

STATEMENT ON A NONPROPRIETARY NAME ADOPTED BY THE USAN COUNCIL

USAN (JK-188)	LORIGERLIMAB
PRONUNCIATION	lor" i jer' li mab
THERAPEUTIC CLAIM	Treatment of cancer

CHEMICAL NAMES

1. Immunoglobulin, anti-(human programmed cell death protein 1) (human-Mus musculus clone MG13.78 κ -chain V-J region) fusion protein with peptide linker (GGGSGGGG) fusion protein with immunoglobulin anti-(human cytotoxic T-lymphocyte-associated protein 4) (human clone 4B6 heavy chain V-D-J region) fusion protein with peptide linker (GGCGGG) fusion protein with peptide (synthetic E-coil) fusion protein with immunoglobulin G4 [10-proline,34-tyrosine,36-threonine,38-glutamic acid,de-229-lysine] (human γ 4-chain C region C-terminal fragment), disulfide with immunoglobulin anti-(human cytotoxic T-lymphocyte-associated protein 4) (human clone 4B6 κ -chain V-J region) fusion protein with linker peptide (GGGSGGGG) fusion protein with immunoglobulin anti-(human programmed cell death protein 1) (human-Mus musculus clone MG13.78 heavy chain V-D-J region) fusion protein with peptide linker (GGCGGG) fusion protein with peptide (synthetic K-coil), dimer (Source: CAS)
2. Immunoglobulin G4 scFv-h-CH2-CH3_scFv, anti-[*Homo sapiens* PDCD1 (programmed cell death 1, PD1, PD-1, CD279)] and anti-[*Homo sapiens* CTLA4 (cytotoxic T-lymphocyte-associated protein 4, CD152)], chimeric, *Homo sapiens* and humanized monoclonal antibody, tetravalent, bispecific; scFv-h-CH2-CH3, chimeric and *Homo sapiens*, anti-PDCD1 and anti-CTLA4 (1-499) [scFv-V-KAPPA-VH (1-237) [V-KAPPA anti-PDCD1 Musmus/Homsap (*Mus musculus* IGKV3-2*01 (80.8%) -(IGHD) -IGKJ2*01 (91.7%) L124>V (108)/*Homo sapiens* IGKV3D-11*02 (77.7%) -(IGHD) -IGKJ4*01 (100%), CDR-IMGT [10.3.9] (27-36.54-56.93-101)) (1-111) -8 mer triglycyl-seryl-tetraglycyl linker (112-119) -VH anti-CTLA4 (*Homo sapiens* IGHV3-30*01 (93.9%) -(IGHD) -IGHJ4*01 (100%), CDR-IMGT [8.8.11] (145-152.170-177.216-226)) (120-237)] -E-coil linker (238-271) [6-mer diglycyl-cysteinyl-triglycyl linker (238-243) -28-mer (glutamyl-valyl-dialanyl-cysteinyl-glutamyl-lysyl)-tris(glutamyl-valyl-dialanyl-leucyl-glutamyl-lysyl) E-coil motif (244-271)] - [scFc *Homo sapiens* IGHG4*01 h-CH2-CH3, G4v5 h P10, G4v21 CH2 Y15.1, T16, E18 (hinge 1-12 S10>P (281) (272-283), CH2 M15.1>Y (305), S16>T (307), T18>E (309) (284-393), CH3 (394-498), CHS K2>del (499)) (272-499)] (248'-246')-disulfide with scFv-V-KAPPA-VH, *Homo sapiens* and humanized, anti-CTLA4 and anti-PDCD1 (1'-269') [V-KAPPA anti-CTLA4 (*Homo sapiens* IGKV3-20*01 (99.0%) -IGKJ1*01 (100%), CDR-IMGT [7.3.9] (27-33.51-53.90-98)) (1'-108') -8 mer triglycyl-seryl-tetraglycyl linker (109'-116') -VH anti-PDCD1 (*Homo sapiens* IGHV1-46*01 (81.6%) -(IGHD) -IGHJ4*01 (92.9%), CDR-IMGT [8.8.12] (142-149.167-174.213-224)) (117'-235')] -K-coil linker (236'-269') [6-mer diglycyl-cysteinyl-triglycyl linker (236'-241') -28-mer (lysyl-valyl-dialanyl-cysteinyl-lysyl-glutamyl)-tris(lysyl-valyl-dialanyl-leucyl-lysyl-

glutamyl) K-coil motif (242'-269')]; dimer (279-279":282-282")-bisdisulfide, produced in a Chinese hamster ovary (CHO)-S cell line, glycoform alfa (Source: WHO pINN list 125)

STRUCTURAL FORMULA

Heavy chain X, X''

EIVLTQSPAT	LSLSPGERAT	LSCRASESVD	NYGMSFMNWF	QQKPGQPPKL	50
LIHAASNQGS	GVPSRFSGSG	SGTDFTLTIS	SLEPEDFAVY	FCQQSKEVPY	100
TFGGGTKVEI	KGGGSGGGGQ	VQLVESGGGV	VQPGRSLRLS	CAASGFTFSS	150
YTMHWVRQAP	GKLEWVTFI	SYDGSNKHYA	DSVKGRFTVS	RDNSKNTLYL	200
QMNSLRAEDT	AIYYCARTGW	LGPFDYWGQG	TLVTVSSGGC	GGGEVAACEK	250
EVAALEKEVA	ALEKEVAALE	KESKYGPPCP	PCPAPEFLGG	PSVFLFPPKP	300
KDTLYITREP	EVTCVVVDVS	QEDPEVQFNW	YVDGVEVHNA	KTKPREEQFN	350
STYRVVSVLT	VLHQDWLNGK	EYKCKVSNKG	LPSSIEKTIS	KAKGQPREPQ	400
VYTLPPSQEE	MTKNQVSLTC	LVKGFYPSDI	AVEWESNGQP	ENNYKTTTPV	450
LDSDGSPFLY	SRLTVDKSRW	QEGNVFSCSV	MHEALHNHYT	QKSLSLSLG	499

Light chain X', X'''

EIVLTQSPGT	LSLSPGERAT	LSCRASQSVS	SSFLAWYQQK	PGQAPRLLIY	50'
GASSRATGIP	DRFSGSGSGT	DFTLTISRLE	PEDFAVYQC	QYGSSPWTFG	100'
QGTKVEIKGG	GSGGGGQVQL	VQSGAEVKKP	GASVKVSCKA	SGYSFTSYWM	150'
NWVRQAPQG	LEWIGVIHPS	DSETWLDQKF	KDRVTTITVDK	STSTAYMELS	200'
SLRSEDTAVY	YCAREHYGTS	PFAYWGQGTL	VTVSSGGCGG	GKVAACKEKV	250'
AALKEKVAAL	KEKVAALKE				269'

Disulfide bridges

23'-89'	23'''-89'''	23-92	23''-92''	138'-212'	138'''-212'''	141-215	141''-215''
240-238'	240''-238''	248-246'	248''-246''	279-279''	282-282''	314-374	314''-374''
420-478	420''-478''						

Glycosylation sites (N)

350	350''
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MOLECULAR FORMULA C₇₃₉₀H₁₁₃₈₆N₁₉₇₆O₂₂₉₉S₅₂ (EXPASY)

MOLECULAR WEIGHT 166.4 kDa (EXPASY)

TRADEMARK None as yet

SPONSOR MacroGenics, Inc.

CODE DESIGNATIONS MGD019, AEX1344

CAS REGISTRY NUMBER 2416595-46-3

UNII JM802R57SA

WHO NUMBER 11771

gbk