

Quantitation Report (QT Reviewed)

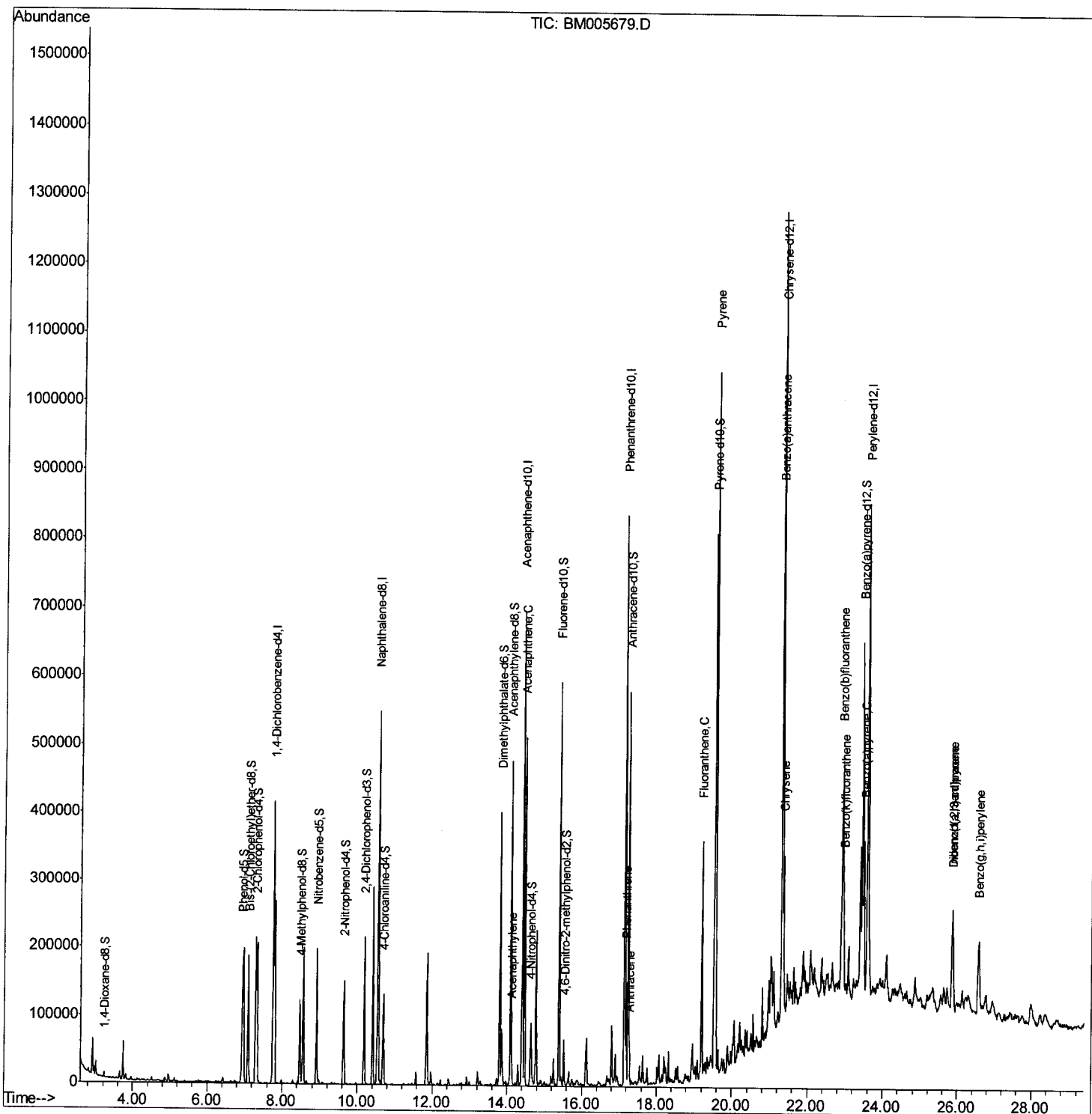
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060616\
 Data File : BM005679.D
 Acq On : 06 Jun 2016 20:51
 Operator : UM/SJ
 Sample : H3412-06MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XS5MSD

Manual Integrations
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Quant Time: Jun 07 05:02:07 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 07 04:25:20 2016
 Response via : Initial Calibration



Quantitation Report (Qedit)

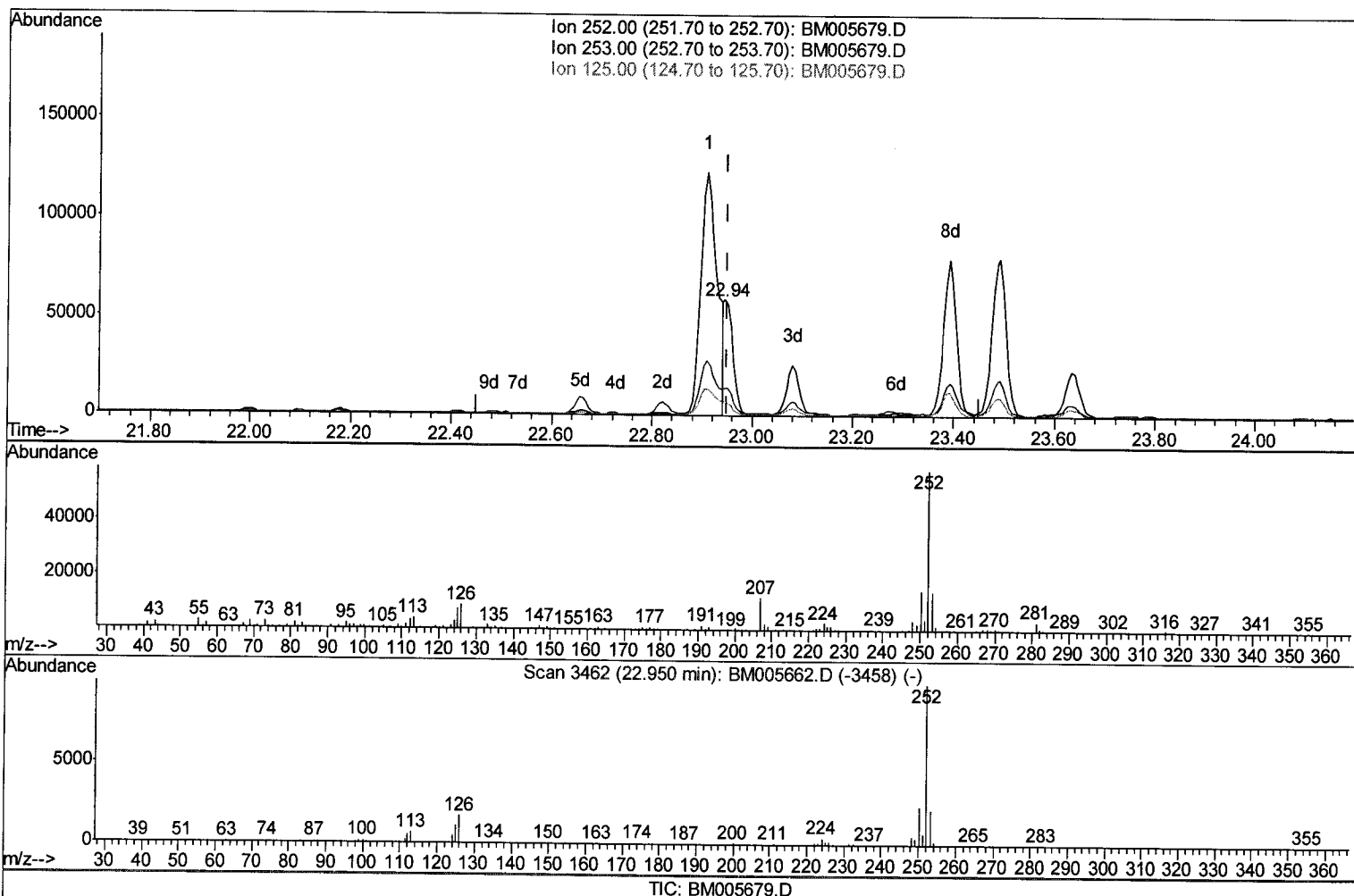
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(36) Benzo(k)fluoranthene

22.944min (-0.006) 2.72ng/ul m

response 75367

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	24.09
125.00	9.60	12.66#
0.00	0.00	0.00

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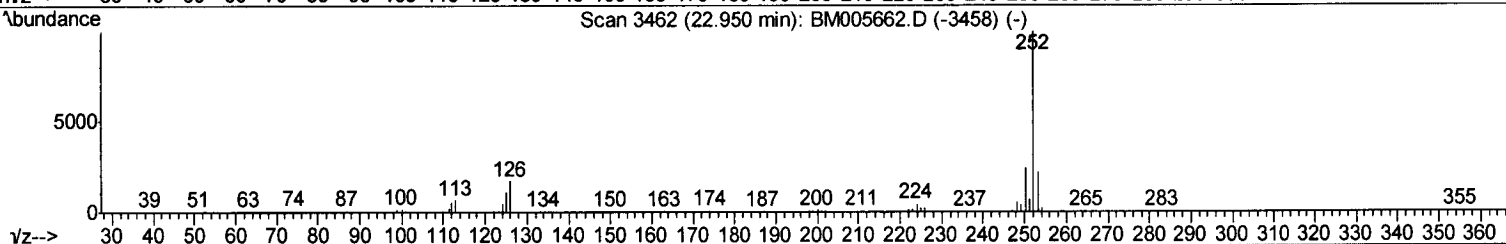
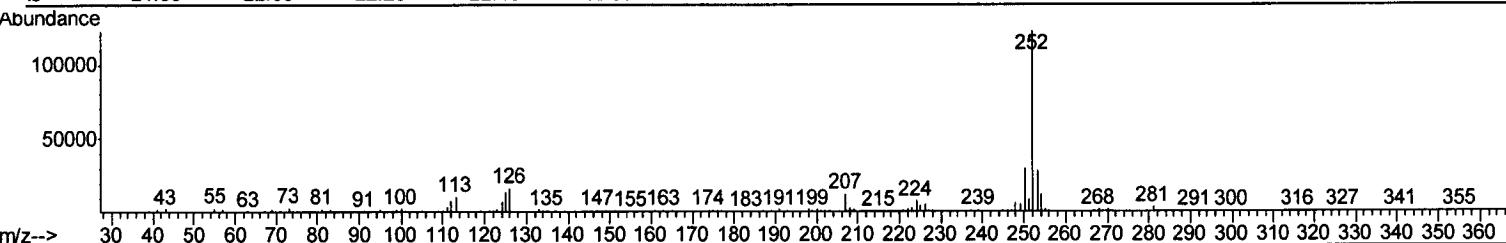
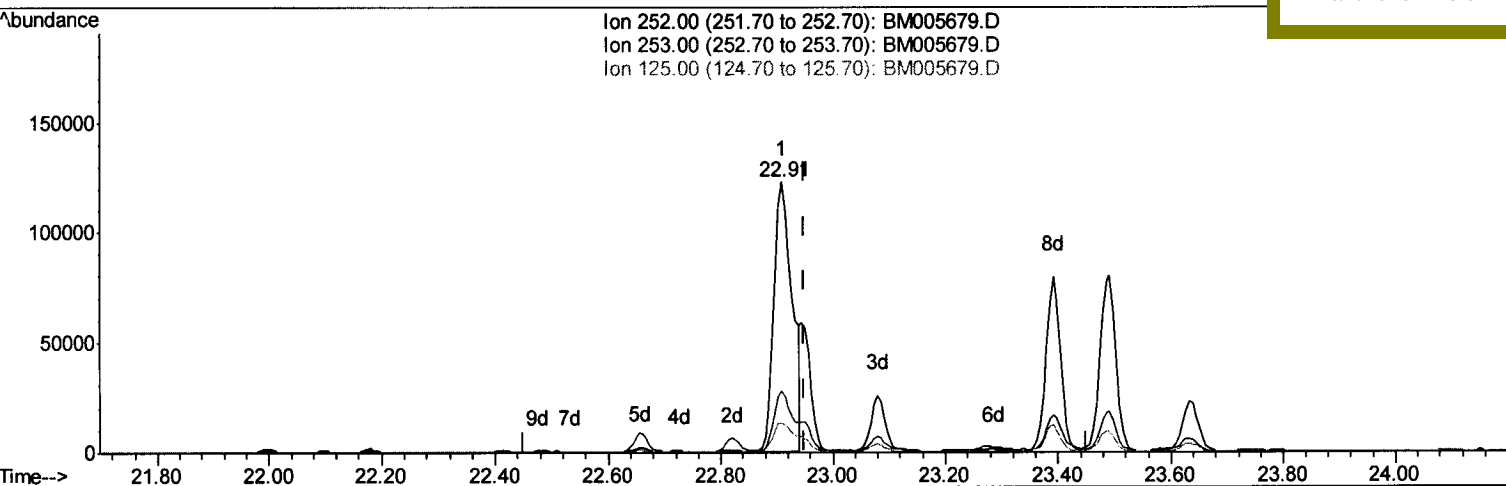
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TIC: BM005679.D

(36) Benzo(k)fluoranthene

22.909min (-0.041) 9.92ng/ul

response 274738

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	22.71
125.00	9.60	10.96
0.00	0.00	0.00

Quantitation Report (Qedit)

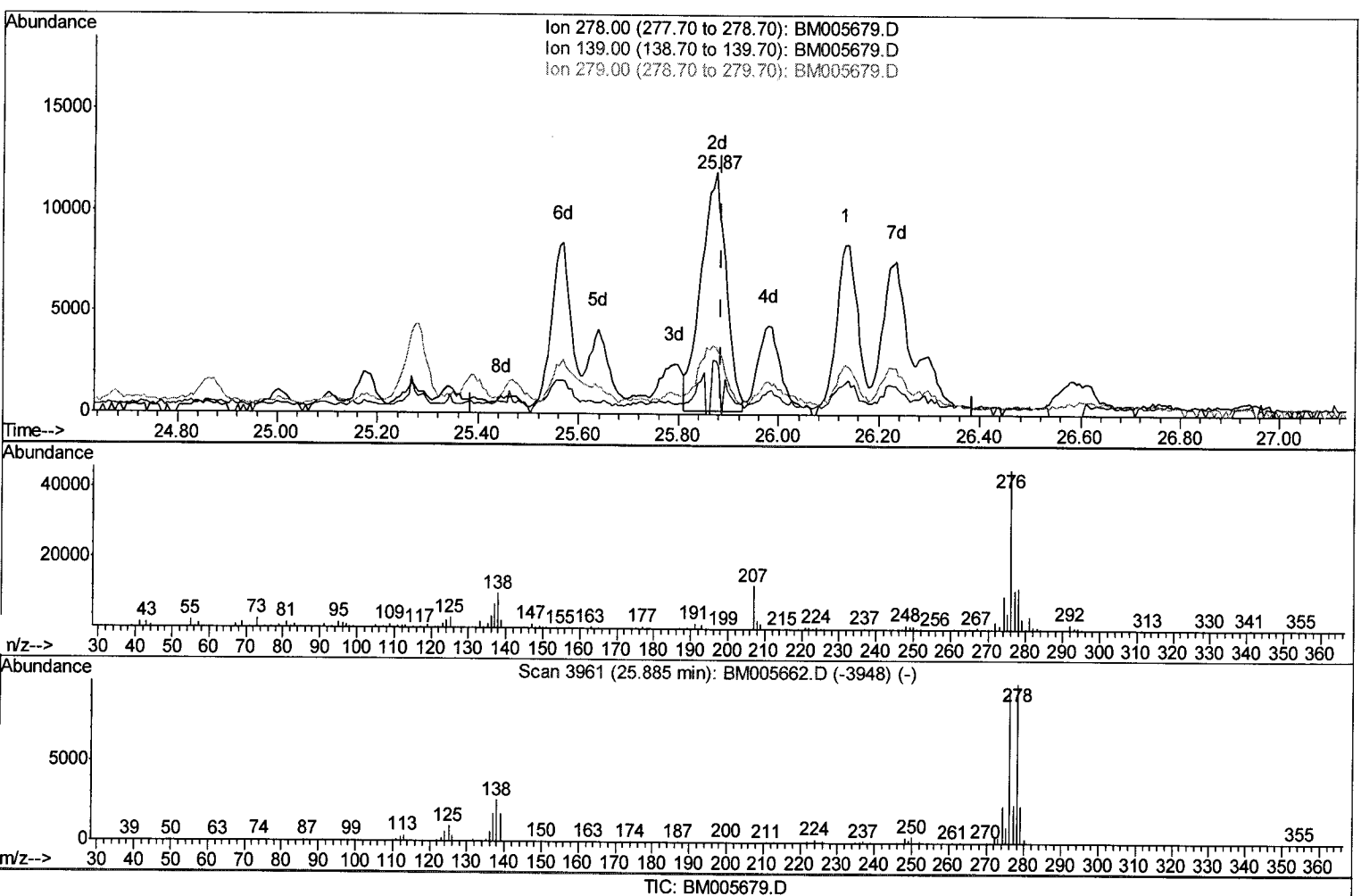
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(40) Dibenzo(a,h)anthracene

25.874min (-0.012) 1.45ng/ul m

response 40859

Ion	Exp%	Act%
278.00	100	100
139.00	15.10	21.92#
279.00	24.80	27.44
0.00	0.00	0.00

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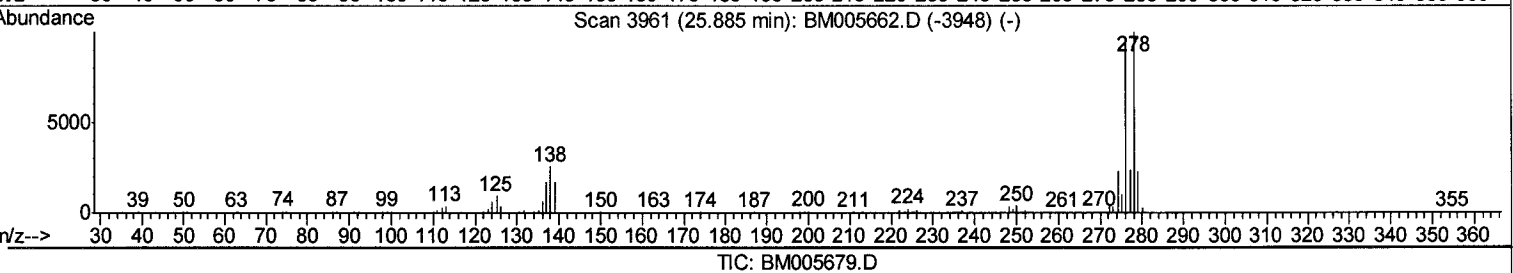
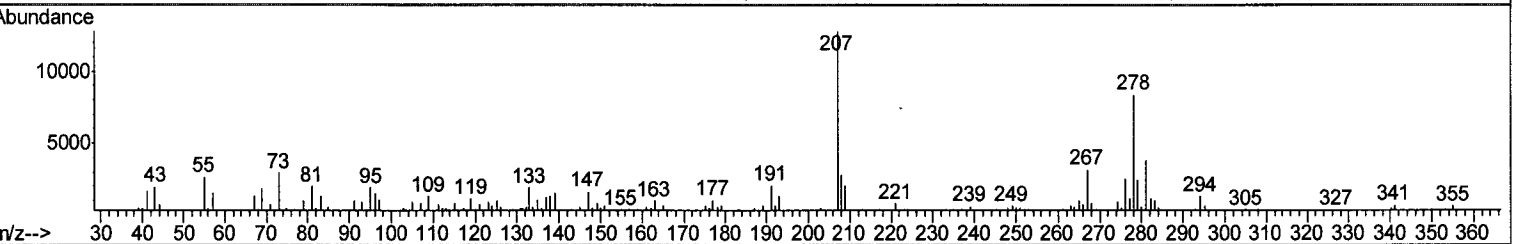
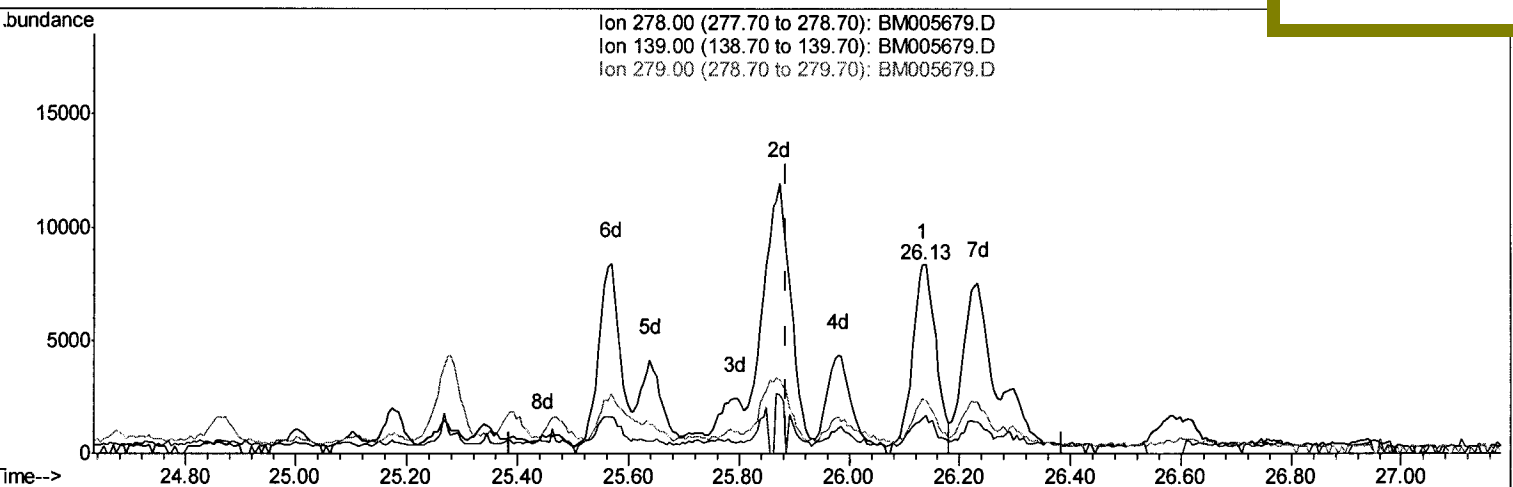
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(40) Dibenzo(a,h)anthracene

26.133min (+0.247) 0.90ng/ul

response 25241

Ion	Exp%	Act%
278.00	100	100
139.00	15.10	17.90
279.00	24.80	29.40
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	7.76	152	111119	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	459404	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	228865	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	461756	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	469823	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	489301	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.22	96	3707	1.40	ng/uL	0.00
3) Phenol-d5	6.94	99	106082	10.86	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.09	67	82434	13.87	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	100145	12.43	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	46733	6.05	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	49366	13.99	ng/ul	0.00
9) 2-Nitrophenol-d4	9.64	143	50945	13.50	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.18	165	80531	12.00	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	78662	9.87	ng/ul	0.00
16) Dimethylphthalate-d6	13.81	166	260805	14.82	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	332856	14.55	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	15476	6.14	ng/ul	0.00
21) Fluorene-d10	15.39	176	228063	15.09	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	13616	6.26	ng/ul	0.00
26) Anthracene-d10	17.24	188	320504	14.62	ng/ul	0.00
30) Pyrene-d10	19.54	212	343513	14.55	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	333136	14.79	ng/ul	0.00

Target Compounds

						Qvalue
18) Acenaphthylene	14.12	152	29770	1.26	ng/ul	97
19) Acenaphthene	14.46	153	209227	13.61	ng/ul	94
25) Phenanthrene	17.19	178	89258	3.50	ng/ul	98
27) Anthracene	17.28	178	28093	1.08	ng/ul	99
28) Fluoranthene	19.20	202	194412	7.46	ng/ul	99
31) Pyrene	19.57	202	580248	19.57	ng/ul	97
32) Benzo(a)anthracene	21.31	228	121739	4.42	ng/ul	96
33) Chrysene	21.36	228	144811	5.56	ng/ul	96
35) Benzo(b)fluoranthene	22.91	252	274738	9.31	ng/ul	99
36) Benzo(k)fluoranthene	22.94	252	75367m	2.72	ng/ul	
38) Benzo(a)pyrene	23.49	252	153373	5.40	ng/ul	97
39) Indeno(1,2,3-cd)pyrene	25.87	276	130276	3.88	ng/ul	96
40) Dibenzo(a,h)anthracene	25.87	278	40859m	1.45	ng/ul	
41) Benzo(g,h,i)perylene	26.57	276	124649	4.42	ng/ul#	93

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