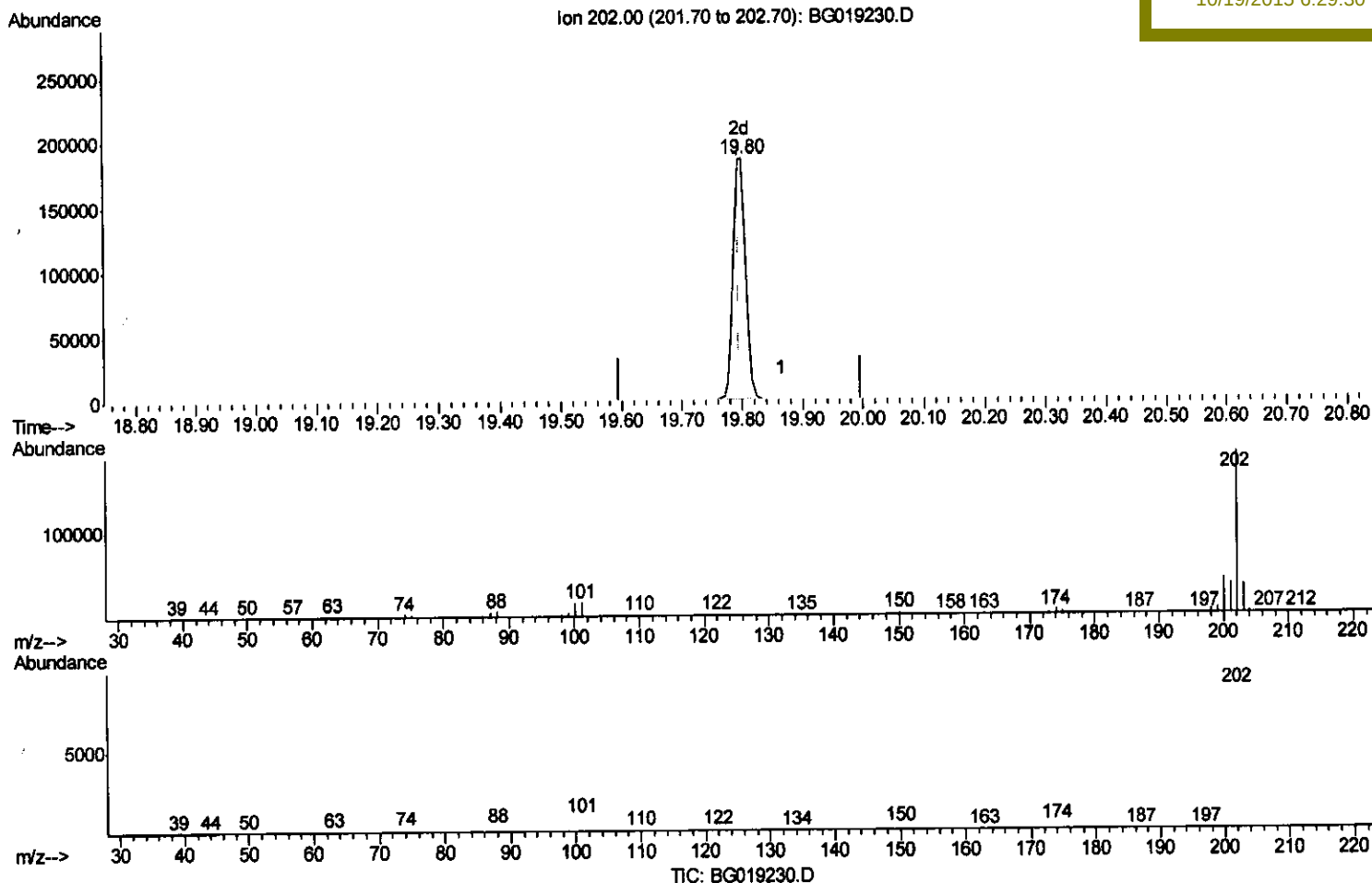
Data Path : Z:\HPCHEM1\BNA_G\Data\BG101615\
 Data File : BG019230.D
 Acq On : 16 Oct 2015 19:27
 Operator : UM/NP
 Sample : G3996-23MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E5LL0MSD

Quant Time: Oct 17 07:19:36 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG101515.M
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(80) Pyrene

19.799min (+0.004) 25.84ng/ul m

response 266292

Ion	Exp%	Act%
202.00	100	100
101.00	12.70	8.67#
100.00	10.60	8.27#
0.00	0.00	0.00

Quantitation Report (Qedit)

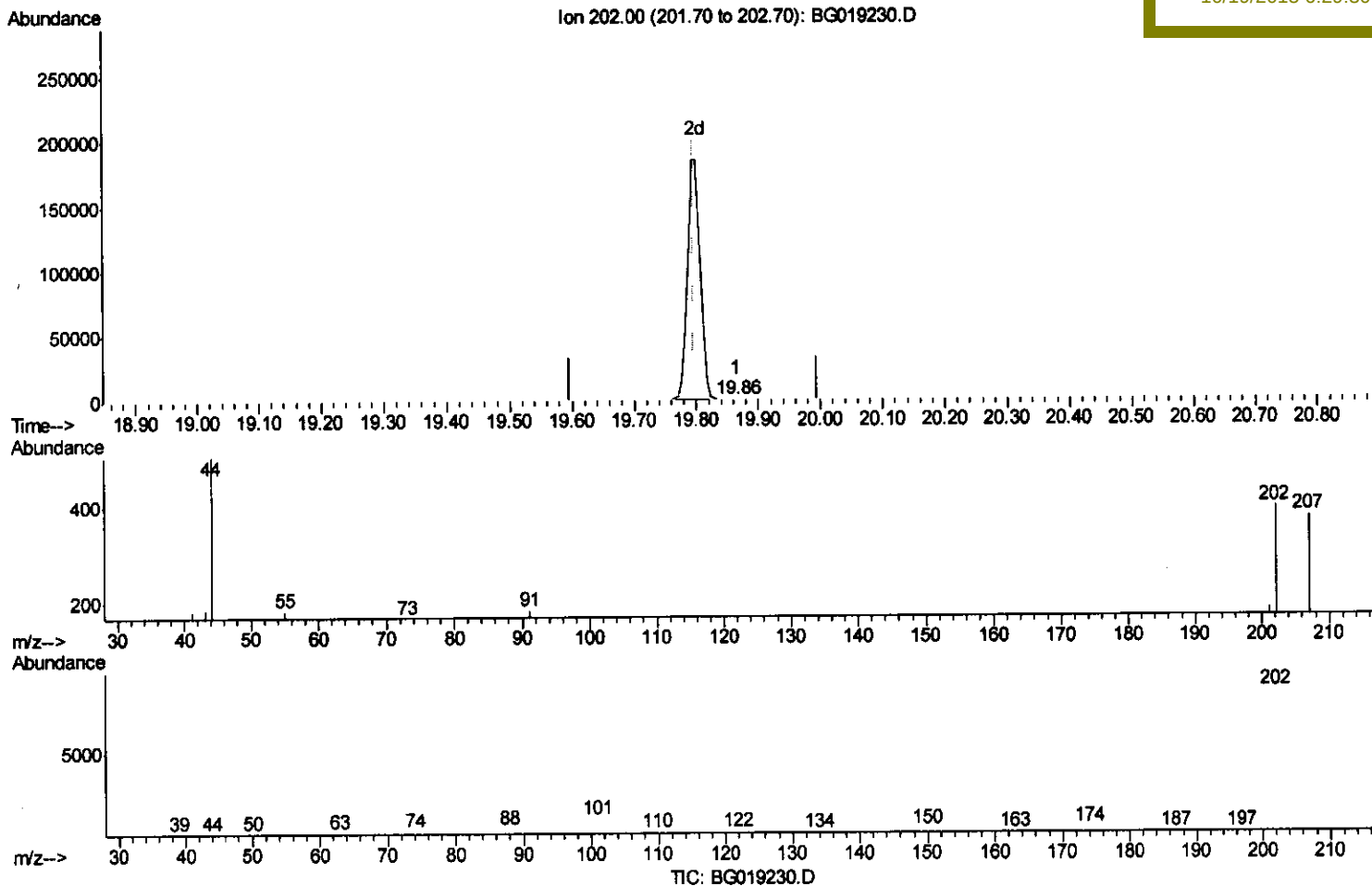
Data Path : Z:\HPCHEM1\BNA_G\Data\BG101615\
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 Acq On : 16 Oct 2015 19:27
 Operator : UM/NP
 Sample : G3996-23MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
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(80) Pyrene

19.864min (+0.068) 0.05ng/ui

response 485

Ion	Exp%	Act%
202.00	100	100
101.00	12.70	0.00#
100.00	10.60	0.00#
0.00	0.00	0.00

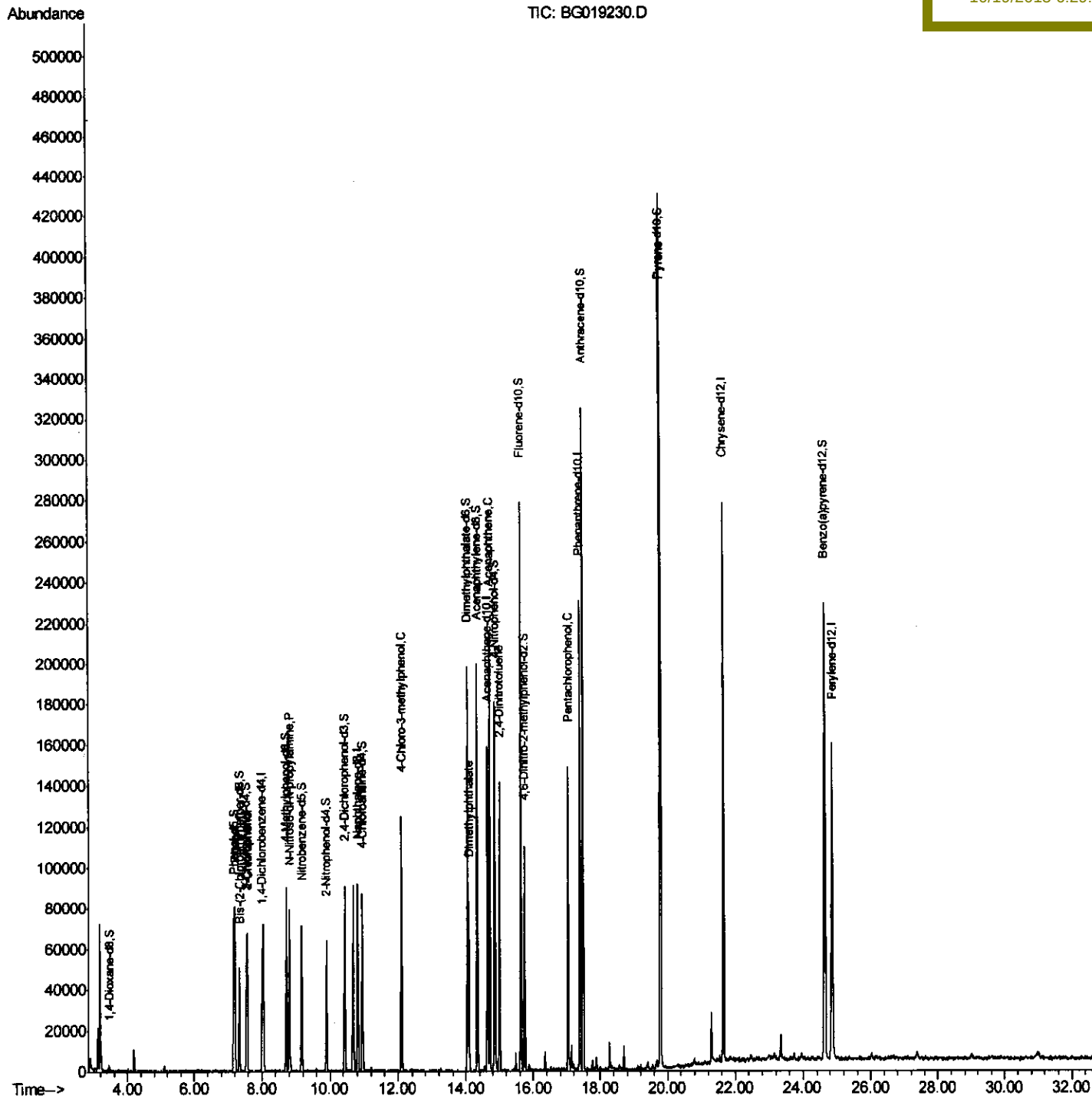
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Data File : BG019230.D
Acq On : 16 Oct 2015 19:27
Operator : UM/NP
Sample : G3996-23MSD
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ESLL0MSD

Quant Time: Oct 19 09:18:26 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG101515.M
Quant Title : SVOA CALIBRATION
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Data Path : Z:\HPCHEM1\BNA_G\Data\BG101615\
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	15421	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	72334	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.63	164	52790	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.38	188	138110m	20.00	ng/ul	0.00
78) Chrysene-d12	21.65	240	173607m	20.00	ng/ul	0.00
86) Perylene-d12	24.85	264	166879	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	1267	4.58	ng/uL	0.00
5) Phenol-d5	7.17	99	36532	27.15	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.33	67	23534	27.17	ng/ul	0.00
9) 2-Chlorophenol-d4	7.54	132	26943	27.47	ng/ul	0.00
13) 4-Methylphenol-d8	8.71	113	30656	26.88	ng/ul	0.00
19) Nitrobenzene-d5	9.16	128	15240	27.35	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	17453	29.51	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.43	165	32016	27.33	ng/ul	0.00
29) 4-Chloroaniline-d4	10.95	131	42762	28.34	ng/ul	0.00
44) Dimethylphthalate-d6	14.04	166	126800	27.46	ng/ul	0.00
47) Acenaphthylene-d8	14.33	160	136449	27.34	ng/ul	0.00
52) 4-Nitrophenol-d4	14.84	143	24709	29.77	ng/ul	0.00
58) Fluorene-d10	15.63	176	106025	26.87	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.74	200	22539	26.45	ng/ul	0.00
71) Anthracene-d10	17.48	188	182094	27.91	ng/ul	0.00
79) Pyrene-d10	19.77	212	225995	28.70	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.63	264	247322	29.34	ng/ul	0.00

Target Compounds

						Qvalue
6) Phenol	7.20	94	35645	25.43	ng/ul#	79
10) 2-Chlorophenol	7.57	128	26787	26.59	ng/ul	99
15) N-Nitroso-di-n-propylamine	8.81	70	25804	26.24	ng/ul	96
33) 4-Chloro-3-methylphenol	12.09	107	42016	28.20	ng/ul	90
45) Dimethylphthalate	14.09	163	49466	10.63	ng/ul	99
50) Acenaphthene	14.70	153	91695	24.88	ng/ul	98
53) 4-Nitrophenol	14.85	109	26533	29.24	ng/ul#	72
55) 2,4-Dinitrotoluene	14.99	165	36995	25.71	ng/ul	97
69) Pentachlorophenol	17.03	266	25768	27.71	ng/ul	93
80) Pyrene	19.80	202	266292m	25.84	ng/ul	

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