

TITLE PAGE

20. Selten vorkommende Stoffe im  
Steinkohlenteer.  
Compounds rarely occurring  
in bituminous coal tar.

Frame Nos. 339 - 350

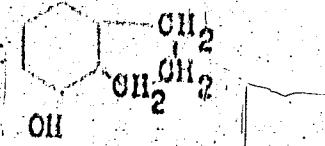
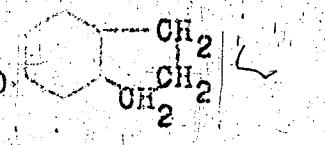
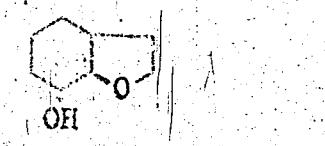
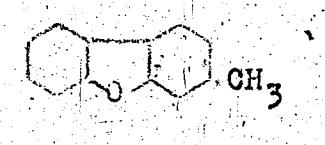
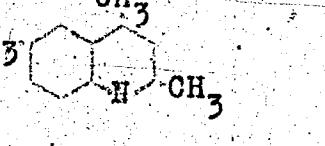
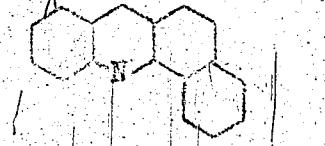
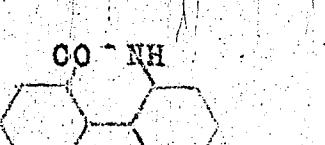
(20)

Aus: Kruber, Über die Ergebnisse der Steinkohlen-  
forschung der letzten 30 Jahre.

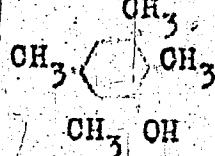
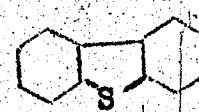
Angew.Chem. 53, 69.

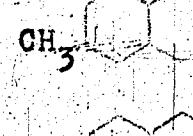
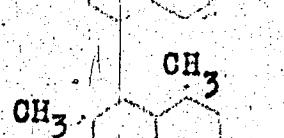
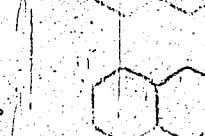
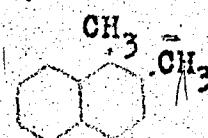
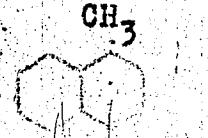
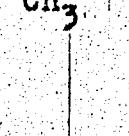
Nach diesem Aufsatz kommen im Steinkohlenoel außer Verbindungen, die bereits in der Arbeit von Fisher, Fuel in Science and Practice, aufgezählt sind, noch folgende Stoffe vor:

Name	Formel	Schmelz-Punkt	Siede-Punkt
Dicyklopentadien		33°	170°
1,5-Dimethylnaphthalin		82°	
Trimethylnaphthaline			
2,3,6-Trimethyl-naphthalin		102°	286°
1,3,7-Trimethyl-naphthalin		13,5°	280°
1-Methylphenanthren		119°	355°
3-Methylphenanthren		65°	350°
9-Methylphenanthren		92°	354°

Name	Formel	Schmelz-Punkt	Siede-Punkt
4-Oxy-hydriden		50°	245°
5-Oxy-hydriden		55°	251°
7-Oxy-cumaron		43°	240°
2-Methyl-diphenylenoxyd		66°	304°
2,4,6-Trimethyl-chinolin		50°	288°
3,4-Benzaoridin		107°	433°
Phensanthridon		286°	435°
Fluoren-nitrile			
Methyl-thionaphthene		52° und 36°	

Isoliert wurden die Verbindungen aus folgenden Fraktionen:

Name	Formel
<u>Leichtöl (Motorenbenzol):</u>	
Offene und cyclische Olefine	
Dihydrobenzol	
Tetrahydrobenzol	
n-Hexylen	C <sub>6</sub> H <sub>12</sub>
n-Heptylen	C <sub>7</sub> H <sub>14</sub>
n-Propionsäure	CH <sub>3</sub> ·CH <sub>2</sub> ·CO <sub>2</sub> H
<u>Benzolvorlauf</u>	
Aethylmercaptan	CH <sub>3</sub> ·CH <sub>2</sub> ·SH
<u>Neutrale Schwerölfaktion (ca. 250°):</u>	
Indol (4-5 %)	
Durenol	
<u>Rohnaphthalin des Steinkohlenteers:</u>	
Thionaphthen (4 %)	
<u>Methylnaphthalin-Faktion:</u>	
2 verschiedene Methyl-thionauethene	
<u>Rohphenanthren:</u>	
Diphenylensulfid (3-5 %)	

Name	Formel
<u>Teerölfaktion 240-270°</u>	
$\alpha$ -Methylnaphthalin	
$\beta$ -Methylnaphthalin	
$\gamma$ -Methylnaphthalin	
1.6-Dimethylnaphthalin	
2.6-Dimethylnaphthalin	
2.7-Dimethylnaphthalin	
1.7-Dimethylnaphthalin	
2.3-Dimethylnaphthalin	
1.2-Dimethylnaphthalin	
1.5-Dimethylnaphthalin	
Phenole (1,0-1,5 %)	
Benzoesäure	

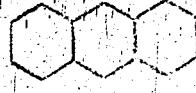
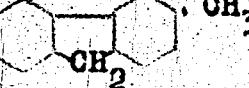
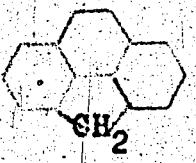
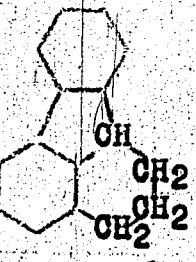
Name	Formel
7-Oxy-cumaron	
4-Oxy-hydrinden	
5-Oxy-hydrinden	
	240-265°:
8-Methyl-chinolin	
3-Methyl-isochinolin	
1-Methyl-isochinolin	
2,8-Dimethyl-chinolin	
7-Methyl-chinolin	
6-Methyl-chinolin	
3-Methyl-chinolin	
1,3-Dimethyl-isochi-nolin	

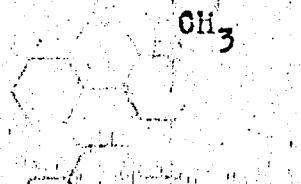
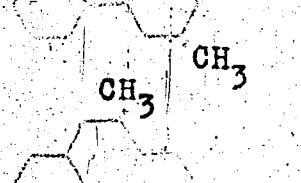
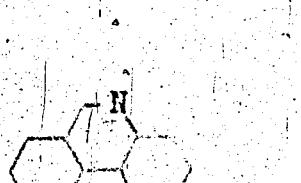
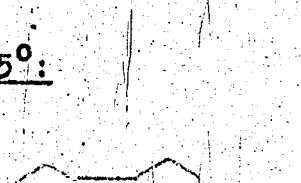
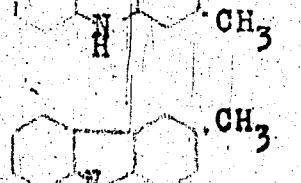
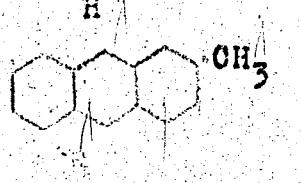
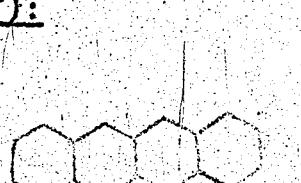
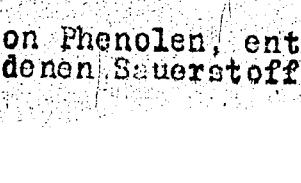
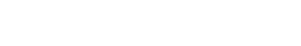
Name	Formel
5-Methyl-chinolin	
4-Methyl-chinolin	
<u>270-290°:</u>	
m-Methyl-diphenyl	
p-Methyl-diphenyl	
4.4'-Dimethyl-diphenyl	
3.4'-Dimethyl-diphenyl	
<u>280-290°:</u>	
Dimethylnaphthaline + Athylnaphthalin (1,0-1,2 % des Gesamtteers)	
Trimethylnaphthaline (0,1% des Gesamtteers)	
2.3.6-Trimethylnaphthalin	
1.3.7-Trimethylnaphthalin	

Name	Formel
	ca. 300°:
primäre Basen (8-10 %)	
$\alpha$ -Naphthylamin	
$\beta$ -Naphthylamin	
2.4.6-Trimethylchinoxalin	
4.5-Benzocindan	
	bis 300° <sup>1)</sup> :
1-Methyl-diphenylenoxyd	
2-Methyl-diphenylenoxyd	
$\alpha$ -Naphthonitril	
$\beta$ -Naphthonitril	
Xylenolfraktion:	
1.3.5-Xylenol	
1.2.6-Xylenol	

<sup>1)</sup> Enthält über 4% O und 0,3% N.

Name	Formel
1.2.3-Xylenol	
m-Äthylphenol	
p-Äthylphenol	
<u>Cumenolfraktion 230°:</u>	
Isopseudocumanol	
sym. m-Methyl-äthylphenol	
<u>250°:</u>	
sym. Hemellitol	
Durenol	
<u>Anthracenöl:</u>	
Phenole 6 %	
4-Oxy-diphenyl	
2-Oxy-diphenylenoxyd	
2-Oxy-fluoren	
Phenanthrol-(2)	

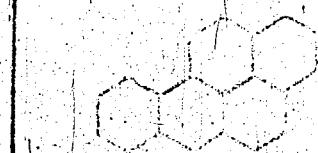
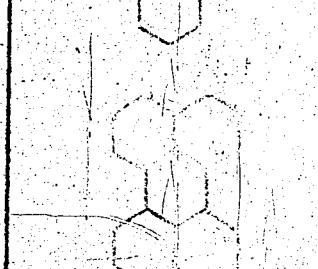
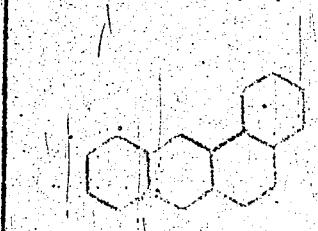
Name	Formel
Phenanthren	
Antracen	
Carbazol	
Fluoren	
Diphenylenoxyd	
	<u>302°:</u>
2-Methylfluoren	
3-Methylfluoren	
Nonadecan	$C_{19}H_{40}$
	<u>Neutrale Anthracenölfraktion 350-360°:</u>
4.5-Methylenphenanthren	
	<u>360-370°:</u>
1.2.3.4-Tetrahydronaphthalen	
2-Phenylnaphthalin	

Name	Formel
1-Methyl-phenanthren	
3-Methyl-phenanthren	
9-Methyl-phenanthren	
Diphenylensulfid	
Fluoren-nitril	
Phenanthridin	
<u>365-375°:</u>	
2-Methyl-carbazol	
3-Methyl-carbazol	
5-Methyl-anthracen	
Naphthacen	

Pech <sup>1)</sup>:

1) Pech ist praktisch frei von Phenolen, enthält aber 1,5-2 % anderweitig gebundenen Sauerstoff = 30 % O-haltige Verbindungen.

Name	Formel
1.2-Benz-carbazol	
3.3'-Naphthylen-phenyleneoxyd	
1.9-Benzoxanthen	
Chrysen	
Triphenylen (1-3 % im Roh- chrysen)	
3.4-Benzaoridin (Chrysenfraktion)	
Phenanthridon	

Name	Formel
	<u>über 440°:</u>
3.4-Benzpyren	
1.2-Benzpyren	
Perylen	
1,2-Benzanthracen	
1,2-Benzo-fluoren	
2.3-Benzo-fluoren	