

Quantitation Report (QT Reviewed)

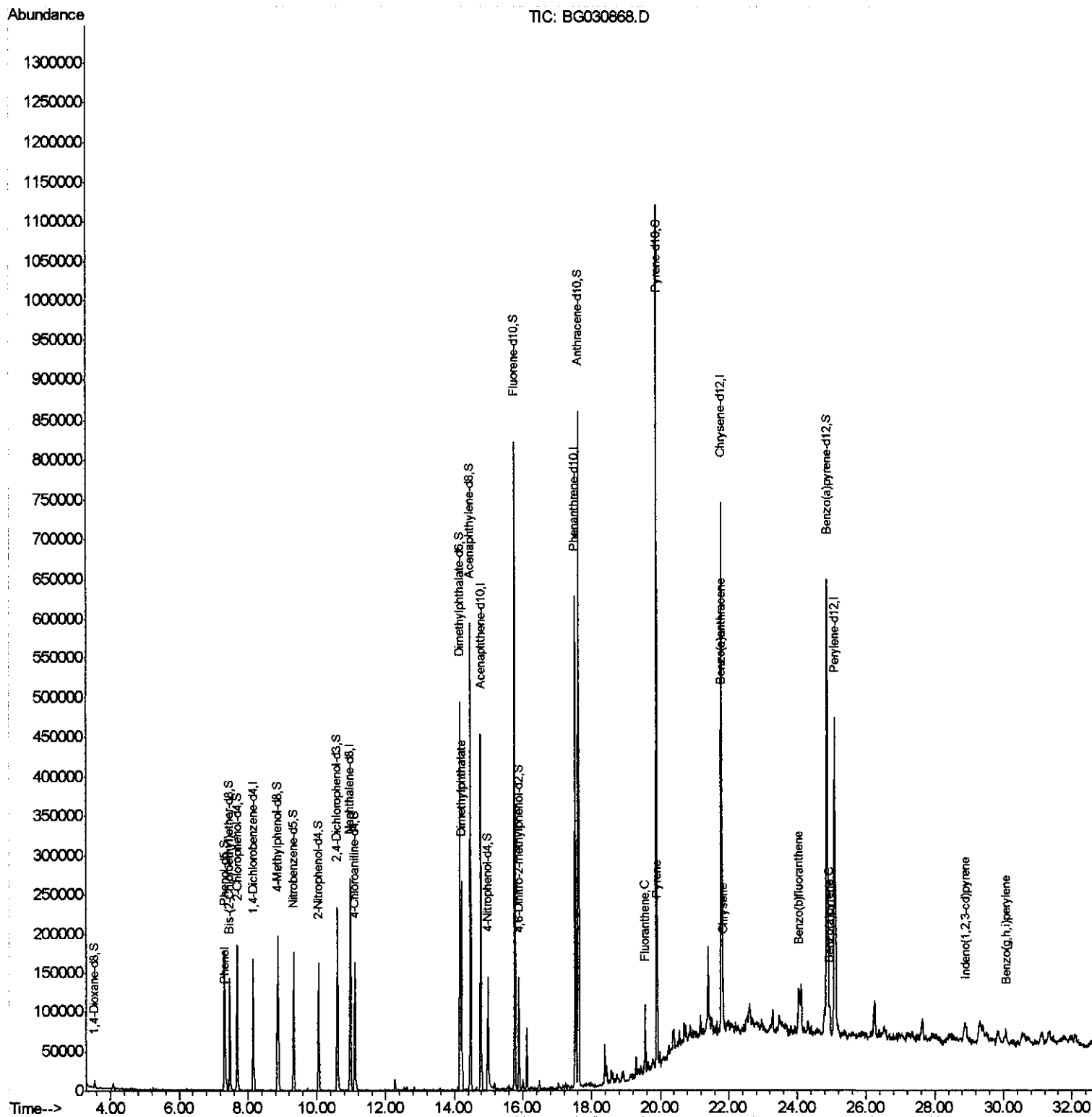
Data Path : Z:\HPCHEM1\BNA G\DATA\BG110817\
 Data File : BG030868.D
 Acq On : 9 Nov 2017 5:06
 Operator : SJ/JU
 Sample : I6245-15
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B0F13

Manual Integrations
 APPROVED

Sohil
 11/9/2017 3:46:36 PM

Quant Time: Nov 09 08:08:41 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG101417.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 09 07:08:58 2017
 Response via : Initial Calibration



Quantitation Report (Qedit)

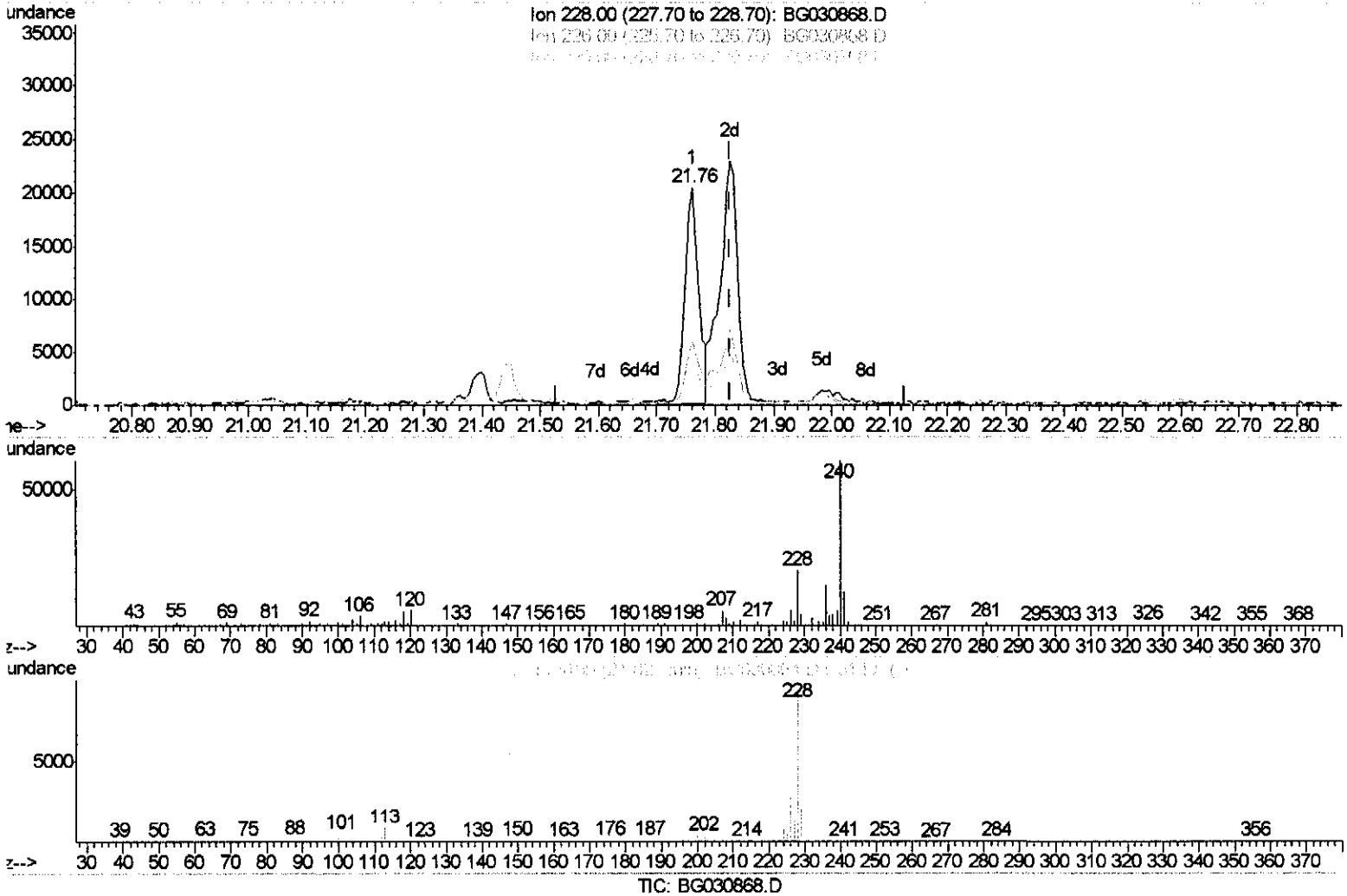
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Quant Time: Nov 09 07:42:00 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG101417.M
 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration



(82) Chrysene

21.760min (-0.066) 1.47ng/ul

response 34415

Ion	Exp%	Act%
228.00	100	100
226.00	30.20	29.43
229.00	19.00	21.48
0.00	0.00	0.00

Quantitation Report (Qedit)

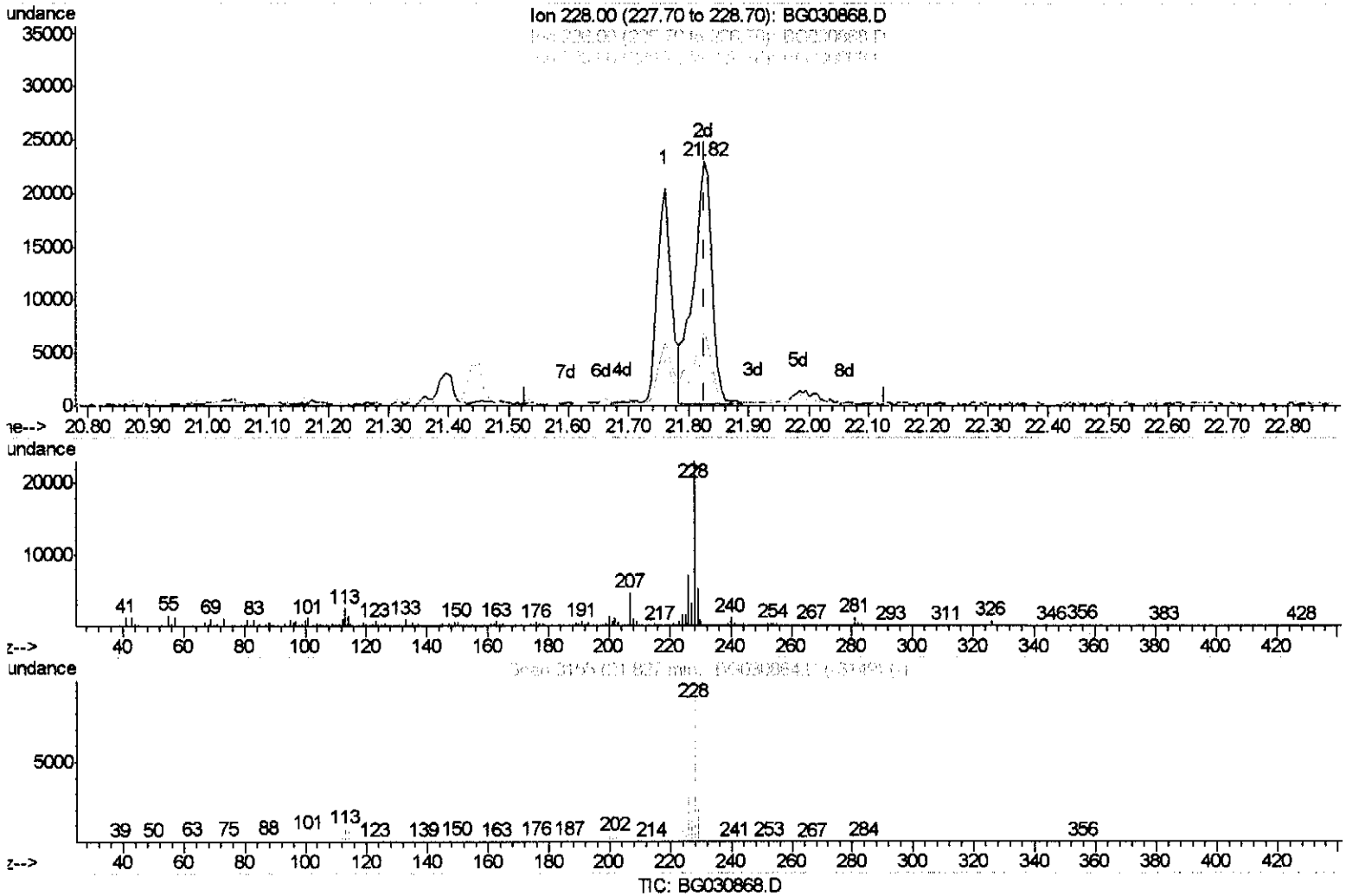
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(82) Chrysene

21.825min (-0.002) 2.09ng/ul m 50 11/09/17
 response 49127

Ion	Exp%	Act%
228.00	100	100
226.00	30.20	31.49
229.00	19.00	23.25#
0.00	0.00	0.00

Quantitation Report (Qedit)

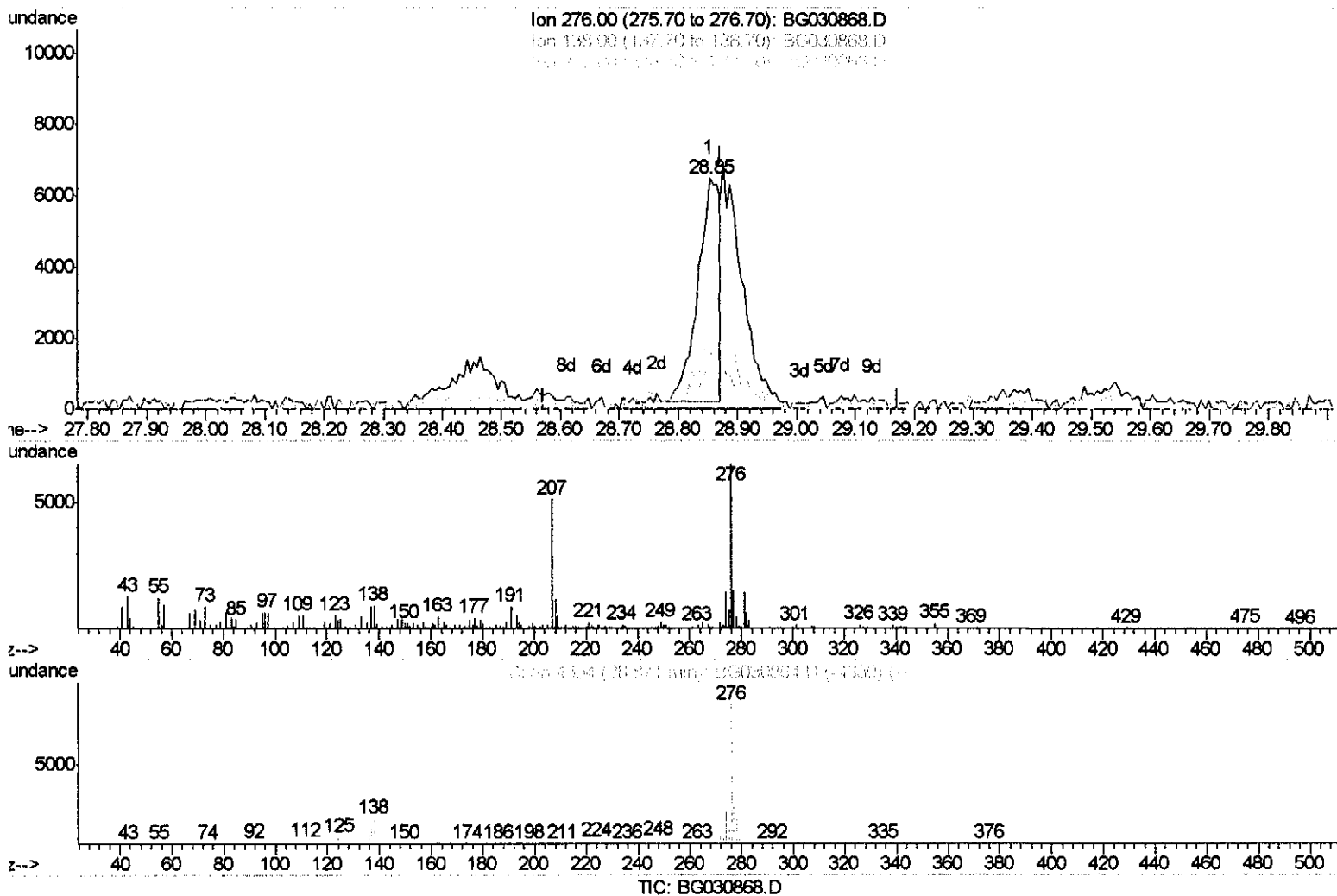
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(89) Indeno(1,2,3-cd)pyrene

28.852min (-0.019) 0.57ng/ul

response 16166

Ion	Exp%	Act%
276.00	100	100
138.00	16.70	15.92
277.00	23.30	25.37
0.00	0.00	0.00

Quantitation Report (Qedit)

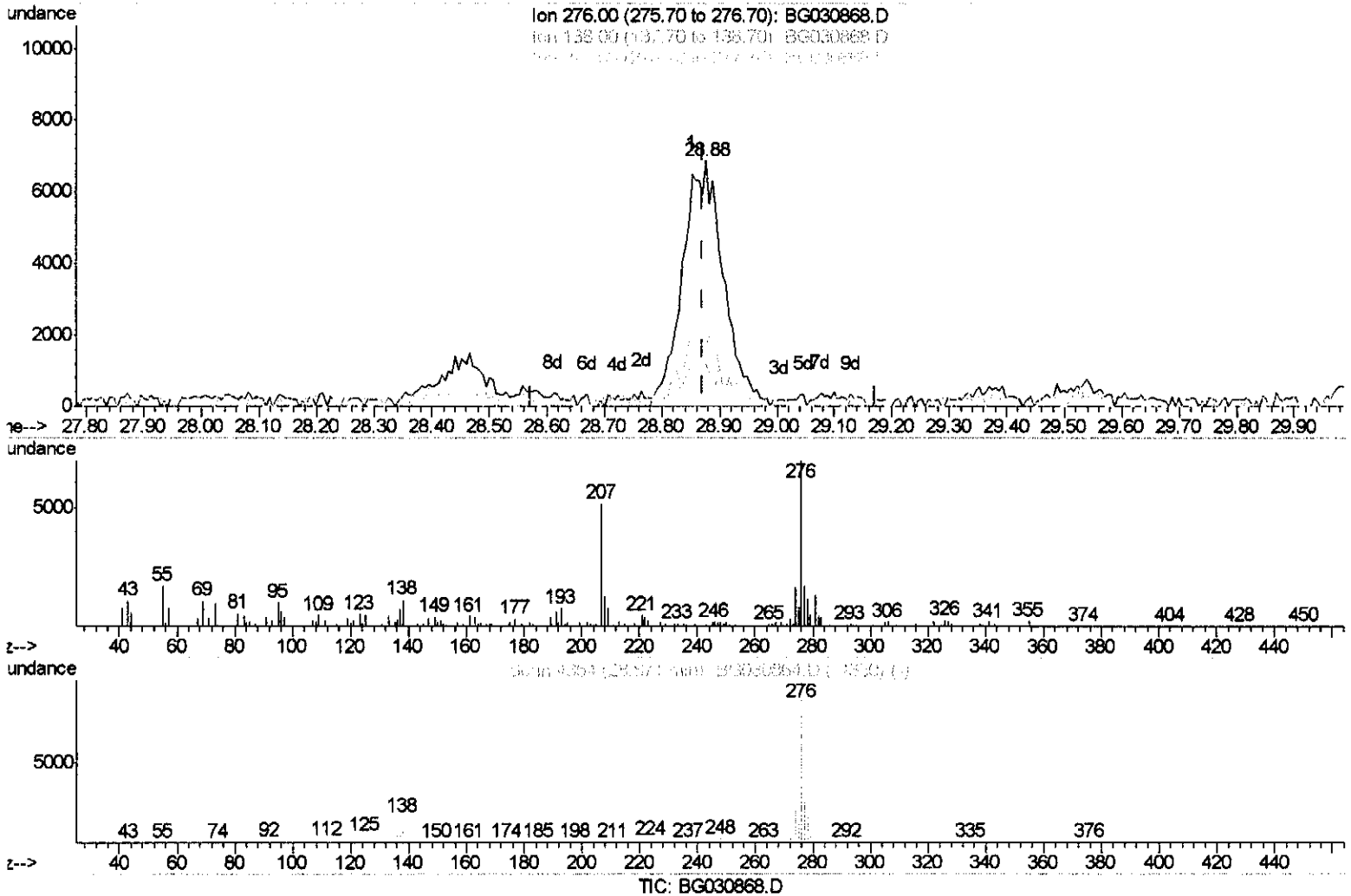
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 Operator : SJ/JU
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(89) Indeno(1,2,3-cd)pyrene

28.876min (+0.004) 1.19ng/ul m SJ 11/09/17

response 33829

Ion	Exp%	Act%
276.00	100	100
138.00	16.70	17.78
277.00	23.30	25.44
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	8.15	152	50551	20.00	ng/ul	0.00
18) Naphthalene-d8	10.97	136	248638	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.77	164	163365	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.51	188	389638	20.00	ng/ul	0.00
75) Chrysene-d12	21.78	240	412014	20.00	ng/ul	0.00
83) Perylene-d12	25.09	264	428871	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.52	96	5517	4.44	ng/uL	0.00
5) Phenol-d5	7.31	99	115773	23.86	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.47	67	67842	22.34	ng/ul	0.00
9) 2-Chlorophenol-d4	7.68	132	94915	27.73	ng/ul	0.00
13) 4-Methylphenol-d8	8.86	113	87420	22.31	ng/ul	0.00
19) Nitrobenzene-d5	9.32	128	50830	28.48	ng/ul	0.00
22) 2-Nitrophenol-d4	10.04	143	63253	34.59	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.59	165	108993	26.15	ng/ul	0.00
29) 4-Chloroaniline-d4	11.10	131	104259	21.53	ng/ul	0.00
43) Dimethylphthalate-d6	14.16	166	370674	26.54	ng/ul	0.00
46) Acenaphthylene-d8	14.46	160	460417	27.11	ng/ul	0.00
51) 4-Nitrophenol-d4	14.98	143	50366	19.53	ng/ul	0.00
57) Fluorene-d10	15.76	176	342994	29.99	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.88	200	40945	21.53	ng/ul	0.00
70) Anthracene-d10	17.61	188	531830	28.86	ng/ul	0.00
76) Pyrene-d10	19.88	212	609762	28.59	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.86	264	604429	29.76	ng/ul	0.00

Target Compounds

						Ovalue
6) Phenol	7.34	94	23607	4.70	ng/ul	90
44) Dimethylphthalate	14.21	163	190219	13.96	ng/ul	99
74) Fluoranthene	19.55	202	52894	2.25	ng/ul	99
77) Pyrene	19.91	202	61667	2.23	ng/ul	97
80) Benzo(a)anthracene	21.76	228	34338	1.33	ng/ul	97
82) Chrysene	21.82	228	49127m	2.09	ng/ul	97
85) Benzo(b)fluoranthene	24.03	252	70669	2.74	ng/ul	97
88) Benzo(a)pyrene	24.93	252	37206	1.48	ng/ul	99
89) Indeno(1,2,3-cd)pyrene	28.88	276	33829m	1.19	ng/ul	-
91) Benzo(g,h,i)perylene	30.06	276	28852	1.23	ng/ul#	94

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

SJ 11/09/17
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