# **Economic Commission for Europe**

## **Inland Transport Committee**

**Working Party on the Transport of Dangerous Goods** 

Joint Meeting of Experts on the Regulations annexed to the European Agreement concerning the International Carriage of Dangerous Goods by Inland Waterways (ADN) (ADN Safety Committee)

Thirtieth session

Geneva, 23–27 January 2017

Item 5 (b) of the provisional agenda

Proposals for amendments to the Regulations annexed to ADN: other proposals

English

23 January 2017

## **Autonomous protective systems**

#### Transmitted by EBU, ESO and ERSTU

### **Initial situation**

Numerous documents were submitted for the twenty-ninth session of the safety committee concerning the requirements governing explosion-potential groups for non-electrical equipment.

Because it was not possible for reasons of time to amend the ADN 2017, the safety committee agreed during the initial phase to settle problems by means of multilateral agreements.

The inland navigation industry is grateful for this. However, multilateral agreement ADN / M 018 imminently requires measures necessitating careful preparation in respect of those ships whose certificate of approval need to be renewed after 31st December 2018.

Consequently it is necessary to address a number of the questions raised by the shipping industry in a timely manner.

## Question

The minutes of the twenty-ninth session of the safety committee ADN/WP.15/AC.2/60 noted under section 44 that the informal "substances" working group has now been requested to look into various facts.

What progress have these investigations made?

## **Enquiry concerning the minutes of the twenty-ninth session**

The minutes of the twenty-ninth session of the safety committee ADN/WP.15/AC.2/60 noted under section 44 – second sub-point – that the industry is required to provide the substances working group with relevant information. This wording in the minutes requires clarification.

Without doubt the <u>inland navigation</u> industry has the greatest interest in a systematic review and, if applicable, provision of the entries in column 16 of table C of the ADN.

The <u>inland navigation</u> industry never acquires title to the cargo. The inland navigation industry does not therefore see itself as being in a position to commission the experimental determination of explosion-potential groups or subgroups based on cargo samples.

Moreover, in the case of N.O.S entries, the inland navigation industry typically does not possess the necessary information that would allow the required explosion-potential groups or subgroups to be determined.

The <u>shipping</u> industry as well is affected by findings in respect of explosion-potential groups or subgroups. The <u>shipping</u> industry is far better informed about the characteristics of the goods to be shipped than the inland navigation industry.

The inland navigation associations therefore ask the safety committee to clarify exactly how it sees the information in question being provided "by the industry".

## **Determination of the explosion-potential group or subgroups**

The inland navigation industry provided the twenty-ninth session of the safety committee with an indication of a theoretical approach to determining explosion-potential groups or subgroups in document INF. 21. A calculation for UN 1170 ethanol was submitted by way of an example.

Given the complexity of the substance, the inland navigation industry appreciates that this aspect cannot be addressed in detail at the twenty-ninth session.

The inland navigation industry is now proposing to undertake a systematic review of explosion-potential groups or subgroups in accordance with the theoretical approach in standards EN-IEC 60079-20-1 and ISO/IEC DIS 80079-20-1.

Following the shipping industry's investigations it should be emphasised that the N.O.S. items in UN 1268 and 3295, of which many millions of tonnes of cargo are transported, are to be prioritised for investigation and probably reclassified as explosion-potential group IIA.

As a prelude to this discussion, the inland navigation industry has already conducted extensive comparative calculations that frequently reveal that a "lighter" explosion-potential group or subgroups is sufficient instead of explosion-potential group IIIB.

These results are summarised in the accompanying table.

Owing to the importance of this matter, the inland navigation industry has already signalled that additional industry experts would like to participate in the informal "substances" working group's future deliberations.

	ESG	FORMULA for I	MIXTUR	ES			ording to ISO-EIC-DIS 80079-20-1
G	roup IIA: MESG ≥ 0,9 mm.					a = classified	according to MESG determination.
G	roup IIB: 0.5 mm < MESG< 0.9 mm.	MESC <sub>1/</sub>				b = classified	according to MIC ratio.
	roup IIC: MESG ≤ 0.5 mm.						•
	oup no. MESO'S 0,0 mm.	$\Sigma_{l}(X_{l}/M)$	ESG <sub>I</sub> )				G and MIC ratio have been determined.
M	IC						according to similarity of chemical structure (provisional
G	roup IIA: MIC > 0,8.					classification).	
	roup IIB: 0,45 ≤ MIC ≤ 0,8.					Classificatie v	an gas mengsels
	roup IIC: MIC < 0,45.						
	UBĠROUP IIB					###	According to the formula in ISO-EIC-DIS 80079-20-1 for mixures and the mixture doesn't contain
	roup IIB: MESG > 0,5 mm.						components for what IIB of IIC is described or the mixture contains less 30% of these components, IIB3 complies
G	roup IIB3: MESG ≥ 0,65 mm.						
	roup IIB2: MESG ≥ 0,75 mm.						
G	roup IIB1: MESG ≥ 0,85 mm.						
G	roup IIA: MESG ≥ 0,9 mm.	Norm IS	O-EIC-L	IS 80079-20-1		according to	
U	N Naam	MESG	MIC	method.	ADN 2015	this norm	
	1088 ACETAL	0.8	14	d			Circles to 4.0 Circles with any
		0,0		0.98 a	IIB <sup>4</sup>	IIB2 IIA	Simular to 1,2 Diethoxyethane
	1089 ACETALDEHYDE (ETHANAL)						
	1090 ACETONE 1092 ACROLEINE	1,0 0,7	77	1 c	IIA IIB	IIA IIB3	(ADN WP.15/AC.2/2016/4 says IIB3)
	1093 ACRYLONITRILE	0.1		0,78 c	IIB	IIB3	(ADN WP.15/AC.2/2016/4 says IIBS)
	1098 ALLYL ALCOHOL	0,0		U,78 C	IIB	IIB2	(ADN WP.15/AC.2/2016/4 says IIB2)
	1100 ALLYI CHIORIDE	1.1	17	1.33 a	IIA	IIA	(ADIV WF. 15/AO.2/2016/4 Says libb)
	1105 PENTANOLS (n-PENTANOL)	0.0		1,35 a	IIA	IIA	
	1106 AMYLAMINE (n-AMYLAMINE)	U,i	98	d	IIA <sup>7</sup>	IIA	all amines products MESG_IIA.
	1107 AMYLCHLORIDES				ΙΆ	IIA	all AMYL CHLORIDES MESG IIA,
	1108 1-PENTENE (n-AMYLENE)	0.0	16	а		IIA	according to AMYLENE IIA
	1114 BENZENE	0.0			IIB <sup>4</sup>	IIA	according to AMYLENE IIA
	1120 BUTANOLS (n-BUTYL ALCOHOL)	0,0		a	IIB	IIA	(ADN WP.15/AC.2/2016/4 says IIB2)
		U,i	91				
	1120 BUTANOLS (sec BUTYLALCOHOL)			d says IIA,	IIB <sup>7</sup>	IIA	(ADN WP.15/AC.2/2016/4 says IIB2)
	1120 BUTANOLS (tert BUTYLALCOHOL)				IIA <sup>7</sup>	IIA	(ADN WP.15/AC.2/2016/4 says IIB2)
	1123 BUTYLACETATES (n-BUTYL ACETATE)	1,0	)4	1,08 c	IIA_	IIA	
	1123 BUTYLACETATES (sec-BUTYL ACETATE)				IIA <sup>7</sup>	IIA	
	1125 n-BUTYLAMINE	0,0		1,13 c	IIA	IIA	
	1127 CHLORORBUTANES (1-CHLORO BUTANE)	1,0		a	IIA	IIA	
	1127 CHLORORBUTANES (2-CHLORO BUTANE)	1,1	16	a	IIA	IIA	
	1127 CHLORORBUTANES (1-CHLORO-2-METHYLPROPANE)	13		a	IIA	IIA	
	1127 CHLORORBUTANES (2-CHLORO-2-METHYLPROPANE)	14		a	IIA	IIA	
	1129 BUTYRALDEHYDE	0,0			IIA	IIA	
	1131 CARBON DISULFIDE	0,3	34	0,39 c	IIC	IIC	
	1134 CHLOORBENZENE (phenylchloride)			d says IIA	IIA	IIA	
	1135 ETHYLENE CHLOROHYDRIN (2-chloroethanol)			d says IIA	IIA <sup>®</sup>	IIA	
	1143 CROTONALDEHYDE	0,0		a	IIB	IIB2	(ADN WP.15/AC.2/2016/4 says IIB2)
	1145 CYCLOHEXANE	0,0		а	IIA	IIA	
	1146 CYCLOPENTANE	1,0		d	IIA	IIA	
	1150 (trans) 1,2-DICHLOROETHYLENE	3,9		a	IIA	IIA	
	1150 (cis)1,2-DICHLOROETHYLENE	3,9	91	a	IIA	IIA	
	1153 ETHYLENE GLYCOL DIETHYLETHER				IIB	IIB2	(ADN WP.15/AC.2/2016/4 says IIB2)
	1154 DIETHYLAMINE	1,		a	IIA	IIA	
	1155 DIETHYL ETHER	0,0	5/	0,88 a	IIB	IIB1	ADN WP.15/AC.2/2016/4 says IIB1
	1157 DIISOBUTYL KETONE			a	IIB4	IIA	all xxxYL KETONES products MESG_IIA,
	1159 DIISOPROPYL ETHER	0,0		а	IIA	IIA	
	1160 DIMETHYLAMINE AQUEOUS SOLUTION	1.	15	a	IIA	IIA	0.1.1.1.10.10.10.10.10.10.10.10.10.10.10
	1163 DIMETHYLHYDRAZINE	0,0		a	IIC	IIB1	Simular to 1,1-Dimethylhydrazine ?
	1165 DIOXANE	0,7		0,19 a	IIB	IIB3	Simular to 1,4-Dioxane and ADN WP.15/AC.2/2016/4 says IIB3
	1167 DIVINYL ETHER	0.0		0.88 c	IIB IIB	IIB3 IIB1	simular to Divinylene oxide) ADN WP.15/AC.2/2016/4 says IIB1
	1170 ETHANOL (ETHYLALCOHOL) > 70% ALCOHOL			U,88 C			
	1170 ETHANOL (ETHYLALCOHOL) 24 < ALCOHOL%, 70	>0	,8		IIB	IIA	according Norm ISO-EIC-DIS 80079-20-1 ≥ 4 vol-% water
	AATA ETINA ENE OLVOOL MONOETINA ETIED				un.	upo.	ADN WP.15/AC.2/2016/4 says IB1 <sup>4</sup> ADN WP.15/AC.2/2016/4 says IB2
	1171 ETHYLENE GLYCOL MONOETHYL ETHER 1172 ETHYLENE GLYCOL MONETHYLETHER-ACETATE	0,0	16 17	a	IIB IIA	IIB2 IIA	ALIN WF. TUMO. 2/2016/4 Says IIBZ
	1172 ETHYLENE GLYCOL MONETHYLETHER-AGETATE	0.0		a a	IIA	IIA	
	1173 ETHYL ACETATE 1175 ETHYL BENZENE	0,0	90	a d	IIA IIA	IIA	
	11/5 ETHYL BENZENE 11/7 2-ETHYLBUTYL ACETATE			d d	IIA	IIA	
	1179 ETHYLBUTYLETHER (ETHYL-tert-BUTYLETHER)			u	IIB4	IIA	Simular to UN 2398 MTBE ?
	1184 ETHYLENE DICHLORIDE (1,2-dichloorethane)	1,8	on.	a	IIA	IIA	Simular to UN 2396 MTBE ?
	1188 ETHYLENE DICHLORIDE (1,2-dichlooremane)	0.8			IIB	IIB1	
		0,0	50	а	IIA <sup>7</sup>		
	1191 OCTYL ALDEHYDES (2-ethylcapronaldehyde)					IIA IIA	
	1191 OCTYL ALDEHYDES (n-OCTALDEHYDE)			a	IIB4		n-octaldehyde
	1193 ETHYLMETHYLKETONE or METHYLETHYLKETONE	0,0		0,92 b	IIA	IIA	classified by MIC
	1198 FORMALDEHYDE SOLUTION	0,8	0/		IIB	IIB	
	1199		00	_	IID.	IID4	ADMIND 45/AC 2004CV - THE HOL
	FURALDEHYDES (a-FURALDEHYDE) or FURFURALDEHYDES (a-FURFURYLALDEHYDE)	0,0	58	а	IIB	IIB1	ADN WP.15/AC.2/2016/4 says IIB1
	1203 MOTORSPIRIT or GASOLINE or PETROL 1208 HEPTANES	0.9	11	0.88 c	IIA IIA	IIA IIA	petrol or gasoline isnt classified in this norm
	1200 HEPTANES 1208 HEXANES	0.0		0,88 c	IIA	IIA	
	1212 ISOBUTANOL or ISOBUTYLALCOHOL	0.0		U,88 C	IIA	IIA	
	1212 ISOBUTANOL OF ISOBUTYLALCOHOL 1213 ISOBUTYLACETATE	1,0		a	IIA <sup>7</sup>	IIA	simular to n-butylacetate
				a	IIA <sup>7</sup>	IIA	
	1214 ISOBUTYLAMINE	1,1		_			simular to DIISOBUTYLAMINE
	1216 ISOOCTENES 1218 ISOPRENE	0,9 0.7	0	a	IIB <sup>4</sup>	IIA IIB2	simular to n-octene
							simular to 1,3-butadiëne and ADN WP.15/AC.2/2016/4 says IIB2
	1219 ISOPROPYLALCOHOL of ISOPROPANOL	1,0	Ju	a	IIA	IIA	

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Methode according to ISO-EIC-DIS 80079-20-1
MESG
                                                                                                                                 FORMULA for MIXTURES
Group IIA: MESG ≥ 0,9 mm.
Group IIB: 0,5 mm < MESG< 0,9 mm.
                                                                                                                                 MESC<sub>1/</sub>
                                                                                                                                                                                                                b = classified according to MIC ratio.
Group IIC: MESG ≤ 0,5 mm.
                                                                                                                                                                                                                c = both MESG and MIC ratio have been determined.
                                                                                                                                         Σ.(X./MESG.)
                                                                                                                                                                                                                d = classified according to similarity of chemical structure (provisional
Group IIA: MIC > 0,8.
Group IIB: 0,45 ≤ MIC ≤ 0,8.
                                                                                                                                                                                                                classification).
Classificatie van gas mengsels
Group IIC: MIC < 0,45.
SUBGROUP IIB
Group IIB: MESG > 0,5 mm.
                                                                                                                                                                                                                                 According to the formula in ISO-EIC-DIS 80079-20-1 for mixtures and the mixture doesn't contain components for what IIB off IIC is described or the mixture contains less 30% of these components. IIB3 complies
Group IIB3: MESG ≥ 0,65 mm.
Group IIB2: MESG ≥ 0,75 mm.
Group IIB1: MESG ≥ 0,85 mm.
Group IIA: MESG ≥ 0,9 mm.
                                                                                                                                         Norm ISO-EIC-DIS 80079-20-1
                                                                                                                                                                                                                 according to
UN Naam
                                                                                                                                         MESG MIC
                                                                                                                                                                                               ADN 2015
   1220 ISOPROPYLE ACETATE
1221 ISOPROPYLAMINE
                                                                                                                                                1,05
                                                                                                                                                                                               IIA<sup>2</sup>
                                                                                                                                                                                                                                  simular to DIISOPROPYLAMINE
                                                                                                                                                1.02
    1223 KEROSINE
                                                                                                                                                                                               IIA7
                                                                                                                                                                                                                                  simular to Diesel Oil No. 1
    1224 KETONES N.O.S
                                                                                                                                                0,84
                                                                                                                                                             0.92 b
                                                                                                                                                                                               IIB<sup>4</sup>
                                                                                                                                                                                                                                  Simular to 2-Butanone (= Ethyl methyl ketone) with the smallest MESG but MIC = 0,92, other ketones all MESG > 0,9
                                                                                                                                                0,93
0,92
0,97
1,10
0,94
    1229 MESITYL OXYDE
                                                                                                                                                                                               IIB<sup>4</sup>
IIA
IIA
IIA
    1230 METHANOL
1231 METHYL ACETATE
                                                                                                                                                              0,82 c
   1235 METHYLAMINE, AQUEOUS solution
1243 METHYL FORMATE
   1244 METHYL HYDRAZINE
1245 METHYL ISOBUTYL KETONE
                                                                                                                                                                                               IIA
IIA
IIA
                                                                                                                                                                                                                                  Simular to dimethyl hydrazine with a MESG = 0,85
                                                                                                                                                1,01
   1247 METHYLMETHACRYLATE
1262 OCTANES
                                                                                                                                                0,95
   1202 OCTANES
1204 PARALDEHYDE
1205 n-PENTANES
1205 PENTANES (2-METHYLBUTANE)
1207 PETROLEUM CRUDE OIL with more than 10% BENZENE
                                                                                                                                                                                               IIA<sup>7</sup>
IIA
IIA
IIB<sup>4</sup>
                                                                                                                                                1,01
0,93
                                                                                                                                                              0,97 c
                                                                                                                                                0.98
                                                                                                                                                                                                                IIΑ
   1267 PETROLEUM CRUDE OIL
1268 PETROLEYUMDESTILLATES NOS. (NAPHTA)
                                                                                                                                                                                               IIB<sup>4</sup>
IIA
IIB<sup>4</sup>
                                                                                                                                                                                                                IIΑ
    1268 PETROLEYUMDESTILLATES NOS
   1288 PETROLEYUMDESTILLATES NOS, with more than 10% BENZENE
1268 PETROLEYUMDESTILLATES NOS. (BENZENE HEARTCUT)
   1274 PROPYLALCOHOL
1275 PROPIONALDEHYDE
1276 n-PROPYLACIFATE
1277 PROPYLAMINE (1-AMINOPROPANE)
1278 1-CHLOROPROPANE (propylchloride)
                                                                                                                                                0,89
0,86
1,04
1,13
                                                                                                                                                                                              IIB
IIB
IIA
IIA
                                                                                                                                                                                                                IIB1
IIB3
                                                                                                                                                                                                                IIA
                                                                                                                                                  _
   1279 1,2-DICHLOROPROPANE (PROPYLENE DICHLORIDE)
1280 PROPYLENE OXIDE
                                                                                                                                                                                               IIA8
                                                                                                                                                0.70
                                                                                                                                                                                               IIB
IIA<sup>8</sup>
                                                                                                                                                                                                                IIR3
   1282 PYRIDINE
1289 SODIUM METHYLATE SOLUTION in alcohol
                                                                                                                                                                                               IIA
IIA
IIA<sup>8</sup>
    1294 TOLUENE
                                                                                                                                                1,06
    1296 TRIETHYLAMINE
                                                                                                                                                                                                                IIA
    1300 TURPENTINE SUBSTITUTE
                                                                                                                                                                                               IIB<sup>4</sup>
IIA
IIA
                                                                                                                                                                                                                                  simular to turpentine oil?
    1301 VINYLACETATE
1307 XYLENES (m-XYLENE) XYLENNE (o-XYLENE)
                                                                                                                                                0,94
1,09
    1545 ALLYL ISOTHIO CYANATE
                                                                                                                                                                                               IIB4
                                                                                                                                                                                                                                  according to METHANENITRIL MESG = 0,8 as nitril with the smallest MESG
    1578 CHLORONITROBENZENES (P-CHLORONITROBENZE)
                                                                                                                                                                                                                                  NITOBENZENE MESG = 0,94 and chloro connections MESG > 0,9
                                                                                                                                                                                               IIB<sup>4</sup>
   1604 ETHYLENEDIAMINE
1648 ACETONITRIL (methylcyanide)
                                                                                                                                                1,18
                                                                                                                                                                                               IIA
                                                                                                                                                1,50
   1662 NITROBENZENE
1663 NITROPHENOLS
1715 ACETIC ANHYDRIDE
                                                                                                                                                                                                                                  (ADN WP.15/AC.2/2016/4 savs IIB1)
                                                                                                                                                                                                                                  according to NITROBENZENE (MESG= 0,94) (phenol= HYDROXYBENZENE AND BENZENE-connections ≥ IIA)
                                                                                                                                                                                               IIB<sup>4</sup>
IIA
IIA<sup>8</sup>
                                                                                                                                                1,23
    1717 ACETYL CHLORIDE
                                                                                                                                                                                                                ПΔ
    1738 BENZYL CHLORIDE
                                                                                                                                                                                               IIA<sup>8</sup>
IIA
IIA
IIB<sup>4</sup>
IIB<sup>4</sup>
   1738 BENZYL CHLORIDE
1750 CHLOROACETIC ACID SOLUTION
1764 DICHLOROACETIC ACID
1779 FORMIC ACID WITH MORE THAN 85% acid by mass
                                                                                                                                                                                                                                 all acetic connections MESG > 0,9
all acetic connections MESG > 0,9
                                                                                                                                                1,86
                                                                                                                                                                                                                ПΔ
    1783 HEXAMETHYLENEDIAMINE solution
                                                                                                                                                                                                                                  according to simular diamineconnections and cyclohexane (hexamethylene) MESG = > 0,9
    1863 FUEL, AVIATION, TURBINE ENGINE
                                                                                                                                                                                                                                 simular to. Kerosine and benzene connections MESG > 0,9
   1863 FUEL, AVIATION, TURBINE ENGINE (with more than 10% BENZEEN 1915 CYCLOHEXANONE
                                                                                                                                                                                               IIB<sup>4</sup>
                                                                                                                                                0,95
                                                                                                                                                                                                                                  ADN WP.15/AC.2/2016/4 says IIB1
    1917 FTHYLACRYLATE
                                                                                                                                                0,86
                                                                                                                                                                                              IIB
IIA<sup>8</sup>
                                                                                                                                                                                                                IIB1
    1918 ISOPROPYLBENZEEN (cumeen)
                                                                                                                                                1.05
                                                                                                                                                                                                                IIA
                                                                                                                                                                                                                                 according to MIC clasification IIA
(ADN WP.15/AC.2/2016/4 says IIB1)
    1919 METHYLACRYLATE
                                                                                                                                                0,85
                                                                                                                                                              0,98 b
                                                                                                                                                                                               IIB
    1920 NONANES
                                                                                                                                                                                                                IIA
    1922 PYRROLIDINE
                                                                                                                                                                                               IIA7
                                                                                                                                                                                                                                  cyclic amine connection MESG > 0,9
                                                                                                                                                                                                                IIA
    1988 ALCOHOLS, FLAMABLE, TOXIC, NOS
    1987 ALCOHOLS NOS
                                                                                                                                                                                               IIB<sup>4</sup>
IIA
IIB<sup>4</sup>
IIB<sup>4</sup>
                                                                                                                                                                                                                                  simular to xxx-anol group not toxic
   1987 ALCOHOLS, NOS (cyclohexanol)
1987 ALCOHOLS, NOS (90 mas-% tert.-BUTANOL en 10 mas-% METHANOL)
    1989 ALDEHYDES NOS
                                                                                                                                                                                                                                  IF non-formaldehyde or connections of formaldehyde present in the mixture, else UN 1198
                                                                                                                                                                                                                IIB2
   1991 CHLOROPRENE
1992 FLAMABLE LIQUID TOXIC, NOS
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	MESG	FORMULA for MIXTURE	ES		Methode acc	cording to ISO-EIC-DIS 80079-20-1
	Group IIA: MESG ≥ 0,9 mm.	MESC <sub>1/</sub>				I according to MESG determination.  I according to MIC ratio.
	Group IIB: 0,5 mm < MESG< 0,9 mm. Group IIC: MESG ≤ 0,5 mm.					•
	MIC MESOS 0,5 mm.	$\Sigma_i(X_i/MESG_i)$				SG and MIC ratio have been determined.
	Group IIA: MIC > 0,8.				d = classified classification	l according to similarity of chemical structure (provisional
	Group IIB: $0.45 \le MIC \le 0.8$ .				Classificatie v	van gas mengsels
	Group IIC: MIC < 0,45.					
	SUBGROUP IIB Group IIB: MESG > 0.5 mm.				***	According to the formula in ISO-EIC-DIS 80079-20-1 for mixures and the mixture doesn't contain components for what IIB off IIC is described or the mixture contains less 30% of these components, IIB3 complied.
	Group IIB3: MESG ≥ 0,65 mm.					components for what no on no to described of the flux tale contains less 50% of these components, indo complie
	Group IIB2: MESG ≥ 0,75 mm.					
	Group IIB1: MESG ≥ 0,85 mm. Group IIA: MESG ≥ 0,9 mm.	Norm ISO-EIC-D	ue onn7g 2n			
					according to	
	UN Naam	MESG MIC	method	ADN 2015	this norm	
	1993 FLAMABLE LIQUID , NOS.(CYCLOHEXANONE / CYCLOHEXANOL MIXTURE)			IIA	IIA	
	1993 FLAMABLE LIQUID , NOS.(with more than 10% BENZENE) 1993 FLAMABLE LIQUID , NOS			IIB <sup>4</sup>	IIB3 IIB3	### ###
	1999 TARS, LIQUID, including road oils and cutback butimens			IIA <sup>7</sup>	IIA	THE STATE OF THE S
	2021 CHLOROPHENOLS, LIQUID (2-CHLOROPHENOL)			IIA <sup>7</sup>	IIA	
	2022 CRESILIC ACID			IIA7	IIA	
	2023 EPICHLOROHYDRINE	0,74	a	IIB	IIB3	ADN WP.15/AC.2/2016/4 says IIB3
	2046 ISOBUTYRALDEHYDE (ISOBUTYL ALDEHYDE) 2046 CYMENES		100	IIA <sup>7</sup>	IIA IIA	simular to butyraldehyde MESG = 0,92
	2040 CYMENES 2047 1.3-DICHLOROPROPENE	_	d	IIA'	IIA	
	2047 1,3-DICHLOROPROP-1-ENE			IIA <sup>7</sup>	IIA	0.008410035
	2047 MIXTURE of 2,3-DICHLOROPROP-1-ENE and 1,3-DICHLOROPROPENE)			IIA <sup>7</sup>	IIA	
	2048 DICYCLOPENTADIENE	0,91	a	IIA	IIA	
	2050 DIISOBUTYLENE			IIA <sup>7</sup>	IIA	simular to isobutylene MESG = 1,00
	2051 2-DIMETHYLAMINO ETHANOL 2053 METHYLISOBUTYLCARBINOL (methylamylalcohol)	1.01	_	IIA <sup>7</sup> IIB <sup>4</sup>	IIA IIA	
	2054 MORPHOLINE	0.92	a a	IIA	IIA	
	2055 STYRENE, monomer		1,21 b	IIA	IIA	
	2056 TETRAHYDROFURAN	0,87	а	IIB	IIB1	ADN WP.15/AC.2/2016/4 says IIB1
	2057 TRIPROPYLENE (nonene) 2078 TOLUENE DIISOCYANATE			IIB <sup>4</sup>	IIA IIA	according to octene ( IIA), and decene isnt even classified.  TOLUENE connection MESG > 0,9 and simular to methylisocyanate MESG = 1,21
	2205 ADIPONITRILE			IIB <sup>4</sup>	IIA?	simular to cyanides without dubble connections and butane?
	2215 MALEIC ANHYDRIDE, MOLTEN			IIB4	IIA?	simular to ACETIC ANHYDIDE gives MESG > 1.00
	2218 ACRYLIC ACID, STABILIZED	0,86	a	IIIB	IIB1	ADN WP.15/AC.2/2016/4 says IIB1
	2227 n-BUTYL METHACRYLATE	0,95	a	IIA	IIA	
	2238 p, m or o CHLOROTOLUENE 2241 CYCLOHEPTANE	_	d d	IIA <sup>7</sup>	IIA IIA	
	2247 OTOLOHEFTANE 2247 n-DECANE	1,05	a	IIA	IIA	
	2248 DI-n-BUTYLAMINE	.,	-	IIA <sup>7</sup>	IIA	simular to n-butylamine with the smallest MESG (0,92) from all butylamines
	2259 TRIETHYLENETETRAMINE			IIB <sup>4</sup>	IIA	simular to the smallest MESG (0,92) from all amineconnections
	2283 DIMETHYLCYCLOHEXANEN		d	IIA <sup>7</sup>	IIA	
	2264 N,N-DIMETHYLCYCLOHEXYLAMINE 2265 N,N-DIMETHYLFORMAMIDE	1,08	d	IIB <sup>4</sup>	IIA	simular to all N,N xxx Amines (MESG > 0,9)
	2268 DIMETHYL-N-PROPYLAMINE	1,00		IIA <sup>7</sup>	IIA	simular to the smallest MESG (0.92) from all amineconnections
	2276 2-ETHYLHEXYLAMINE			IIA <sup>7</sup>	IIA	simular to the smallest MESG (0.92) from all amineconnections
						simular to all alkenes ≥ C3 (MÈSG > 0,9)
	2278 n-HEPTENE 2280 HEXAMETHYLENEDIAMINE. SOLID.			IIB <sup>4</sup>	IIA IIA	(ADN_WP.15/AC.2/2016/4 says IIB1) simular to the smallest MESG (0,92) from all amineconnections
	2280 HEXANOLS 2282 HEXANOLS			IIA	IIB1	simular to the smallest MESG (0,92) from all amineconnections simular to 1-hexanol (MESG = 0.85)
	2286 PENTAMETHYLHEPTANE		d	IIA <sup>7</sup>	IIA	
	2288 ISOHEXENEN			IIB4	IIA	simular to all alkenes ≥ C3 (MESG > 0,9)
	2289 ISOPHORONEDIAMINE			IIA <sup>7</sup>	IIA	simular to the smallest MESG (0,92) from all amineconnections
	2302 5-METHYLHEXAN-2-ONE 2303 ISOPROPENYLBENZENE	0.88	a	IIA IIB	IIA IIB1	simular to hexan-2-on (MESG =0,98) ADN WP.15/AC.2/2016/4 savs IIB1
	2309 OCTADIENE (1,7-OCTADIENE)	2,00	-	IIB <sup>4</sup>	IIA	simular to penta-1,3-diene (MESG = 0,97)?
	2312 PHENOL, MOLTEN		d	IIA <sup>8</sup>	IIA	and the state of t
	2321 TRICHLOROBENZENES			IIA <sup>7</sup>	IIA	all chloro and benzene connections MESG > 0,9
	2323 TRIETHYL PHOSPHITE			IIB <sup>4</sup>	?	
	2324 TRIISOBUTYLENE 2325 1,3,5-TRIMETHYLBENZENE	0.98	_	IIB <sup>4</sup> IIA <sup>7</sup>	IIA IIA	Simular to isobutylene (MESG = 1,00)
	2323 ALLYLACETATE	0,98	a a	IIA <sup>7</sup>	IIA	
	2348 BUTYLACRYLATEN, n-BUTYLACRYLAAT	0,88	a	iiB	IIB1	ADN WP.15/AC.2/2016/4 says IIB1
	2350 BUTYLMETHYLETHER			IIB <sup>4</sup>	IIA	simular to tert-Butyl methylether (MESG = 1,0)
	2356 2-CHLOROPROPANE	1,32	a	IIA	IIA	
	2357 CYCLOHEXYLAMINE 2362 1,1-DICHLOROETHANE	1,82	a	IIB <sup>4</sup> IIA	IIA IIA	simular to the smallest MESG (0,92) from all amineconnections
	2370 1- HEXENE	1,02	a	IIB4	IIA	simular to all alkenes ≥ C3 (MESG > 0.9)
	2381 DIMÉTHYL DISULPHIDE			IIB	IIA	simular to the group sulfides MESG > 0,9
	2382 DIMETHYL HYDRAZINE SYMMETRICAL	0.05	_	IIC <sup>6</sup>	IIB1?	Simular to dimethyl hydrazine with a MESG = 0,85
	2383 DIPROPYLAMINE 2397 3-METHYLBUTAN-2-ONE	0,95	a	IIA IIA <sup>7</sup>	IIA IIA	
	2398 METHYL-tert-BUTYLETHER (MTBE)	1.00	a	IIA	IIA	
	2404 PROPIONITRILE		_	IIA <sup>7</sup>	IIA	
	2414 THIOPHENE	0,91	a	IIA	IIA	

ESG FORMULA for MIXTURES			Methode according to ISO-EIC-DIS 80079-20-1				
Group IIA: MESG ≥ 0,9 mm.  Group IIB: 0.5 mm < MESG< 0.9 mm.	MESC <sub>1/</sub>				a = classified according to MESG determination. b = classified according to MIC ratio.		
Group IIC: MESG   0.5 mm.	Σ <sub>i</sub> (X <sub>i</sub> /MESG <sub>i</sub> )				SG and MIC ratio have been determined.		
MIC	2 (14.11.233)				d according to similarity of chemical structure (provisional		
Group IIA: MIC > 0,8. Group IIB: 0.45 ≤ MIC ≤ 0,8.				classification	n). van gas mengsels		
Group IIC: MIC < 0.45.				Classificatie	van gas mengsels		
SUBGROUP IIB				###	According to the formula in ISO-EIC-DIS 80079-20-1 for mixures and the mixture doesn't contain		
Group IIB: MESG > 0,5 mm.  Group IIB3: MESG ≥ 0.65 mm.					components for what IIB of IIC is described or the mixture contains less 30% of these components, IIB3 complies		
Group IIB2: MESG ≥ 0,75 mm.							
Group IIB1: MESG ≥ 0,85 mm. Group IIA: MESG ≥ 0,9 mm.	Norm ISO-EIC-DI	90079-20-4					
UN Naam	MESG MIC	method.		according t	0		
	MESO MIC	metriou.	ADN 2015	this norm			
2430 ALKYLPHENOLS SOLID, NOS MOLTEN 2458 HEXADIENES			IIA <sup>7</sup> IIB <sup>4</sup>	IIA IIA	(ADN WP.15/AC.2/2016/4 says IIA)		
2477 METHYL ISOTHIOCYANATE			IIB4	IIA	simular to METHYLISOCYANATE and thio-groups (all MESG >0,9)		
2485 n-BUTYLISOCYANATE or ISOBUTYLISOCYANATE			IIB <sup>4</sup>	IIA	simular to METHYLISOCYANATE(MESG = 1,21)		
2486 ISOBUTYL ISOCYANATE			IIB <sup>4</sup>	IIA	simular to METHYL ISOCYANATE en PHENYL ISOCYANATE (UN 2487)		
2487 PHENYLISOCYANATE 2491 ETHANOLAMINE			IIA IIB <sup>4</sup>	IIA IIA	simular to METHYLISOCYANATE(MESG = 1,21) (ADN WP.15/AC.2/2016/4 says IIA)		
2493 HEXAMETHYLENEIMINE			IIA	IIA	(ALIV VVI. TURO. 220 For Salys III) simular to cyclobran en amines		
2527 ISOBUTYLACRYLATE			IIBº	IIB1	simular to n-Butyl acrylate (MESG = 0,88)		
2528 ISOBUTYLISOBUTYRATE	1,00	a	IIA	IIA			
2531 METHACRYLIC ACID, 2564 TRICHLOROACETIC ACID	0,95	a	IIB <sup>4</sup> IIA <sup>7</sup>	IIA IIA			
2608 NITROPROPANES	0.84	a	IIB <sup>7</sup>	IIB2	ADN WP.15/AC.2/2016/4 says IIB2		
2615 ETHYLPROPYLETHER	-,	-	IIA <sup>7</sup>	IIA			
2618 VINYLTOLUENE			IIB <sup>4</sup>	IIA	simular to vinylbenzene (MIC= 1,21)		
2683 AMMONIUM SULPHIDE SOLUTION 2709 BUTYLBENZENES (n-BUTYLBENZENE)			IIB <sup>4</sup> IIA <sup>7</sup>	IIA IIA	according to NH3 and Sulfides IIA		
2733 2-AMINOBUTANE	0.92	С	IIA <sup>7</sup>	IIA	simular to 1-aminobutane (MESG = 0,92)		
2789 ACETIC ACID SOLUTION, more than 80% acid by mass.	0,02	•	IIA <sup>7</sup>	IIA	on the control of the		
2811 TOXIC SOLID ORGANIC NOS (1,3,5, trichlorobezene, MOLTEN)			IIB <sup>4</sup>	IIA	all chloro and benzene connections MESG > 0,9, and ADN WP.15/AC.2/2016/4 says IIA		
2920 HEXADECYLTRIMETHYLAMMONIUM CHLORIDE (50%) and ETHANOL 35%) 2920 2-PROPANOL and DODECYLDIMETHYLAMMONIUM CHLORIDE			IIB IIA	IIB1 IIA	according to Ethanol IIB1, all chlorides MESG. 0,9 According to propanol IIA, all chlorides MESG > 0,9		
2924 AQUEOUS SOLUTION OF DIALKYL-(C8-C18)-DIMETHYLAMMONIUM CHLORIDE			IIA	IIA	According to property in a second control of the co		
and 2-PROPANOL)							
2924 FLAMMABLE LIQUID, CORROSIVE, NOS 2929 TOXIC ORGANIC LIQUID FLAMMABLE. NOS			IIB <sup>4</sup>	IIB3 IIB3	simular to all ADN class 8 products there is no IIB product, so IIB3 complies		
2935 ETHYL-2-CHLORO-PROPIONATE			IIA	IIA	***		
2947 ISOPROPYL CHLOROACETATE	1,24	a	IIA	IIA			
2983 ETHYLENE OXIDE AND PROPYLENE OXIDE MIXTURE, WITH ≤ 30% ETHYLENE OXIDE			IIB IIB <sup>4</sup>	IIB3 IIB2	According to the formula in ISO-EIG-DIS 80079-20-1 for mixures, ETH oX MESG 0,59 and 30%, prop ox MESG 0,7 and 70%		
3079 METHACRYLNITRILE 3092 1-METHOXY-2-PROPANOL			IIB	IIB1	according to methanenitrile (HCN) MESG-0,8 (ADN WP.15/AC.2/2016/4 says IIB1)		
3175 SOLIDS CONTAINING FLAMMABLE LIQUID, N.O.S., MOLTEN			IIA <sup>7</sup>	IIA	,		
3256 ELEVATED TEMPERATURE LIQUID, FLAMMABLE, N.O.S.			IIB <sup>4</sup>	IIB3	UN 3256 mainly contains hydrocarbons with large alkane chains and connections with cyclo products MESG > 0.9, so IIB3 complies		
3271 ETHERS, NOS. (tertAMYLMETHYLETHER)			IIB <sup>4</sup>	IIB3 IIB3	simular to dietylene ether (MESG=0,7)		
3271 ETHERS, NOS. (tertAMYLMETHYLETHER) 3272 ESTERS. NOS			IIB <sup>4</sup>	IIB1	simular to dietylene ether (MESG=0,7) simular to the products with the smallest MESG within the group acetic acid X esters, ((= Acetoacetic acid methyl ester) MESG= 0.85		
3286 FLAMMABLE LIQUID. TOXIC CORROSIVE, N.O.S.			IIB <sup>4</sup>	IIB3	according to the toxic propertys and coorosive propertys IIB3 complies		
3295 HYDROCARBONS LIQUID NOS, with > 10% BENZENE			IIB <sup>4</sup>	IIB3	###		
3295 HYDROCARBONS LIQUID NOS,(1-OCTEEN)			IIB <sup>4</sup>	IIB3	***		
3295 HYDROCARBONS LIQUID NOS, (mixture of POLYCYCL AROMATES) 3295 HYDROCARBONS LIQUID NOS.CONTAINS ISOPRENE EN PENTADIEN			IIA IIB <sup>4</sup>	IIA IIB3	***		
3412 FORMIC ACID	1,86	a	IIA	IIA			
3429 CHLOROTOLUIDINES, LIQUID			IIA <sup>7</sup>	IIA			
3446 NITROTOLUENES, SOLID, MOLTEN			IIB <sup>4</sup> IIA <sup>8</sup>	IIB3 IIA	simular to NITROBENZENE MESG > 0,9 en TOLUENE MESG > 1, BII3 complies		
3451 TOLUIDINES, SOLID, MOLTEN 3455 CRESOLS, SOLID, MOLTEN			IIA <sup>8</sup>	IIA			
3463 PROPIONIC ACID with not less than 90% acid by mass			IIA <sup>7</sup>	IIA			
3475 ETHANOL AND GASOLINE MIXTURE with > 10% but < 90% ethanol			IIA	IIA	4		
3475 ETHANOL AND GASOLINE MIXTURE met > 90% ethanol 3494 PETROLEUM SOUR CRUDE OIL. FLAMMABLE, TOXIC			IIB IIB <sup>4</sup>	IIB1 IIB3	According to the formula in ISO-EIC-DIS 80079-20-1 for mixtures 90% ethanol		
9001 SUBSTANCESWITH A FLASHPOINT FP > 60 °C, HANDED OVER FOOR CARRIAGE AT A			IIB <sup>4</sup>	IIB3	### ###		
TEMP WITHIN 15 k FROM THE FLASHPOINT							