

Quantitation Report (QT Reviewed)

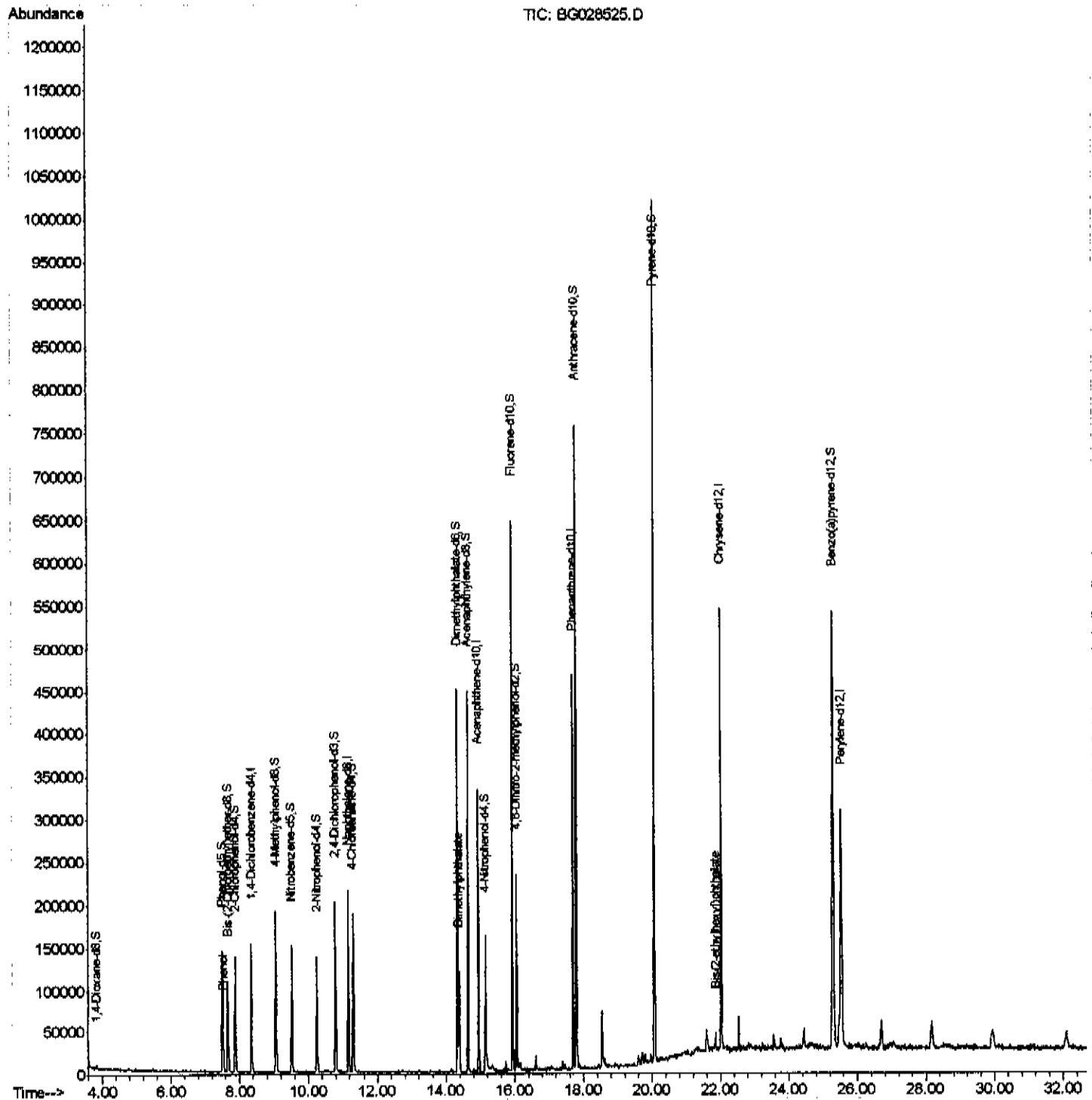
Data Path : Z:\HPCHEM1\BNA G\DATA\BG082817\
 Data File : BG028525.D
 Acq On : 28 Aug 2017 19:13
 Operator : SJ/JU
 Sample : I4905-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampled :
 B0AY5

Manual Integrations
 APPROVED

Sohil
 8/29/2017 6:01:42 PM

Quant Time: Aug 29 04:03:48 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 29 03:57:40 2017
 Response via : Initial Calibration



Quantitation Report (Qedit)

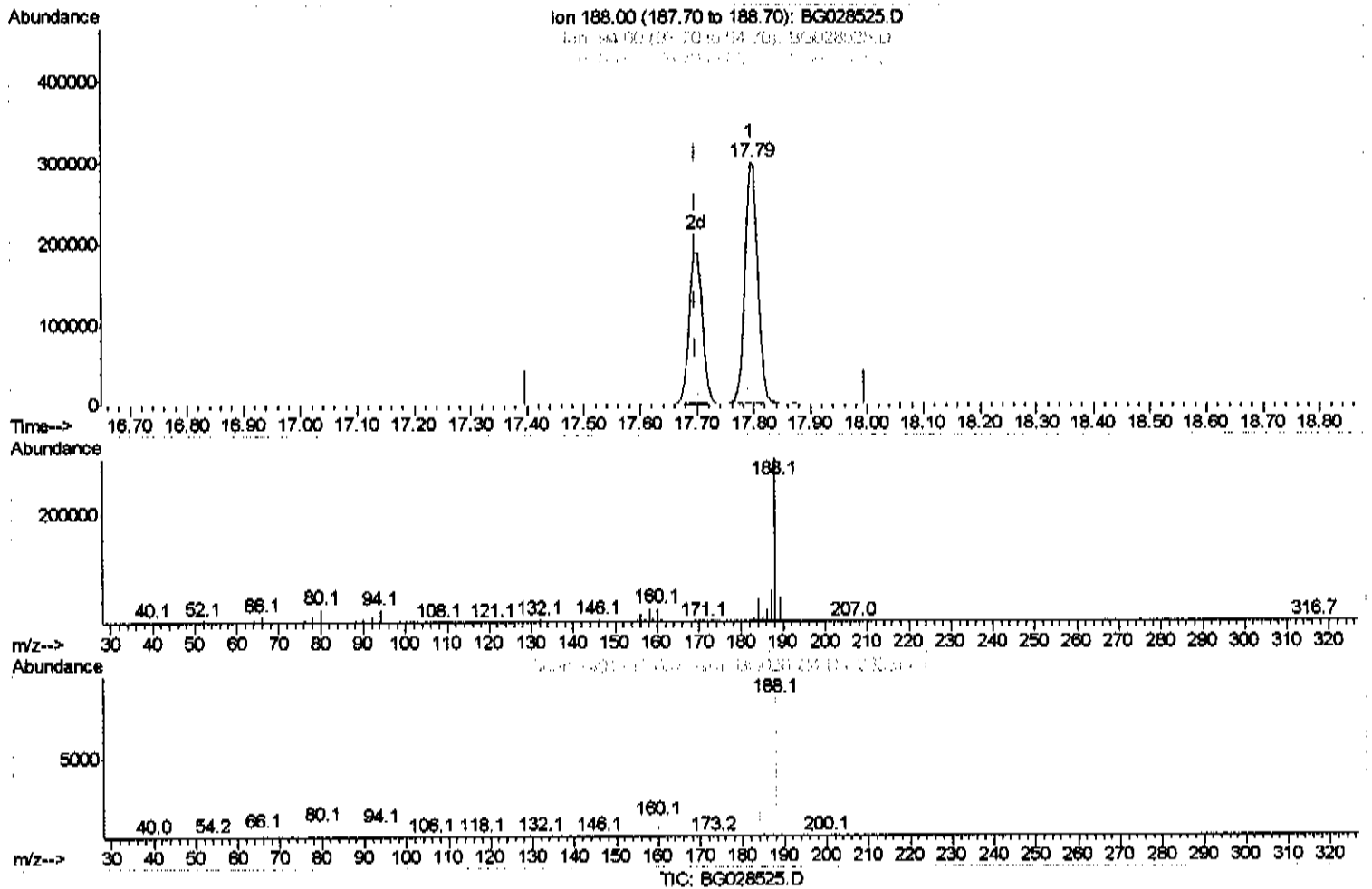
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(61) Phenanthrene-d10 (I)

17.794min (+0.097) 20.00ng/ui

response 473516

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	8.17
80.00	9.50	8.31
0.00	0.00	0.00

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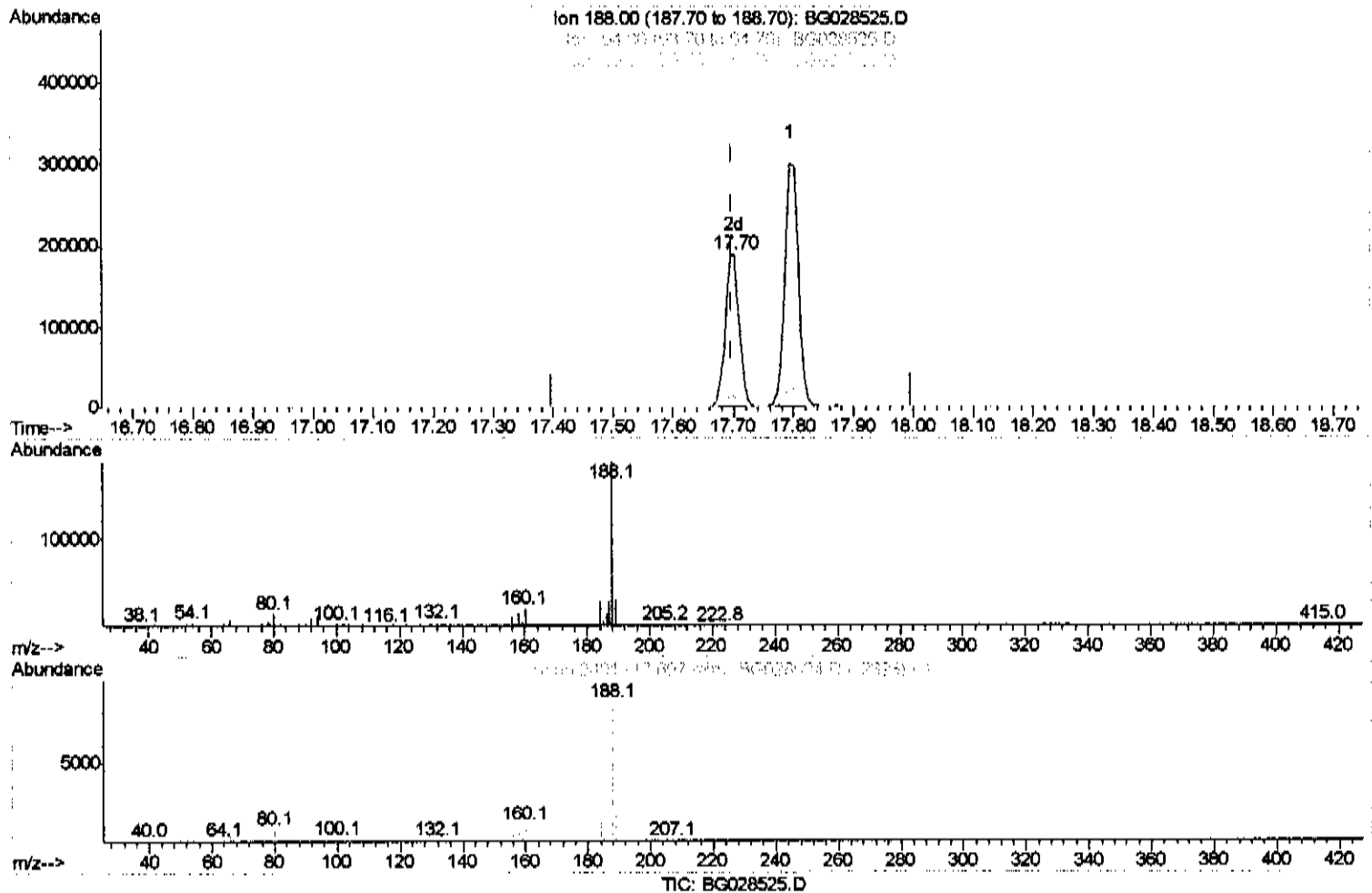
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(61) Phenanthrene-d10 (I)

17.700min (+0.003) 20.00ng/ul m

response 302948

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	6.37%
80.00	9.50	7.48%
0.00	0.00	0.00

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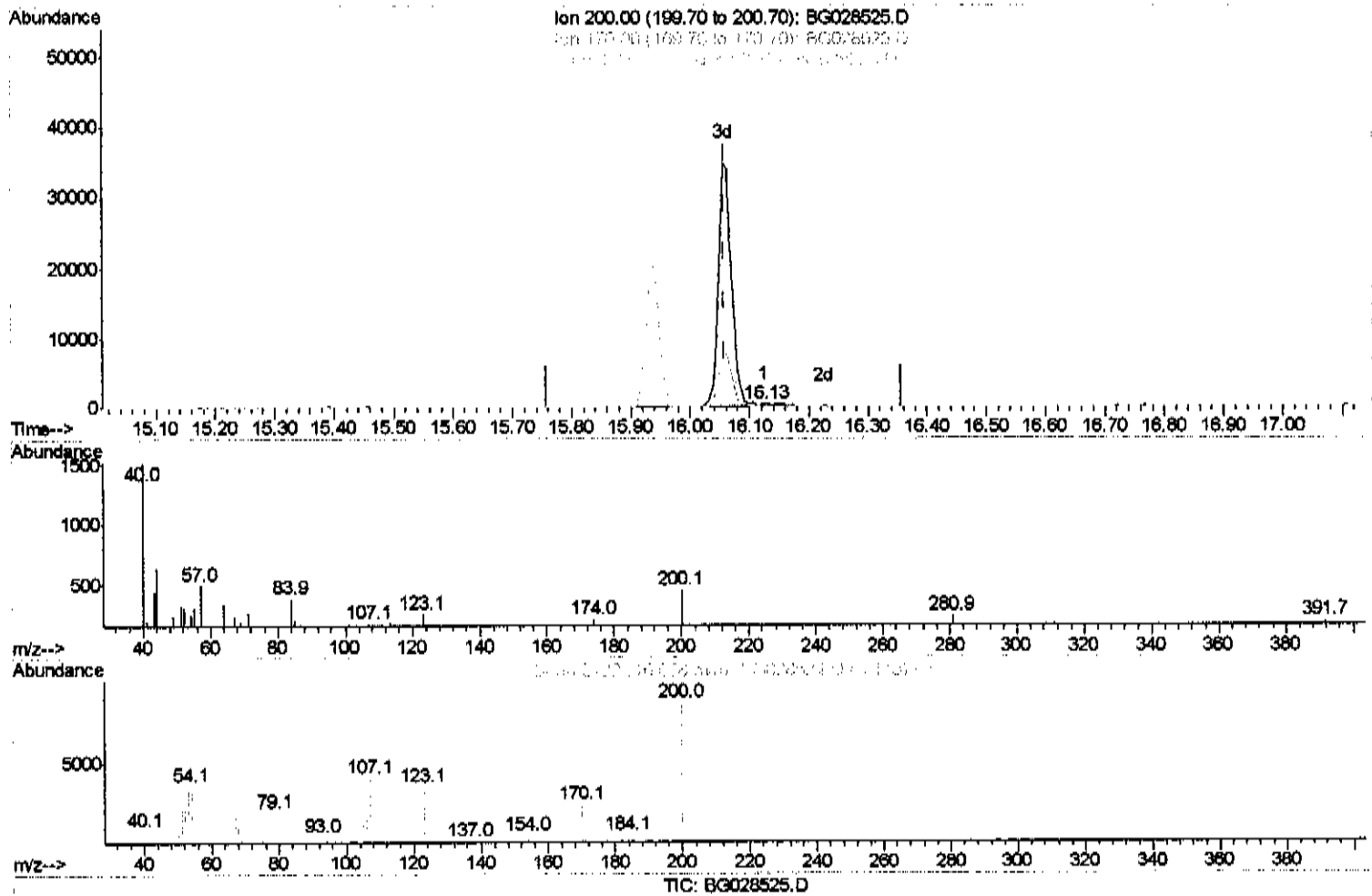
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(62) 4,6-Dinitro-2-methylphenol-d2 (S)

16.128min (+0.068) 0.13ng/ul

response 255

Ion	Exp%	Act%
200.00	100	100
170.00	23.20	0.00#
52.00	47.20	69.47#
0.00	0.00	0.00

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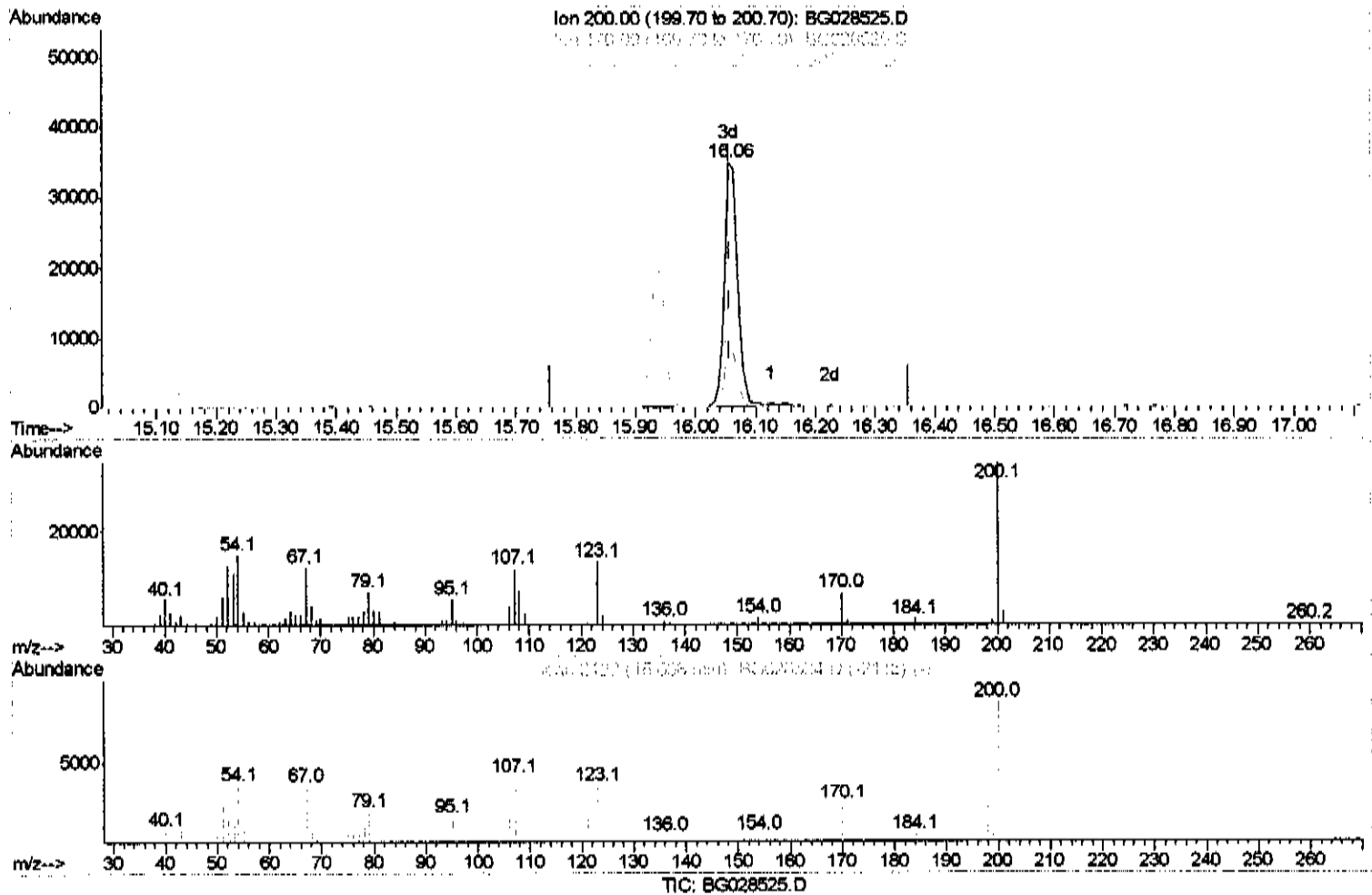
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 Operator : SJ/JU
 Sample : I4905-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
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(62) 4,6-Dinitro-2-methylphenol-d2 (S)

16.055min (-0.003) 27.26ng/ul m

response 53644

Ion	Exp%	Act%
200.00	100	100
170.00	23.20	19.77
52.00	47.20	36.52#
0.00	0.00	0.00

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ClientSampleId :
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.34	152	43321	20.00	ng/ul	0.00
18) Naphthalene-d8	11.16	136	176382	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.94	164	117948	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.70	188	302948m	20.00	ng/ul	0.00
75) Chrysene-d12	22.02	240	357121	20.00	ng/ul	0.00
83) Perylene-d12	25.53	264	355208	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.82	96	3004	3.23	ng/uL	0.00
5) Phenol-d5	7.49	99	84054	17.67	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.65	67	52751	17.51	ng/ul	0.00
9) 2-Chlorophenol-d4	7.87	132	60910	20.29	ng/ul	0.00
13) 4-Methylphenol-d8	9.03	113	77492	19.49	ng/ul	0.00
19) Nitrobenzene-d5	9.51	128	34364	24.92	ng/ul	0.00
22) 2-Nitrophenol-d4	10.23	143	42417	28.05	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.77	165	80756	25.81	ng/ul	0.00
29) 4-Chloroaniline-d4	11.29	131	101744	29.21	ng/ul	0.00
43) Dimethylphthalate-d6	14.33	166	315510	30.08	ng/ul	0.00
46) Acenaphthylene-d8	14.65	160	341935	28.91	ng/ul	0.00
51) 4-Nitrophenol-d4	15.14	143	46599	27.94	ng/ul	0.00
57) Fluorene-d10	15.94	176	269664	28.65	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.06	200	53644m	27.26	ng/ul	0.00
70) Anthracene-d10	17.79	188	473516	32.60	ng/ul	0.00
76) Pyrene-d10	20.07	212	593601	32.41	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.29	264	578781	34.81	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Phenol	7.52	94	5396	1.141	ng/ul#	75
44) Dimethylphthalate	14.38	163	73631	7.623	ng/ul	98
81) Bis(2-ethylhexyl)phthalate	21.86	149	11932	1.051	ng/ul#	97

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