

Quantitation Report (QT Reviewed)

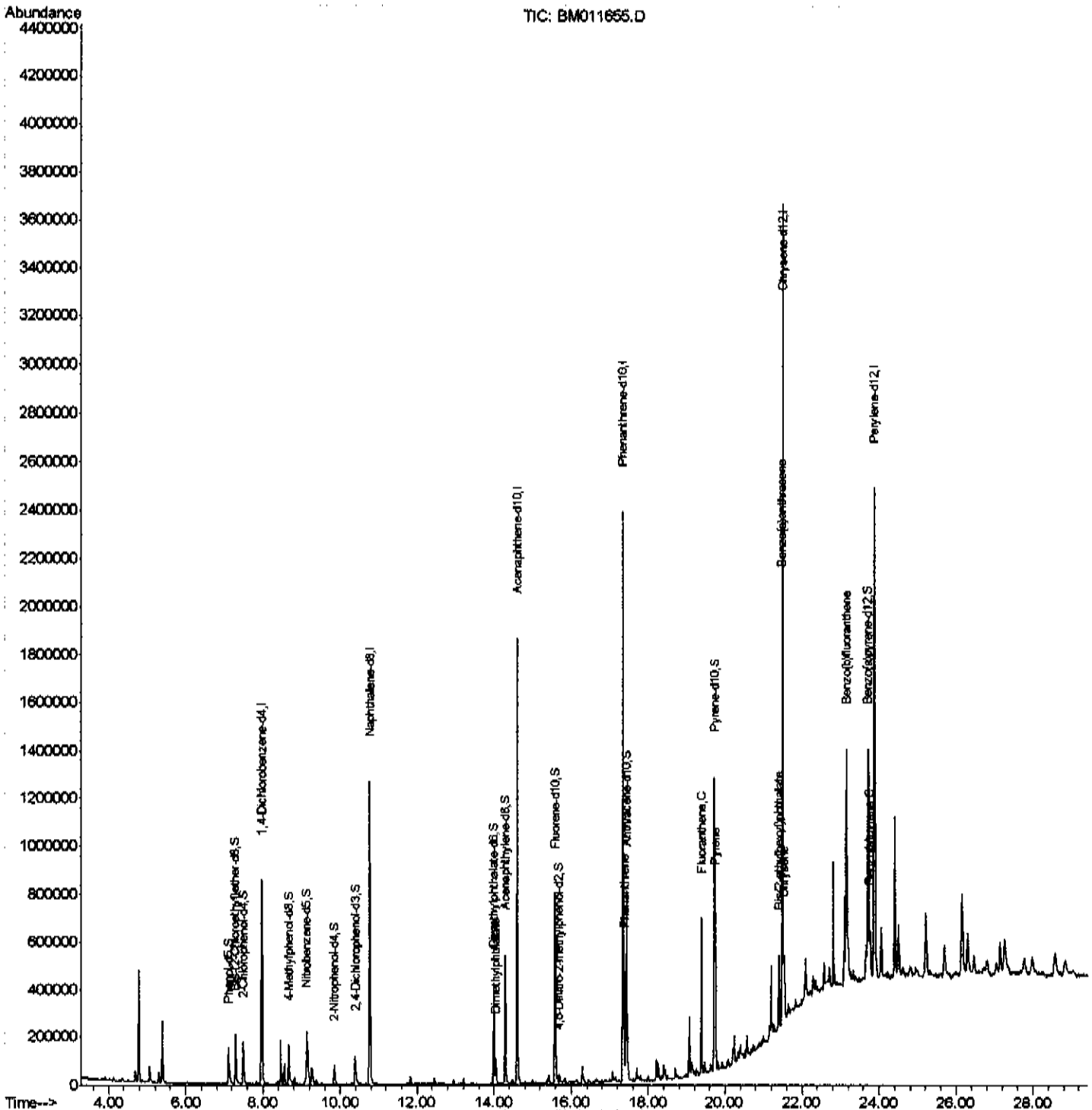
Data Path : Z:\HPCHEM1\BNA M\DATA\BM092117\
 Data File : BM011655.D
 Acq On : 21 Sep 2017 15:59
 Operator : SJ/JU
 Sample : I5284-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 CB3C9

Manual Integrations
 APPROVED

Sohil
 9/22/2017 4:06:05 PM

Quant Time: Sep 22 07:17:44 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 22 03:47:29 2017
 Response via : Initial Calibration



Quantitation Report (Qedit)

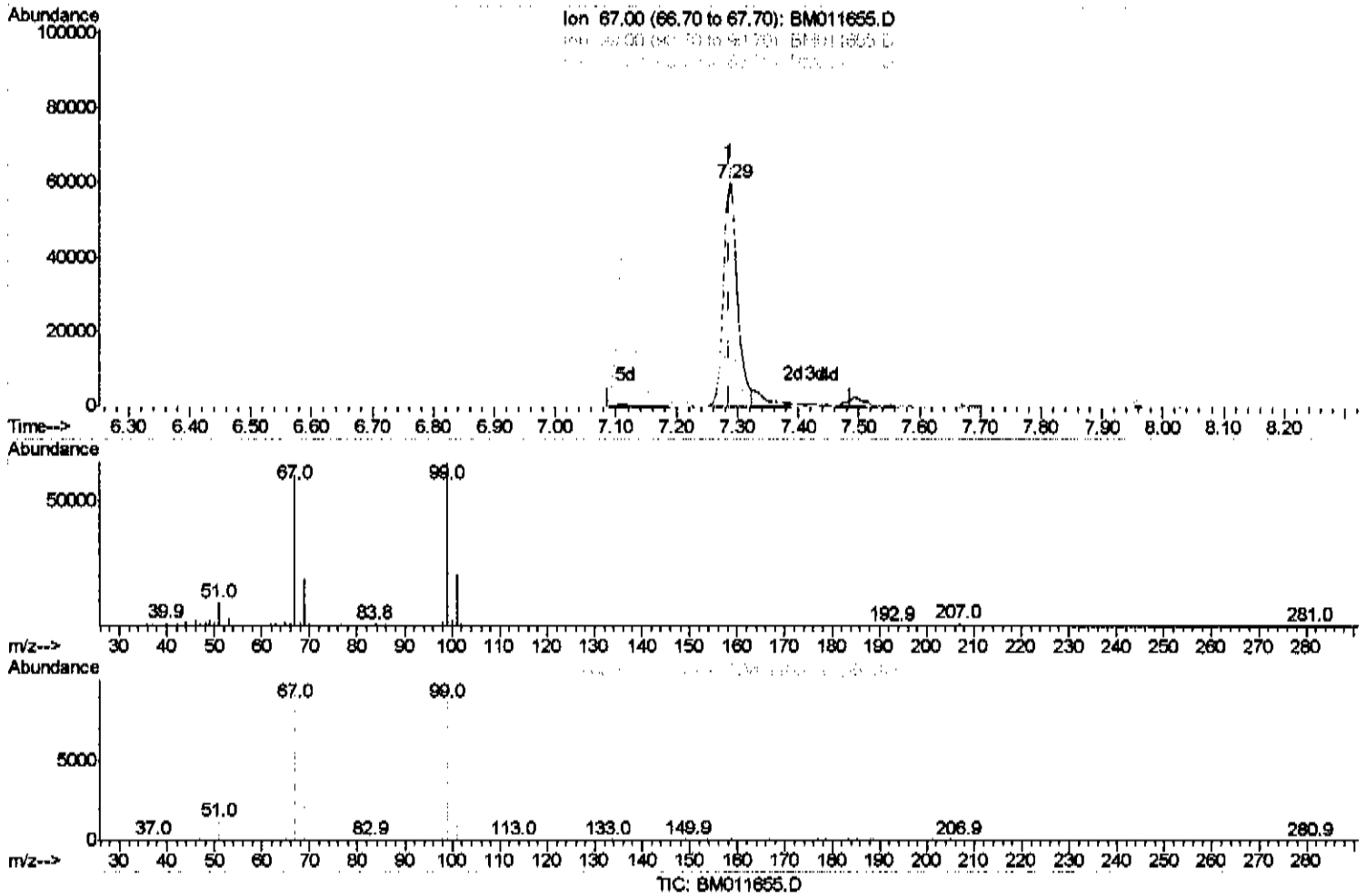
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(7) Bis-(2-Chloroethyl)ether-d8 (S)

7.287min (+0.000) 6.89ng/ul

response 99102

Ion	Exp%	Act%
67.00	100	100
99.00	117.60	108.10
69.00	33.70	31.19
0.00	0.00	0.00

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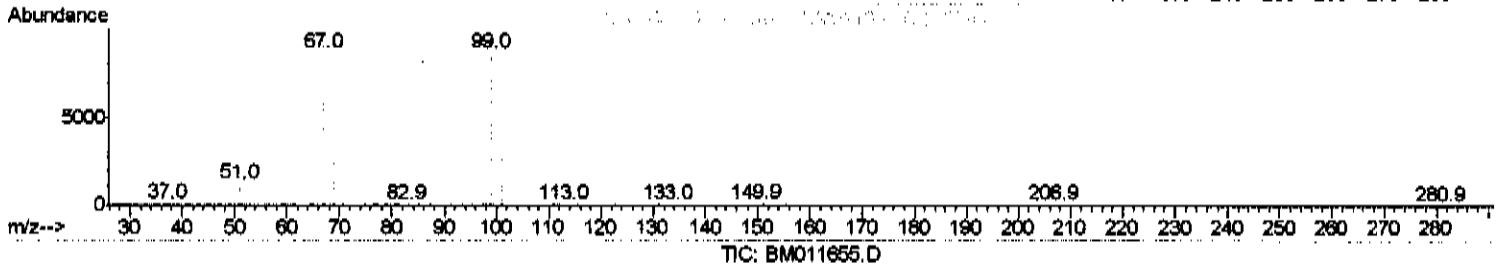
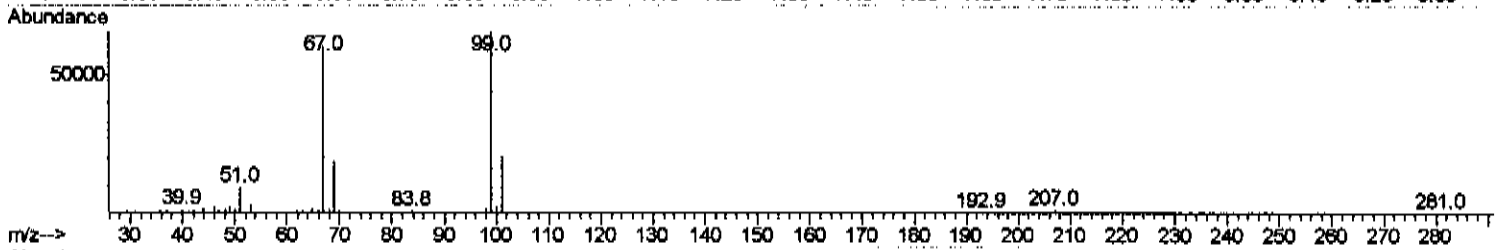
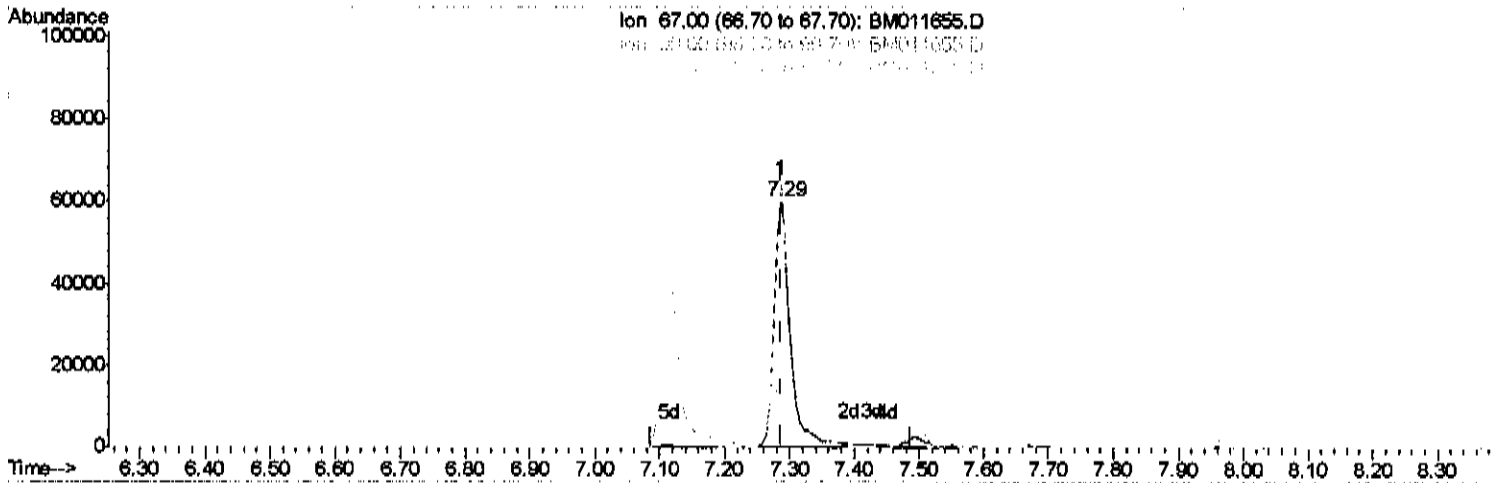
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(7) Bis-(2-Chloroethyl)ether-d8 (S)

7.287min (+0.000) 7.41ng/ul m

response 108669

Ion	Exp%	Act%
67.00	100	100
99.00	117.60	108.10
69.00	33.70	31.19
0.00	0.00	0.00

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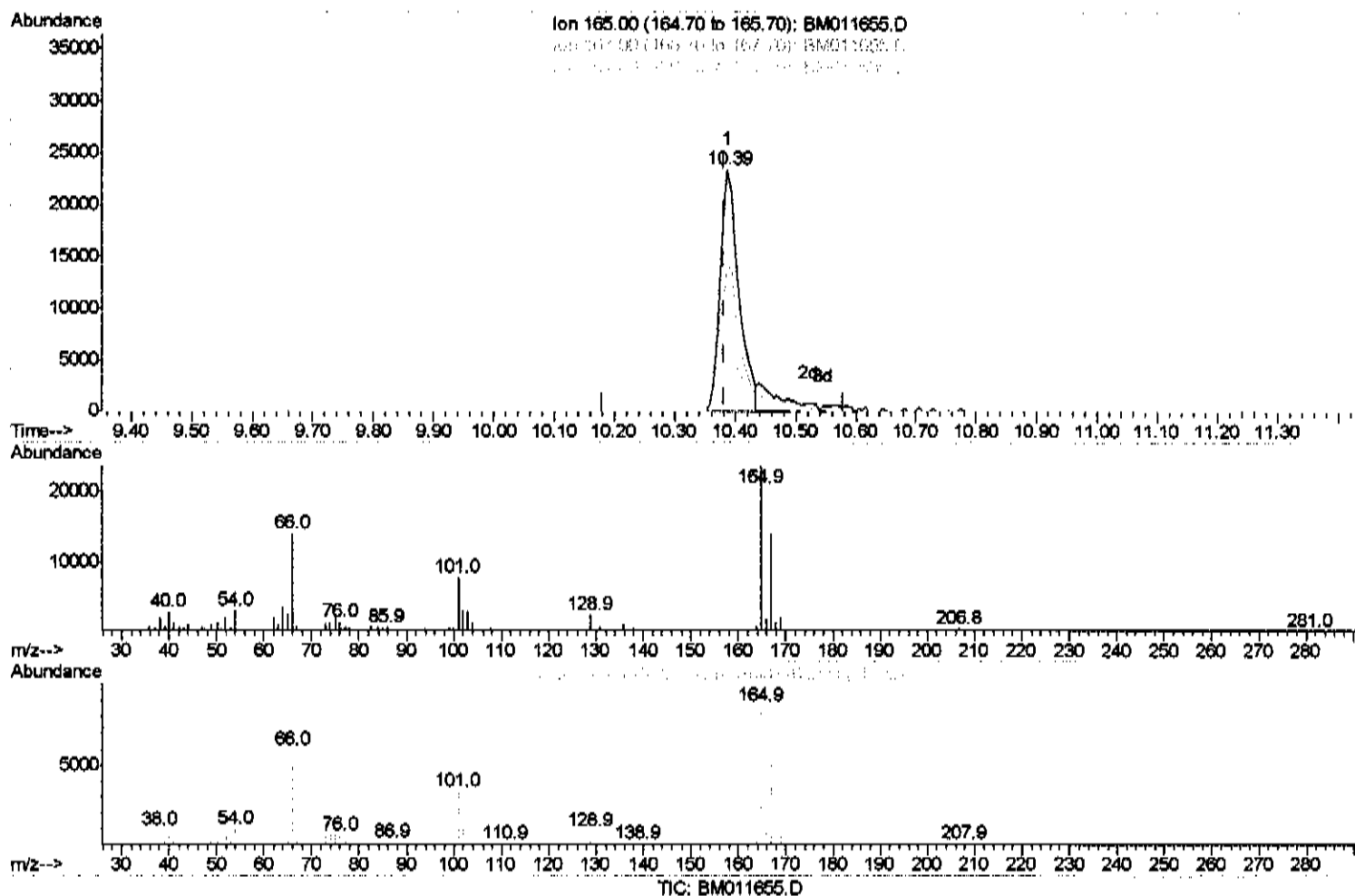
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(26) 2,4-Dichlorophenol-d3 (S)

10.386min (+0.006) 3.18ng/ul

response 50901

Ion	Exp%	Act%
165.00	100	100
167.00	69.50	60.29
101.00	31.20	33.30
0.00	0.00	0.00

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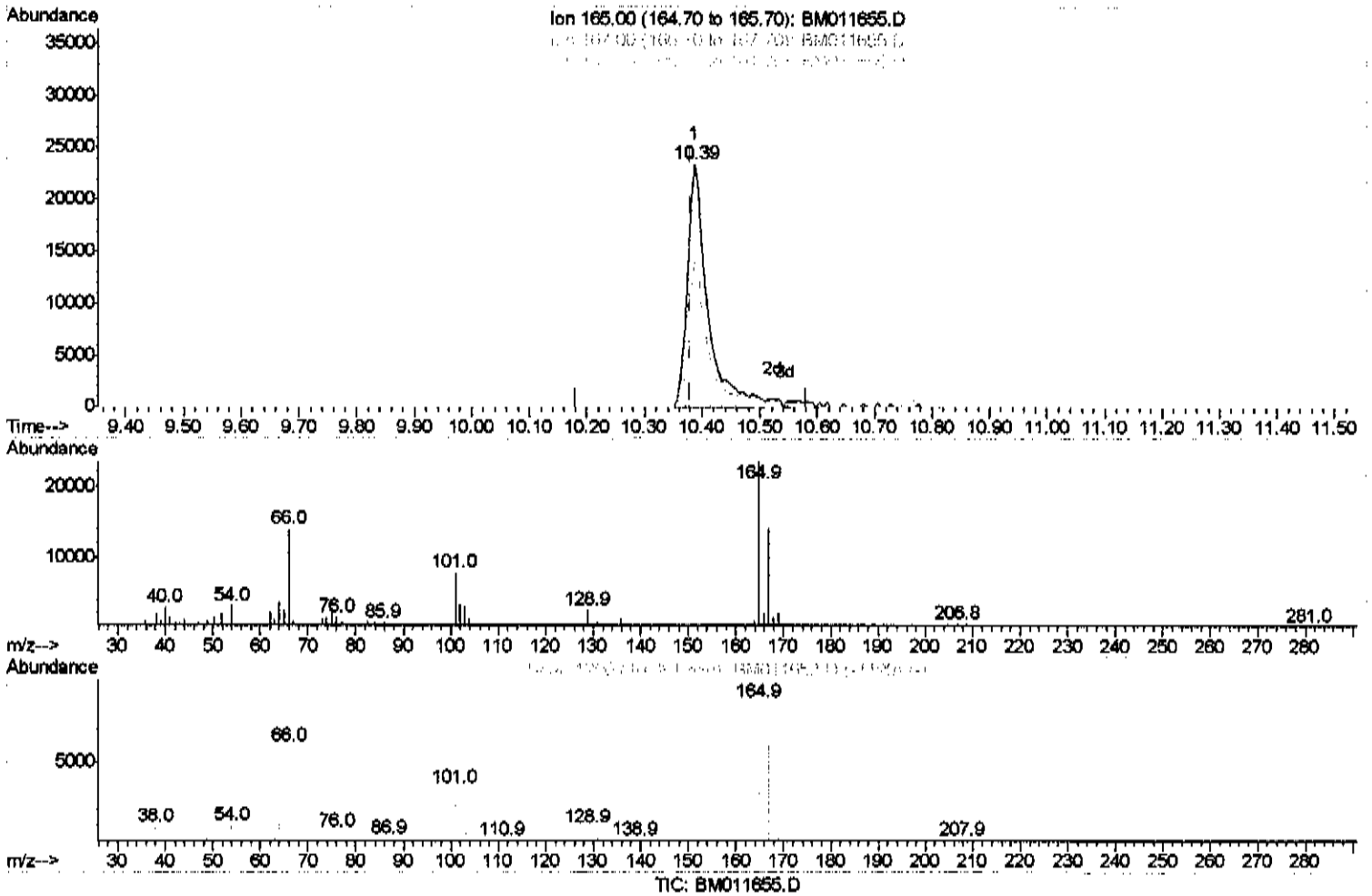
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(26) 2,4-Dichlorophenol-d3 (S)

10.388min (+0.006) 3.66ng/ul m

response 59545

Ion	Exp%	Act%
165.00	100	100
167.00	69.50	60.29
101.00	31.20	33.30
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	7.96	152	222403	20.00	ng/ul	0.00
18) Naphthalene-d8	10.77	136	1006226	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.59	164	582215	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.33	188	1306100	20.00	ng/ul	0.00
78) Chrysene-d12	21.50	240	1462402	20.00	ng/ul	0.00
86) Perylene-d12	23.87	264	1468545	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.39	96	3864	0.78	ng/uL	0.00
5) Phenol-d5	7.12	99	86141	4.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.29	67	106669m>	7.41	ng/ul>	0.00
9) 2-Chlorophenol-d4	7.49	132	77459	4.81	ng/ul	0.00
13) 4-Methylphenol-d8	8.66	113	70074	4.20	ng/ul	0.00
19) Nitrobenzene-d5	9.13	128	46757	6.12	ng/ul	0.00
22) 2-Nitrophenol-d4	9.85	143	23607	2.97	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.39	165	58545m>	3.66	ng/ul>	0.00
29) 4-Chloroaniline-d4	10.95	131	3167	0.18	ng/ul	0.05
44) Dimethylphthalate-d6	13.99	166	247090	5.02	ng/ul	0.00
47) Acenaphthylene-d8	14.29	160	376649	6.33	ng/ul	0.00
52) 4-Nitrophenol-d4	14.80	143	112	0.01	ng/ul	0.03
58) Fluorene-d10	15.58	176	307473	7.21	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.69	200	7703	1.01	ng/ul	0.00
71) Anthracene-d10	17.43	188	459312	7.14	ng/ul	0.00
79) Pvrene-d10	19.72	212	560501	8.19	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.71	264	504870	7.21	ng/ul	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	14.04	163	69270	1.470	ng/ul#	98
70) Phenanthrene	17.37	178	246316	3.383	ng/ul	100
77) Fluoranthene	19.38	202	383297	5.119	ng/ul#	90
80) Pvrene	19.74	202	356851	4.182	ng/ul#	89
83) Benzo(a)anthracene	21.48	228	181400	2.253	ng/ul	97
84) Bis(2-ethylhexyl)phthalate	21.39	149	85816	1.424	ng/ul#	93
85) Chrysene	21.53	228	179565	2.460	ng/ul	98
88) Benzo(b)fluoranthene	23.14	252	226125	2.560	ng/ul#	88
91) Benzo(a)pvrene	23.76	252	132774	1.592	ng/ul#	91

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