

Quantitation Report (QT Reviewed)

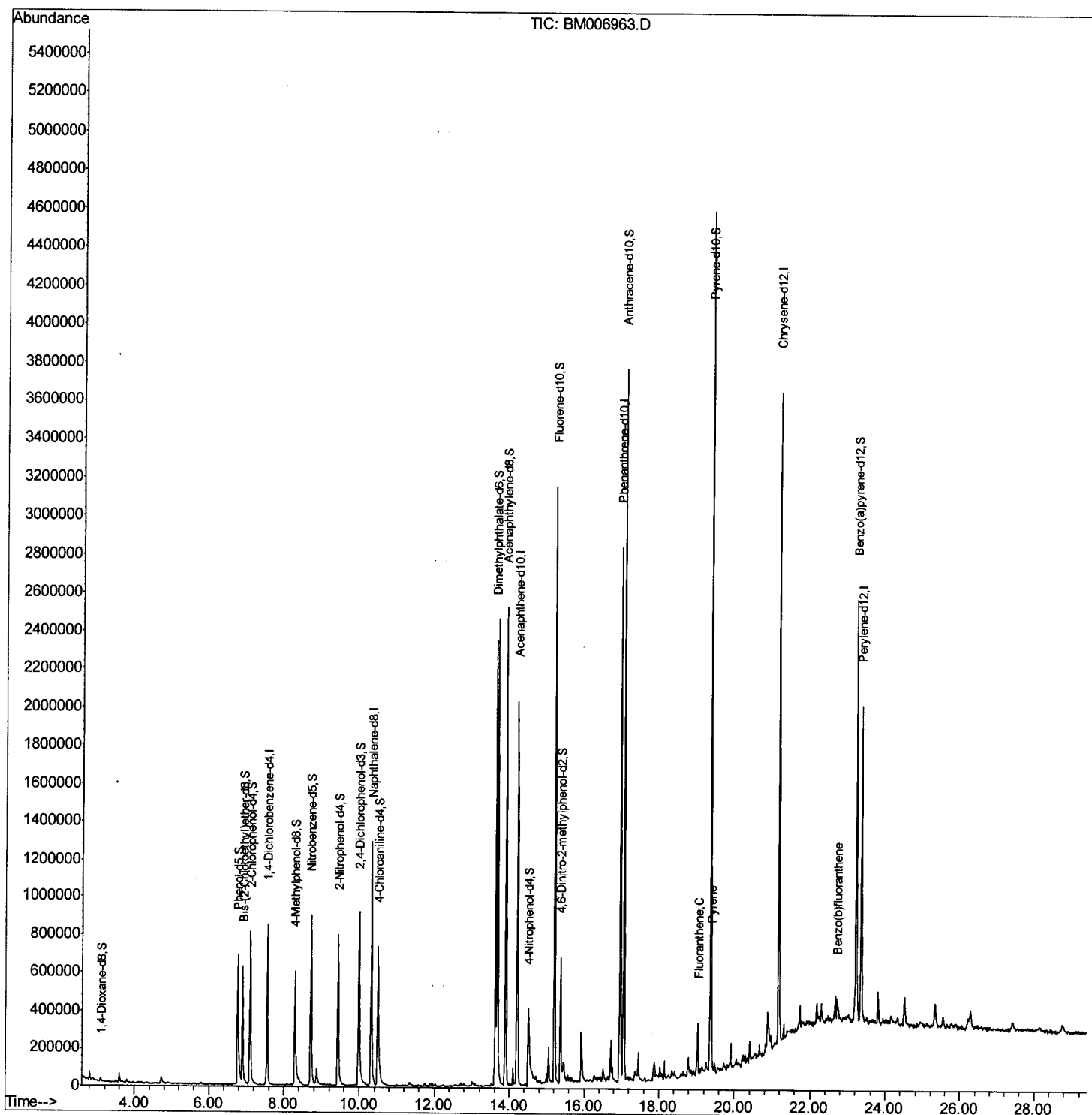
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080616\
 Data File : BM006963.D
 Acq On : 07 Aug 2016 06:55
 Operator : UM/SJ
 Sample : H4301-20
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA0G1

Manual Integrations
 APPROVED

sohil
 8/8/2016 7:06:40 PM

Quant Time: Aug 08 04:38:37 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 08 03:51:46 2016
 Response via : Initial Calibration



Quantitation Report (Qedit)

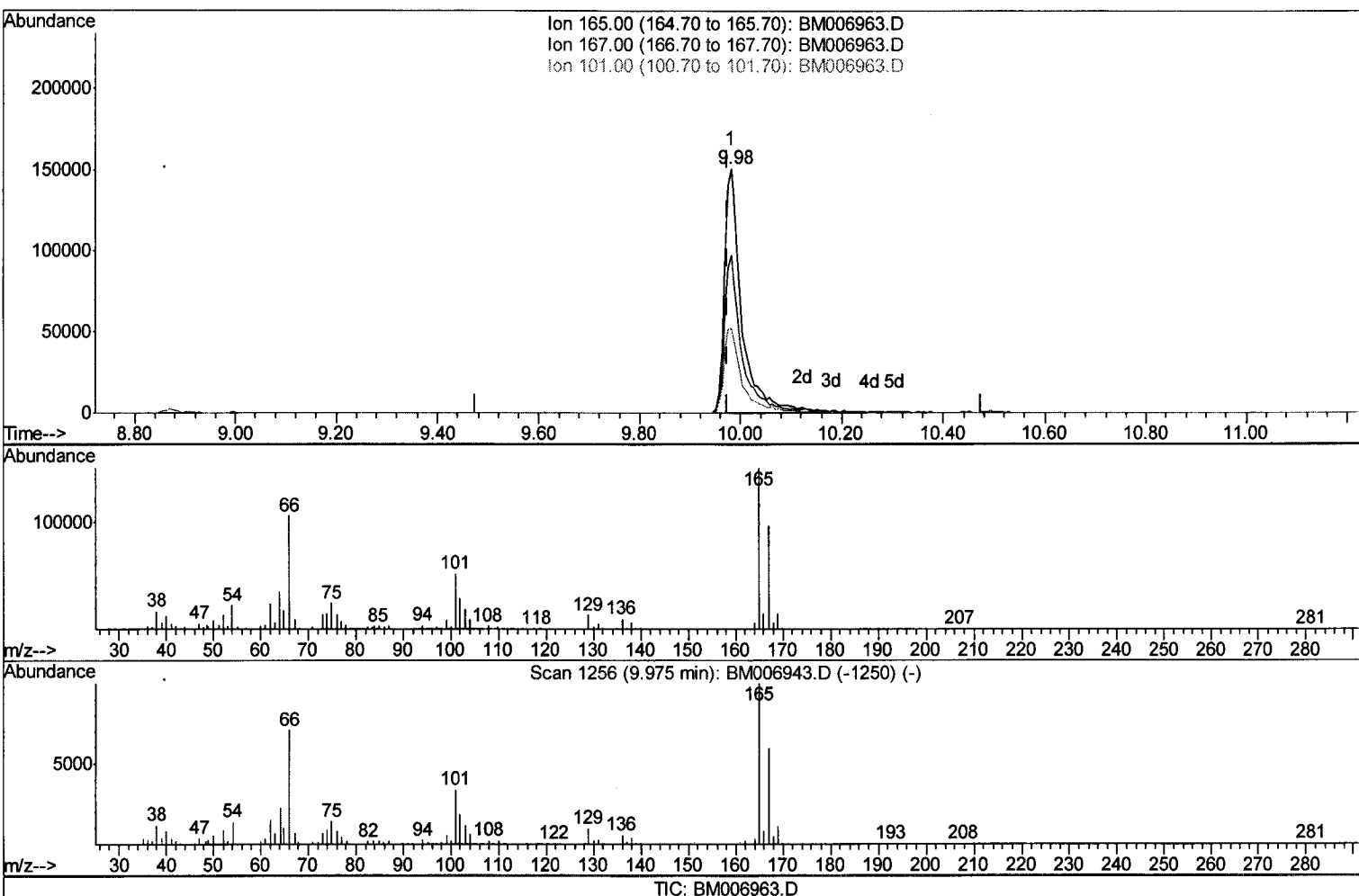
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(10) 2,4-Dichlorophenol-d3 (S)

9.981min (+0.006) 22.41ng/ul m

response 357114

Ion	Exp%	Act%
165.00	100	100
167.00	62.10	64.65
101.00	34.00	34.51
0.00	0.00	0.00

UM
 08/11/16

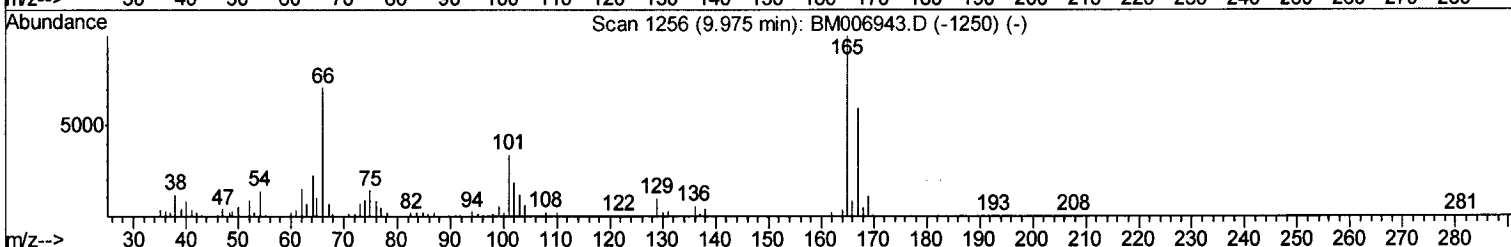
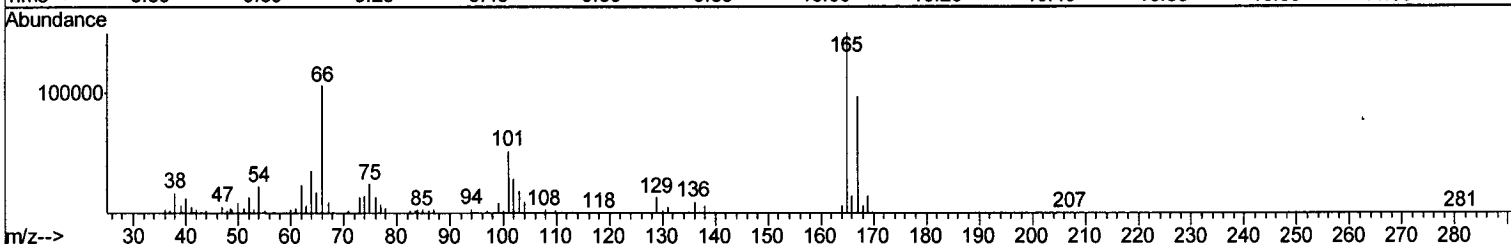
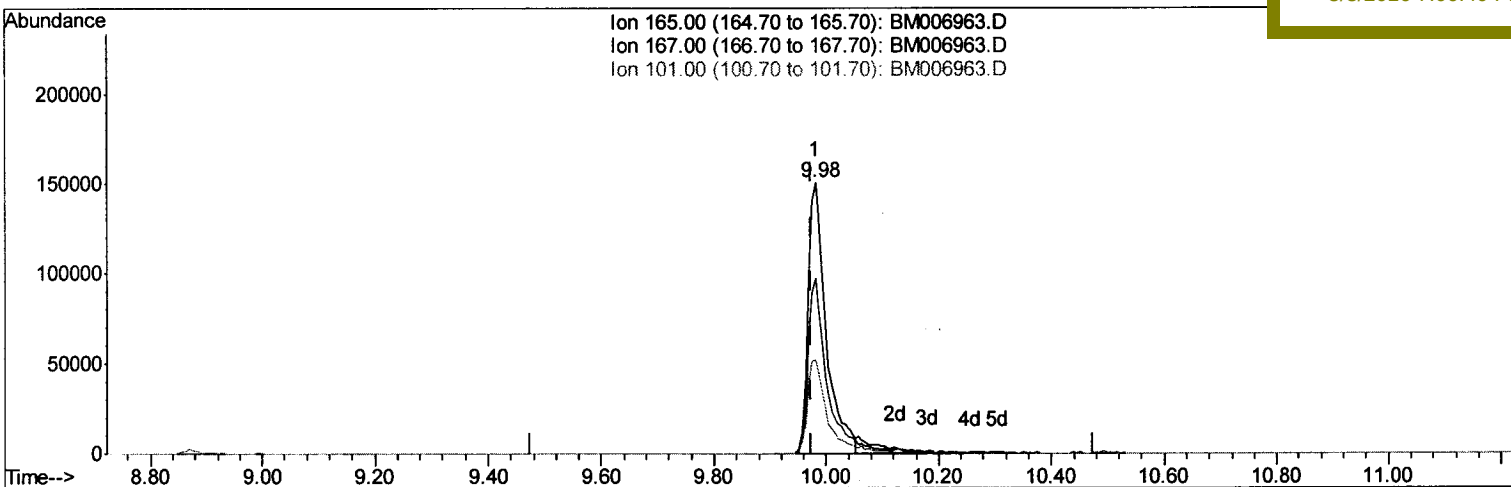
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TIC: BM006963.D

(10) 2,4-Dichlorophenol-d3 (S)

9.981min (+0.006) 21.20ng/ul

response 337823

Ion	Exp%	Act%
165.00	100	100
167.00	62.10	64.65
101.00	34.00	34.51
0.00	0.00	0.00

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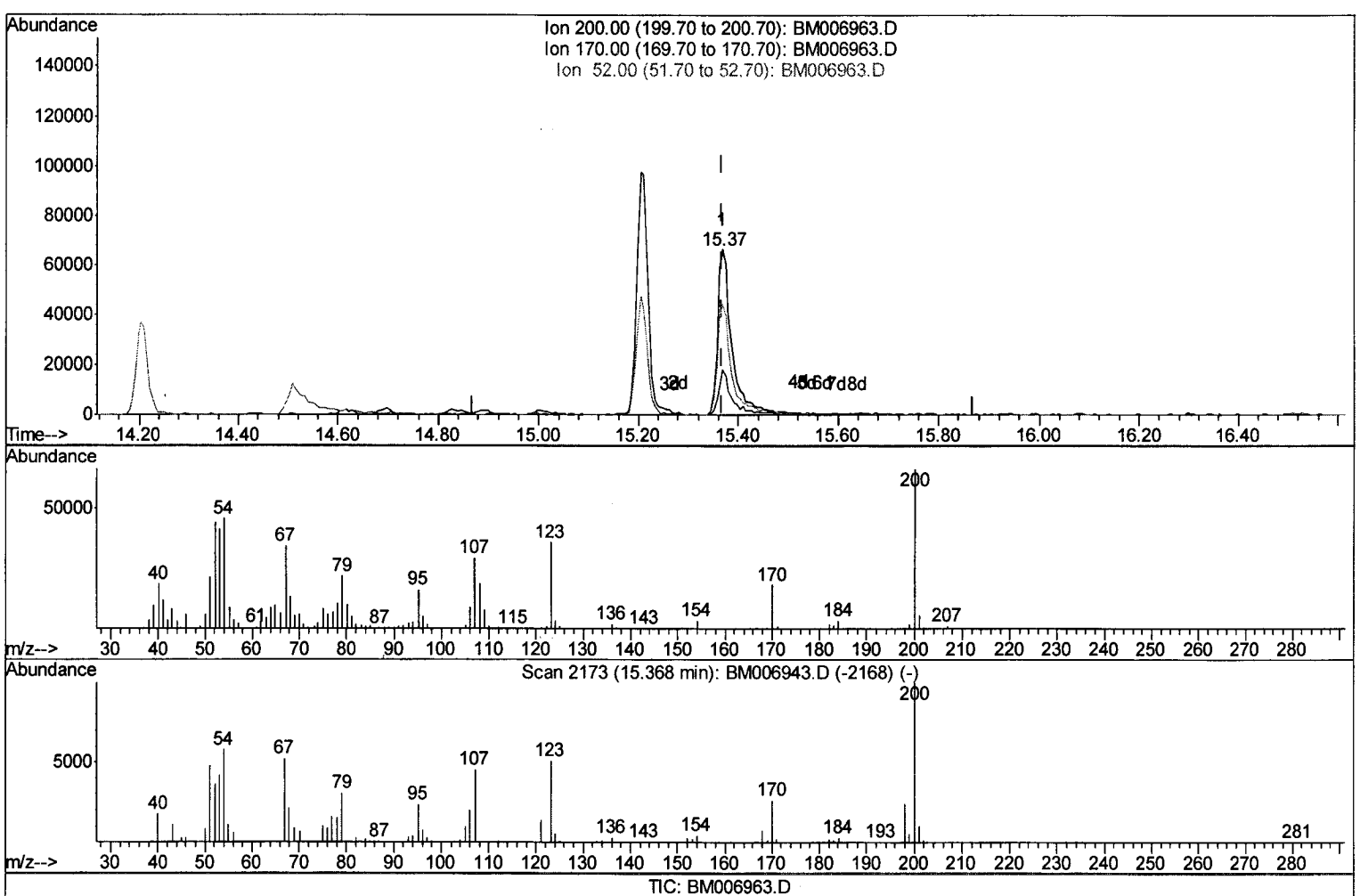
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(24) 4,6-Dinitro-2-methylphenol-d2 (S)

15.369min (+0.000) 21.16ng/ul m

response 131295

Ion	Exp%	Act%
200.00	100	100
170.00	28.10	27.63
52.00	93.50	66.70#
0.00	0.00	0.00

4M
 08/11/16

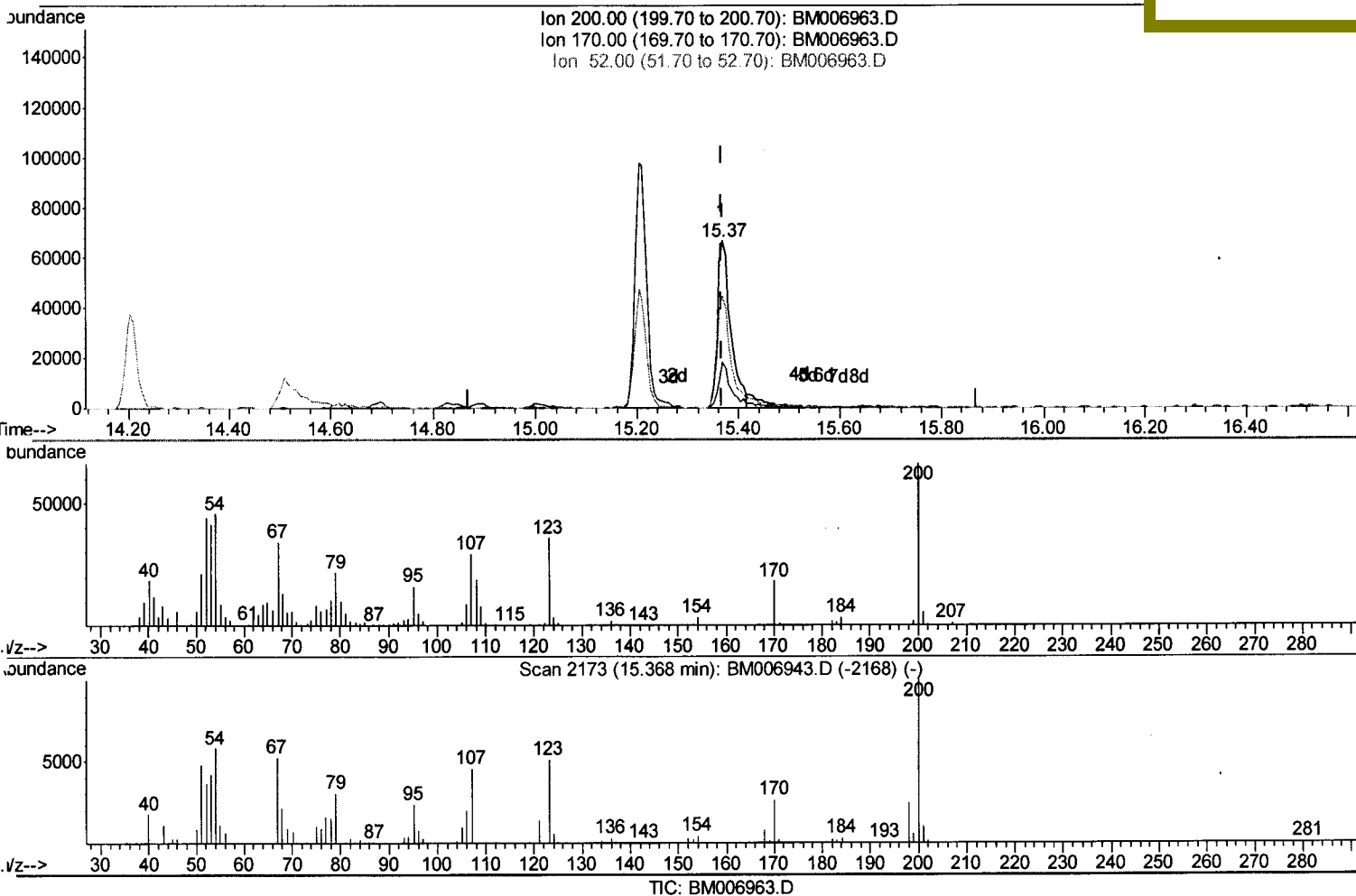
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(24) 4,6-Dinitro-2-methylphenol-d2 (S)

15.369min (+0.000) 19.51ng/ul

response 121029

Ion	Exp%	Act%
200.00	100	100
170.00	28.10	27.63
52.00	93.50	66.70#
0.00	0.00	0.00

Quantitation Report (Qedit)

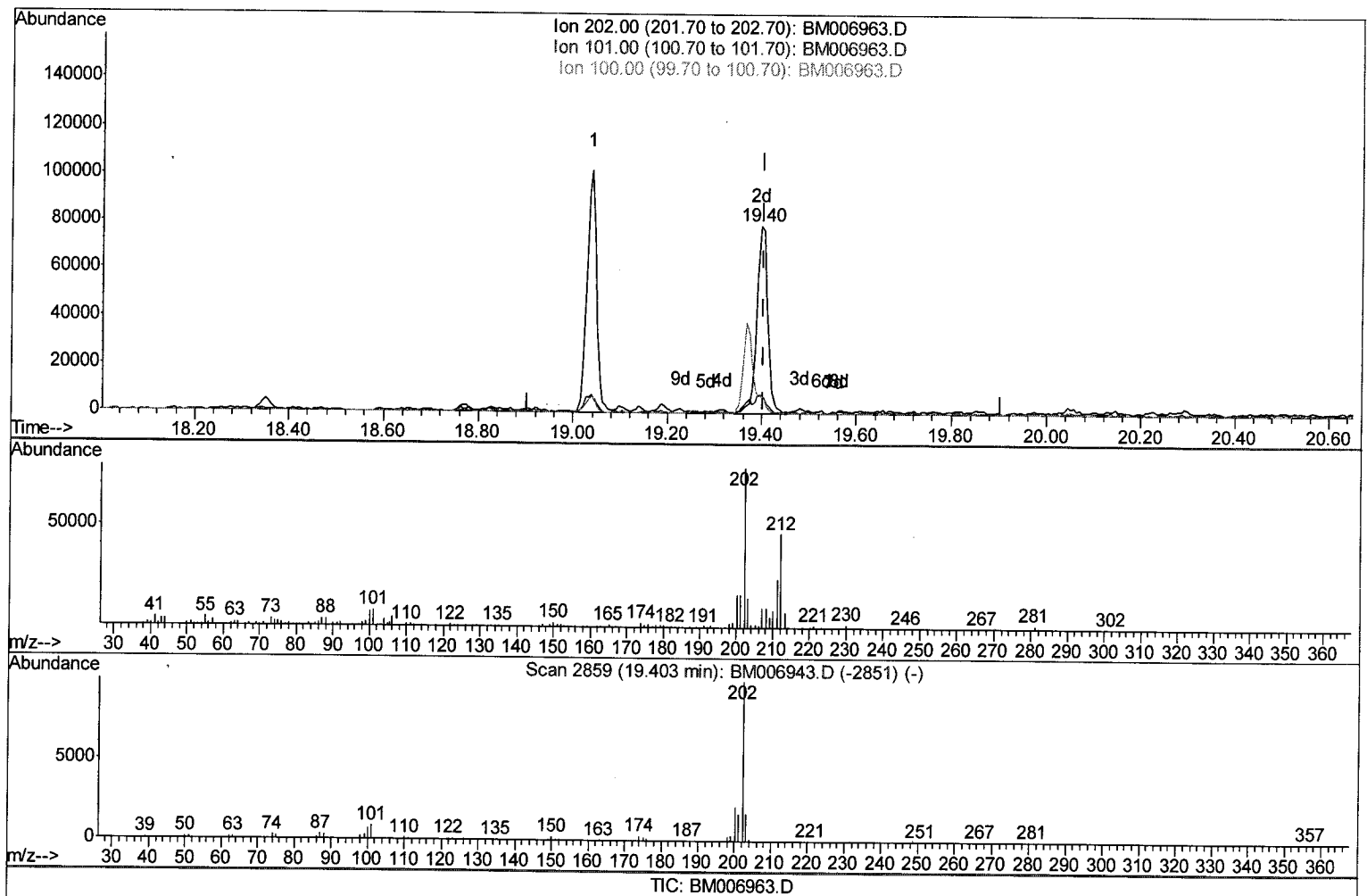
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Instrument :
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(31) Pyrene

19.398min (-0.006) 1.29ng/ul m

response 125486

Ion	Exp%	Act%
202.00	100	100
101.00	8.00	9.99#
100.00	6.80	9.11#
0.00	0.00	0.00

V.M
 08/11/16

Quantitation Report (Qedit)

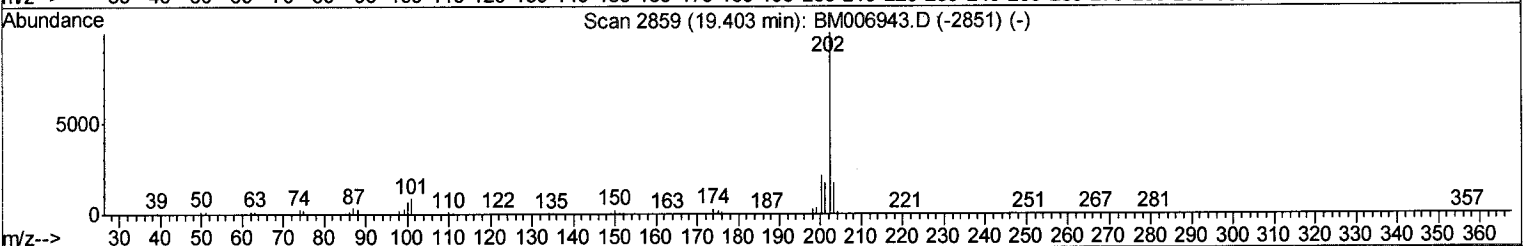
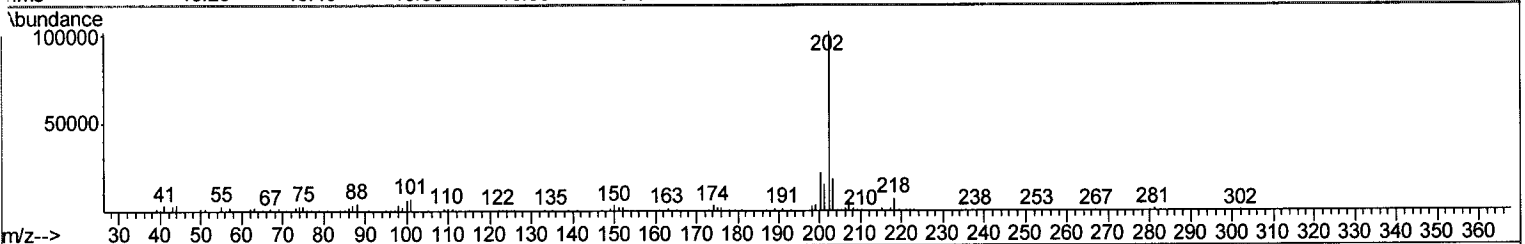
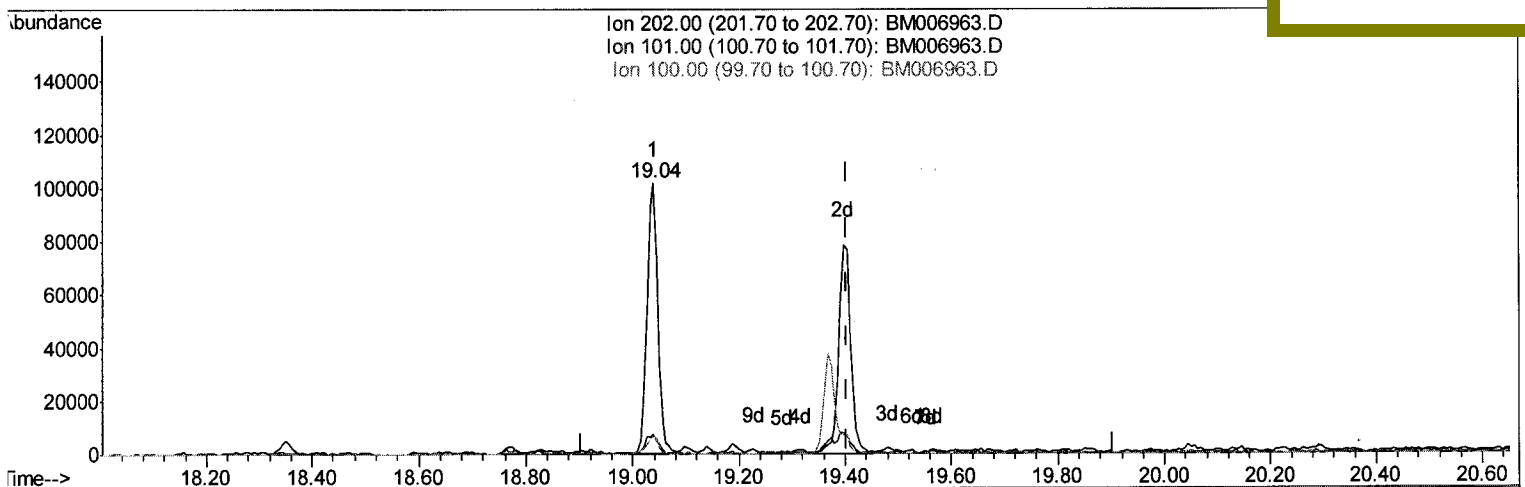
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TIC: BM006963.D

(31) Pyrene

19.039min (-0.364) 1.45ng/ul

response 141008

Ion	Exp%	Act%
202.00	100	100
101.00	8.00	7.34
100.00	6.80	6.27
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	193614	20.00	ng/ul	0.00
7) Naphthalene-d8	10.32	136	944916	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.20	164	623241	20.00	ng/ul	0.00
23) Phenanthrene-d10	16.96	188	1502368	20.00	ng/ul	0.00
29) Chrysene-d12	21.17	240	1566330	20.00	ng/ul	0.00
34) Perylene-d12	23.36	264	1167809	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.09	96	9487	2.24	ng/uL	0.00
3) Phenol-d5	6.77	99	368825	20.09	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.90	67	277869	24.14	ng/ul	0.00
5) 2-Chlorophenol-d4	7.10	132	303918	23.19	ng/ul	0.00
6) 4-Methylphenol-d8	8.29	113	231851	15.13	ng/ul	0.00
8) Nitrobenzene-d5	8.72	128	164682	23.46	ng/ul	0.00
9) 2-Nitrophenol-d4	9.43	143	195936	23.31	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	9.98	165	357114m	22.41	ng/ul	0.00
12) 4-Chloroaniline-d4	10.49	131	410795	21.59	ng/ul	0.00
16) Dimethylphthalate-d6	13.63	166	1366412	24.91	ng/ul	0.00
17) Acenaphthylene-d8	13.90	160	1611866	25.41	ng/ul	0.00
20) 4-Nitrophenol-d4	14.52	143	125700	15.80	ng/ul	0.02
21) Fluorene-d10	15.21	176	1189086	25.51	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.37	200	131295m	21.16	ng/ul	0.00
26) Anthracene-d10	17.06	188	1878345	25.88	ng/ul	0.00
30) Pyrene-d10	19.37	212	2235031	29.93	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.22	264	1488326	26.76	ng/ul	0.00

Target Compounds

						Qvalue
28) Fluoranthene	19.04	202	140309	1.37	ng/ul	98
31) Pyrene	19.40	202	125486m	1.29	ng/ul	
35) Benzo(b)fluoranthene	22.71	252	93318	1.22	ng/ul#	95

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