



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:00 AM UTC

PDB ID : 8W83 / pdb_00008w83
Title : HLA-DQ2.5-alpha1 gliadin peptide in complex with DQN0344AE02
Authors : Irie, M.; Tsushima, T.; Teranishi-Ikawa, Y.; Takahashi, N.; Ishii, S.; Okura, Y.; Fukami, T.A.; Torizawa, T.
Deposited on : 2023-08-31
Resolution : 2.82 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

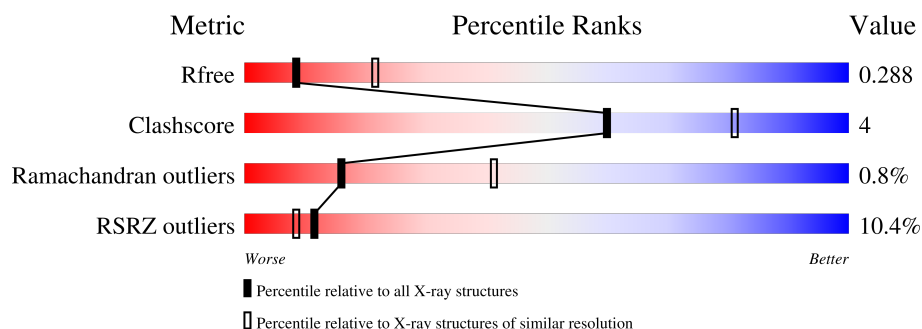
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	228	4% 88% 8% .
1	E	228	7% 82% 13% 5%
1	I	228	11% 84% 11% 5%
1	M	228	24% 80% 10% 10%
2	B	215	4% 89% 11%
2	F	215	20% 83% 15% .
2	J	215	7% 90% 7% .

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Mol	Chain	Length	Quality of chain
2	N	215	
3	C	189	
3	G	189	
3	K	189	
3	O	189	
4	D	226	
4	H	226	
4	L	226	
4	P	226	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 23643 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called DQN0344AE02 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1620	1025	269	317	9			
1	E	217	Total	C	N	O	S	0	0	0
			1605	1017	262	317	9			
1	I	217	Total	C	N	O	S	0	0	0
			1589	1007	261	312	9			
1	M	206	Total	C	N	O	S	0	0	0
			1506	945	254	298	9			

- Molecule 2 is a protein called DQN0344AE02 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	214	Total	C	N	O	S	0	0	0
			1620	1016	266	333	5			
2	F	212	Total	C	N	O	S	0	0	0
			1522	954	253	310	5			
2	J	209	Total	C	N	O	S	0	0	0
			1548	971	252	321	4			
2	N	212	Total	C	N	O	S	0	0	0
			1529	961	245	318	5			

- Molecule 3 is a protein called HLA class II histocompatibility antigen, DQ alpha 1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	179	Total	C	N	O	S	0	0	0
			1407	907	227	271	2			
3	G	179	Total	C	N	O	S	0	0	0
			1357	871	215	269	2			
3	K	179	Total	C	N	O	S	0	0	0
			1354	874	219	259	2			
3	O	178	Total	C	N	O	S	0	0	0
			1370	881	227	260	2			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	47	SER	CYS	engineered mutation	UNP P01909
C	184	LEU	-	expression tag	UNP P01909
C	185	GLU	-	expression tag	UNP P01909
C	186	VAL	-	expression tag	UNP P01909
C	187	LEU	-	expression tag	UNP P01909
C	188	PHE	-	expression tag	UNP P01909
C	189	GLN	-	expression tag	UNP P01909
G	47	SER	CYS	engineered mutation	UNP P01909
G	184	LEU	-	expression tag	UNP P01909
G	185	GLU	-	expression tag	UNP P01909
G	186	VAL	-	expression tag	UNP P01909
G	187	LEU	-	expression tag	UNP P01909
G	188	PHE	-	expression tag	UNP P01909
G	189	GLN	-	expression tag	UNP P01909
K	47	SER	CYS	engineered mutation	UNP P01909
K	184	LEU	-	expression tag	UNP P01909
K	185	GLU	-	expression tag	UNP P01909
K	186	VAL	-	expression tag	UNP P01909
K	187	LEU	-	expression tag	UNP P01909
K	188	PHE	-	expression tag	UNP P01909
K	189	GLN	-	expression tag	UNP P01909
O	47	SER	CYS	engineered mutation	UNP P01909
O	184	LEU	-	expression tag	UNP P01909
O	185	GLU	-	expression tag	UNP P01909
O	186	VAL	-	expression tag	UNP P01909
O	187	LEU	-	expression tag	UNP P01909
O	188	PHE	-	expression tag	UNP P01909
O	189	GLN	-	expression tag	UNP P01909

- Molecule 4 is a protein called MHC class II HLA-DQ-beta-1 - alpha1 gliadin peptide chimeric protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	181	Total 1428	C 917	N 239	O 266	S 6	0	0	0
4	H	178	Total 1347	C 861	N 228	O 252	S 6	0	0	0
4	L	179	Total 1344	C 851	N 239	O 247	S 7	0	0	0
4	P	186	Total 1441	C 916	N 252	O 266	S 7	0	0	0

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	984	GLY	-	linker	PDB ?
D	985	SER	-	linker	PDB ?
D	986	GLY	-	linker	PDB ?
D	987	GLY	-	linker	PDB ?
D	988	GLY	-	linker	PDB ?
D	989	GLY	-	linker	PDB ?
D	990	SER	-	linker	PDB ?
D	991	ILE	-	linker	PDB ?
D	992	GLU	-	linker	PDB ?
D	993	GLY	-	linker	PDB ?
D	994	ARG	-	linker	PDB ?
D	995	GLY	-	linker	PDB ?
D	996	SER	-	linker	PDB ?
D	997	GLY	-	linker	PDB ?
D	998	GLY	-	linker	PDB ?
D	999	GLY	-	linker	PDB ?
D	1000	SER	-	linker	PDB ?
D	1191	LEU	-	expression tag	UNP O19712
D	1192	GLU	-	expression tag	UNP O19712
D	1193	VAL	-	expression tag	UNP O19712
D	1194	LEU	-	expression tag	UNP O19712
D	1195	PHE	-	expression tag	UNP O19712
D	1196	GLN	-	expression tag	UNP O19712
H	984	GLY	-	linker	PDB ?
H	985	SER	-	linker	PDB ?
H	986	GLY	-	linker	PDB ?
H	987	GLY	-	linker	PDB ?
H	988	GLY	-	linker	PDB ?
H	989	GLY	-	linker	PDB ?
H	990	SER	-	linker	PDB ?
H	991	ILE	-	linker	PDB ?
H	992	GLU	-	linker	PDB ?
H	993	GLY	-	linker	PDB ?
H	994	ARG	-	linker	PDB ?
H	995	GLY	-	linker	PDB ?
H	996	SER	-	linker	PDB ?
H	997	GLY	-	linker	PDB ?
H	998	GLY	-	linker	PDB ?
H	999	GLY	-	linker	PDB ?
H	1000	SER	-	linker	PDB ?
H	1191	LEU	-	expression tag	UNP O19712
H	1192	GLU	-	expression tag	UNP O19712

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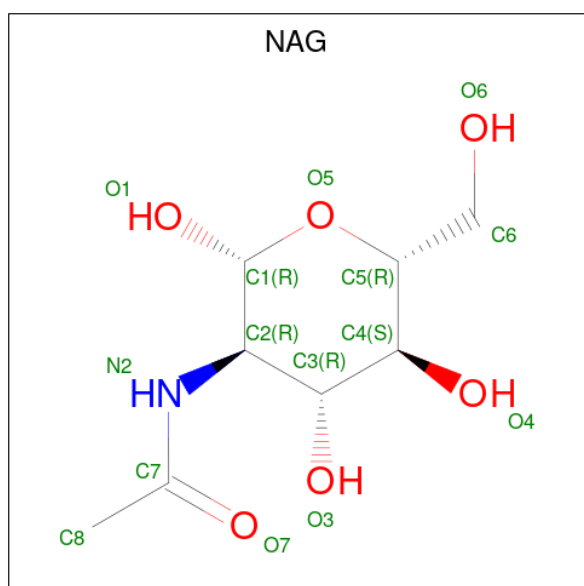
Chain	Residue	Modelled	Actual	Comment	Reference
H	1193	VAL	-	expression tag	UNP O19712
H	1194	LEU	-	expression tag	UNP O19712
H	1195	PHE	-	expression tag	UNP O19712
H	1196	GLN	-	expression tag	UNP O19712
L	984	GLY	-	linker	PDB ?
L	985	SER	-	linker	PDB ?
L	986	GLY	-	linker	PDB ?
L	987	GLY	-	linker	PDB ?
L	988	GLY	-	linker	PDB ?
L	989	GLY	-	linker	PDB ?
L	990	SER	-	linker	PDB ?
L	991	ILE	-	linker	PDB ?
L	992	GLU	-	linker	PDB ?
L	993	GLY	-	linker	PDB ?
L	994	ARG	-	linker	PDB ?
L	995	GLY	-	linker	PDB ?
L	996	SER	-	linker	PDB ?
L	997	GLY	-	linker	PDB ?
L	998	GLY	-	linker	PDB ?
L	999	GLY	-	linker	PDB ?
L	1000	SER	-	linker	PDB ?
L	1191	LEU	-	expression tag	UNP O19712
L	1192	GLU	-	expression tag	UNP O19712
L	1193	VAL	-	expression tag	UNP O19712
L	1194	LEU	-	expression tag	UNP O19712
L	1195	PHE	-	expression tag	UNP O19712
L	1196	GLN	-	expression tag	UNP O19712
P	984	GLY	-	linker	PDB ?
P	985	SER	-	linker	PDB ?
P	986	GLY	-	linker	PDB ?
P	987	GLY	-	linker	PDB ?
P	988	GLY	-	linker	PDB ?
P	989	GLY	-	linker	PDB ?
P	990	SER	-	linker	PDB ?
P	991	ILE	-	linker	PDB ?
P	992	GLU	-	linker	PDB ?
P	993	GLY	-	linker	PDB ?
P	994	ARG	-	linker	PDB ?
P	995	GLY	-	linker	PDB ?
P	996	SER	-	linker	PDB ?
P	997	GLY	-	linker	PDB ?
P	998	GLY	-	linker	PDB ?

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Chain	Residue	Modelled	Actual	Comment	Reference
P	999	GLY	-	linker	PDB ?
P	1000	SER	-	linker	PDB ?
P	1191	LEU	-	expression tag	UNP O19712
P	1192	GLU	-	expression tag	UNP O19712
P	1193	VAL	-	expression tag	UNP O19712
P	1194	LEU	-	expression tag	UNP O19712
P	1195	PHE	-	expression tag	UNP O19712
P	1196	GLN	-	expression tag	UNP O19712

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

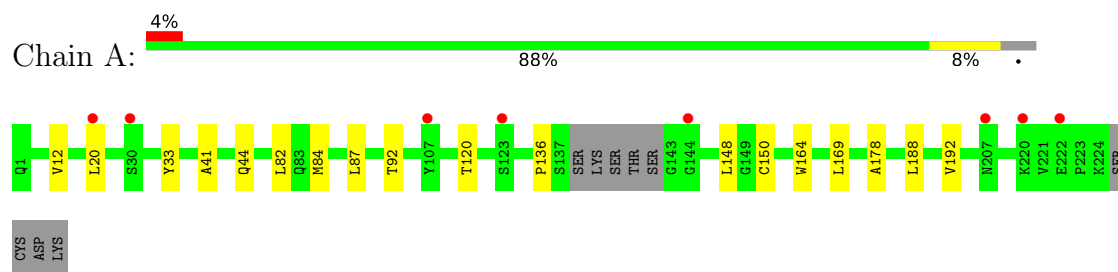


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	K	1	Total	C	N	O	0	0
			14	8	1	5		
5	O	1	Total	C	N	O	0	0
			14	8	1	5		

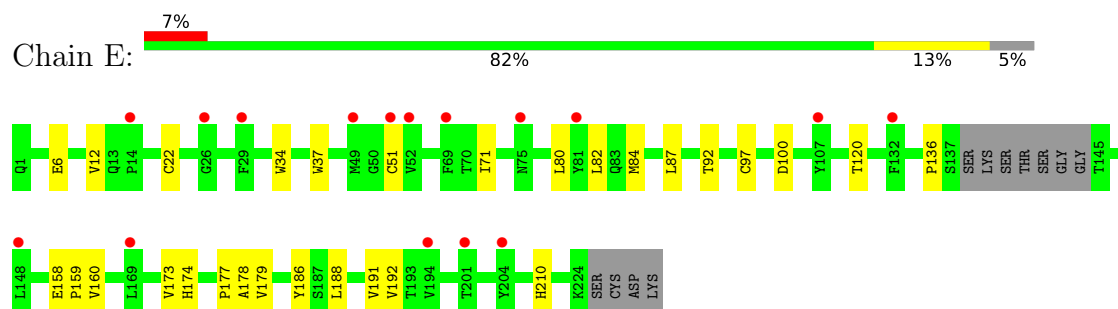
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

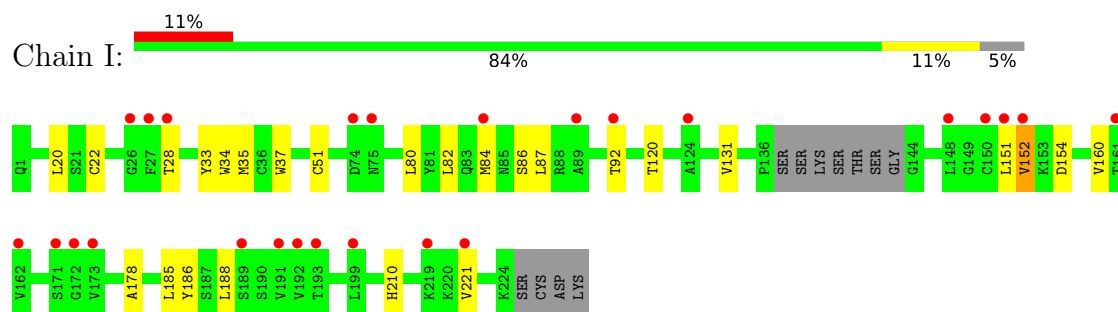
- Molecule 1: DQN0344AE02 Fab heavy chain



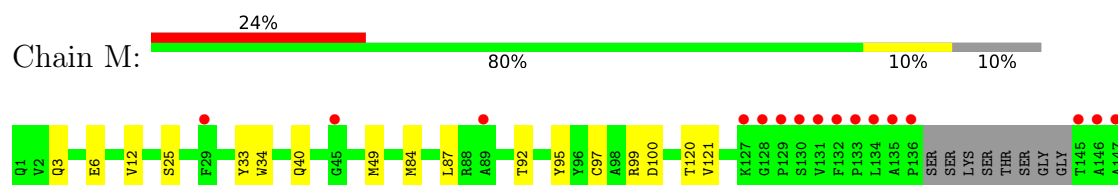
- Molecule 1: DQN0344AE02 Fab heavy chain

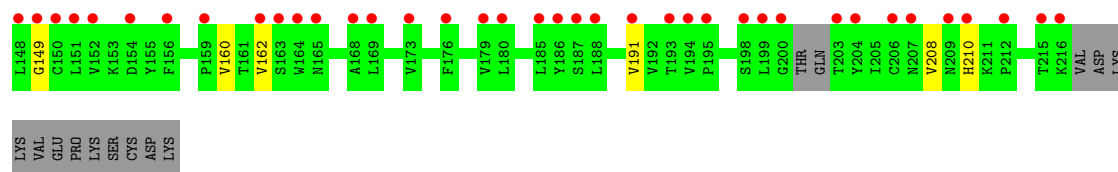


- Molecule 1: DQN0344AE02 Fab heavy chain

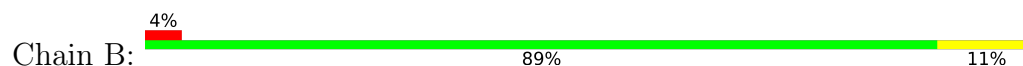


- Molecule 1: DQN0344AE02 Fab heavy chain

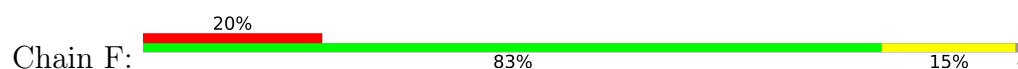




- Molecule 2: DQN0344AE02 Fab light chain



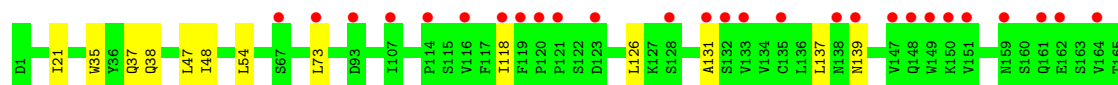
- Molecule 2: DQN0344AE02 Fab light chain



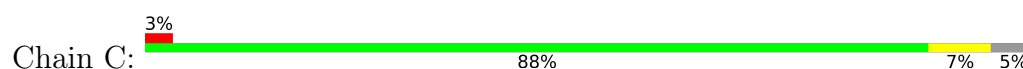
- Molecule 2: DQN0344AE02 Fab light chain



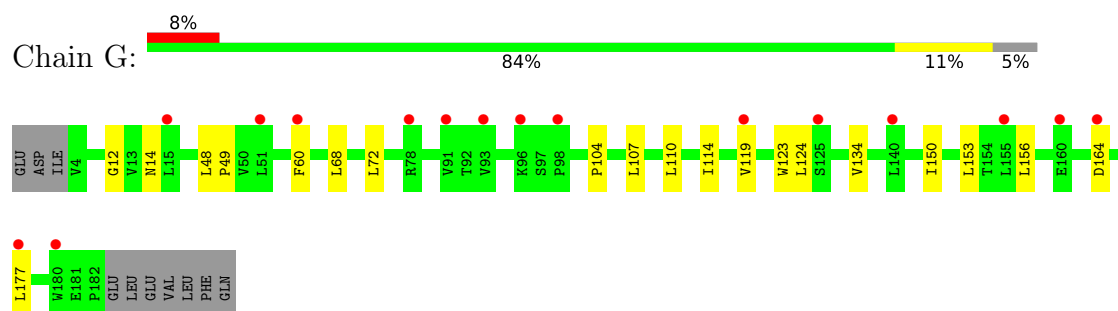
- Molecule 2: DQN0344AE02 Fab light chain



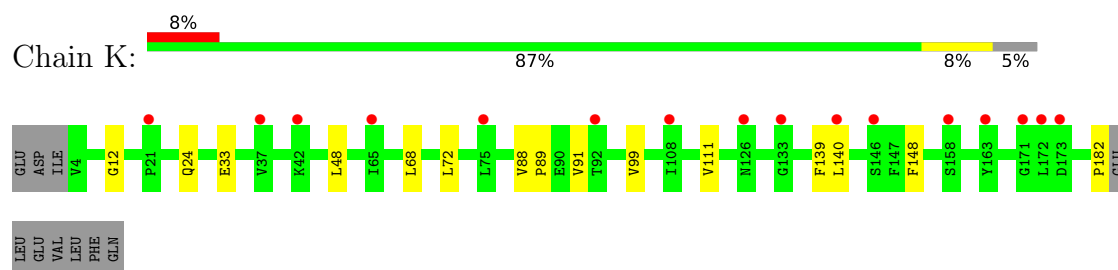
- Molecule 3: HLA class II histocompatibility antigen, DQ alpha 1 chain



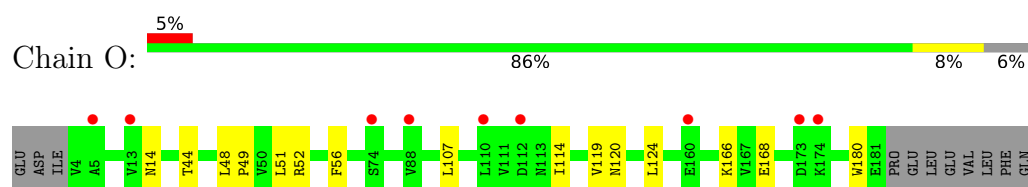
- Molecule 3: HLA class II histocompatibility antigen, DQ alpha 1 chain



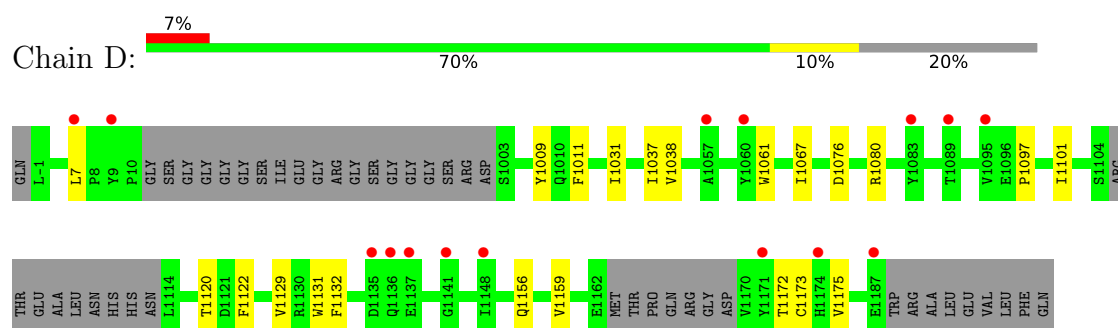
- Molecule 3: HLA class II histocompatibility antigen, DQ alpha 1 chain



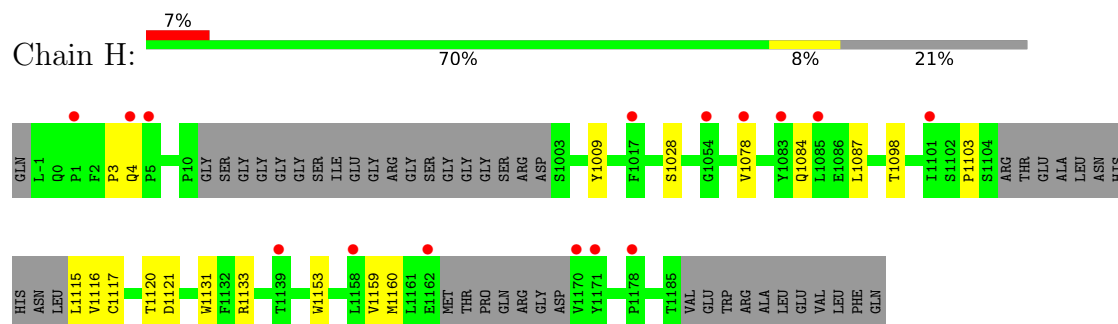
- Molecule 3: HLA class II histocompatibility antigen, DQ alpha 1 chain



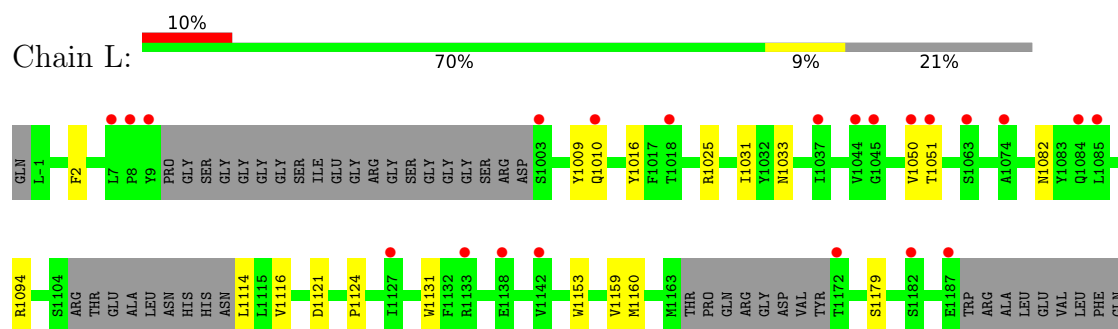
- Molecule 4: MHC class II HLA-DQ-beta-1 - alpha1 gliadin peptide chimeric protein



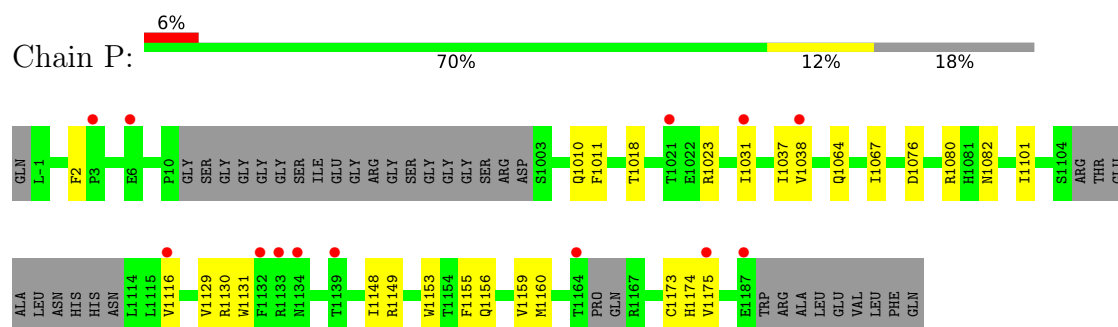
- Molecule 4: MHC class II HLA-DQ-beta-1 - alpha1 gliadin peptide chimeric protein



- Molecule 4: MHC class II HLA-DQ-beta-1 - alpha1 gliadin peptide chimeric protein



- Molecule 4: MHC class II HLA-DQ-beta-1 - alpha1 gliadin peptide chimeric protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.49Å 174.24Å 129.66Å 90.00° 93.17° 90.00°	Depositor
Resolution (Å)	54.86 – 2.82 54.86 – 2.82	Depositor EDS
% Data completeness (in resolution range)	47.7 (54.86-2.82) 46.2 (54.86-2.82)	Depositor EDS
R_{merge}	0.50	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 2.81Å)	Xtriage
Refinement program	BUSTER 2.11.8 (8-JUN-2022)	Depositor
R, R_{free}	0.278 , 0.321 (Not available) , 0.288	Depositor DCC
R_{free} test set	2141 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 19.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	23643	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.45	0/1664	0.77	0/2279
1	E	0.47	0/1649	0.77	0/2260
1	I	0.47	0/1633	0.81	0/2239
1	M	0.48	0/1543	0.77	0/2110
2	B	0.49	0/1655	0.80	0/2253
2	F	0.54	0/1554	0.86	1/2127 (0.0%)
2	J	0.48	0/1583	0.81	2/2163 (0.1%)
2	N	0.48	0/1563	0.80	0/2139
3	C	0.47	0/1449	0.85	0/1985
3	G	0.48	0/1397	0.84	0/1924
3	K	0.47	0/1395	0.87	1/1920 (0.1%)
3	O	0.49	0/1410	0.85	0/1933
4	D	0.48	0/1462	0.79	0/1998
4	H	0.49	0/1381	0.81	0/1893
4	L	0.48	0/1374	0.80	0/1876
4	P	0.48	0/1474	0.80	0/2010
All	All	0.48	0/24186	0.81	4/33109 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	152	ASP	CA-CB-CG	6.69	119.29	112.60
3	K	148	PHE	CA-CB-CG	5.15	118.95	113.80
2	J	142	PRO	CA-C-N	5.08	127.33	120.38
2	J	142	PRO	C-N-CA	5.08	127.33	120.38

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1620	0	1502	12	0
1	E	1605	0	1477	24	0
1	I	1589	0	1446	14	0
1	M	1506	0	1387	14	0
2	B	1620	0	1539	14	0
2	F	1522	0	1392	22	0
2	J	1548	0	1425	7	0
2	N	1529	0	1376	9	0
3	C	1407	0	1330	9	0
3	G	1357	0	1221	16	0
3	K	1354	0	1242	8	0
3	O	1370	0	1278	10	0
4	D	1428	0	1361	15	0
4	H	1347	0	1226	16	0
4	L	1344	0	1234	11	0
4	P	1441	0	1364	17	0
5	C	14	0	13	0	0
5	G	14	0	13	0	0
5	K	14	0	13	0	0
5	O	14	0	13	0	0
All	All	23643	0	21852	196	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (196) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:150:LYS:NZ	2:B:196:GLU:HB2	1.88	0.89
3:C:164:ASP:HB3	3:C:179:HIS:HA	1.56	0.87
1:M:160:VAL:HG22	1:M:210:HIS:CD2	2.11	0.84
1:M:40:GLN:HE22	2:N:38:GLN:HE22	1.22	0.83
3:G:14:ASN:HB3	3:G:68:LEU:HD11	1.66	0.75
4:P:1064:GLN:HB2	4:P:1067:ILE:HB	1.69	0.74
2:B:29:ILE:HB	2:B:92:HIS:HB2	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:37:TRP:HD1	1:E:71:ILE:HD13	1.51	0.74
2:F:151:VAL:HG12	2:F:190:HIS:HD2	1.52	0.72
4:L:2:PHE:H	4:L:1082:ASN:HD21	1.38	0.71
4:P:2:PHE:H	4:P:1082:ASN:HD21	1.40	0.68
4:H:1131:TRP:HE1	4:H:1159:VAL:HG12	1.59	0.68
2:F:114:PRO:HD2	2:F:202:LEU:HG	1.76	0.67
3:K:33:GLU:HB2	3:K:140:LEU:HD11	1.77	0.66
2:B:150:LYS:HZ2	2:B:196:GLU:HB2	1.58	0.65
1:M:160:VAL:HG22	1:M:210:HIS:HD2	1.60	0.65
2:J:149:TRP:CG	2:J:180:LEU:HD13	2.31	0.65
2:N:37:GLN:HB2	2:N:47:LEU:HD11	1.79	0.65
4:P:1101:ILE:HD11	4:P:1173:CYS:HB2	1.80	0.63
2:B:186:ASP:HA	2:B:189:LYS:HD3	1.80	0.62
2:B:37:GLN:HB2	2:B:47:LEU:HD11	1.81	0.62
3:K:48:LEU:HD11	4:L:1153:TRP:CD2	2.35	0.61
1:E:160:VAL:HG12	1:E:210:HIS:ND1	2.16	0.61
3:O:49:PRO:HA	3:O:52:ARG:HE	1.66	0.60
1:E:173:VAL:HG12	1:E:192:VAL:HG22	1.83	0.60
2:F:37:GLN:HB2	2:F:47:LEU:HD21	1.83	0.60
4:D:1097:PRO:HB3	4:D:1122:PHE:HB3	1.82	0.60
1:E:178:ALA:HA	1:E:188:LEU:HB3	1.83	0.60
2:F:65:SER:HB2	2:F:72:THR:HB	1.83	0.59
3:O:49:PRO:HA	3:O:52:ARG:NE	2.17	0.59
4:H:1131:TRP:NE1	4:H:1159:VAL:HG12	2.17	0.59
1:A:20:LEU:HD12	1:A:82:LEU:HD23	1.83	0.59
4:L:1094:ARG:HD3	4:L:1179:SER:HA	1.85	0.59
3:C:145:HIS:CE1	4:D:1031:ILE:HD12	2.38	0.59
2:J:149:TRP:CD1	2:J:180:LEU:HD13	2.38	0.59
4:P:1064:GLN:CB	4:P:1067:ILE:HB	2.33	0.58
3:K:91:VAL:HG22	3:K:111:VAL:HG13	1.85	0.58
4:P:1037:ILE:HG12	4:P:1038:VAL:HG23	1.86	0.57
3:O:124:LEU:HD11	3:O:166:LYS:HD2	1.87	0.57
4:P:1116:VAL:HG22	4:P:1160:MET:HG2	1.87	0.57
2:F:151:VAL:HG12	2:F:190:HIS:CD2	2.36	0.57
2:F:119:PHE:CB	2:F:134:VAL:HB	2.35	0.56
2:B:29:ILE:HB	2:B:92:HIS:CB	2.35	0.56
2:J:121:PRO:HD3	2:J:133:VAL:HG22	1.87	0.56
1:I:131:VAL:HG22	1:I:152:VAL:HG22	1.88	0.56
1:E:22:CYS:HB3	1:E:80:LEU:HB3	1.88	0.55
3:G:107:LEU:HD12	3:G:153:LEU:HD23	1.89	0.55
1:I:84:MET:HB3	1:I:87:LEU:HD21	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:60:PHE:HB3	4:H:3:PRO:HG3	1.89	0.55
4:P:1010:GLN:HB2	4:P:1031:ILE:HB	1.89	0.55
1:I:22:CYS:HB3	1:I:80:LEU:HB3	1.89	0.55
2:N:48:ILE:HG13	2:N:54:LEU:HA	1.89	0.54
1:E:37:TRP:CD1	1:E:71:ILE:HD13	2.37	0.53
3:G:123:TRP:CD1	3:G:134:VAL:HG23	2.43	0.53
2:N:137:LEU:HD21	2:N:197:VAL:HG21	1.90	0.53
1:E:158:GLU:HG2	1:E:159:PRO:HA	1.90	0.53
2:N:191:LYS:HA	2:N:212:ARG:HE	1.73	0.53
4:P:1131:TRP:CD1	4:P:1159:VAL:HG12	2.44	0.53
2:B:146:LYS:HB3	2:B:198:THR:HB	1.89	0.53
4:P:1130:ARG:HB2	4:P:1174:HIS:HB3	1.91	0.53
4:L:1131:TRP:CD1	4:L:1159:VAL:HG12	2.44	0.53
2:F:133:VAL:HB	2:F:180:LEU:HB3	1.90	0.53
3:G:48:LEU:HD12	4:H:1153:TRP:CG	2.44	0.53
4:L:1016:TYR:HB2	4:L:1025:ARG:HB3	1.91	0.52
1:M:162:VAL:HG22	1:M:208:VAL:HG12	1.91	0.52
2:F:11:LEU:HD11	2:F:19:VAL:HG13	1.92	0.52
1:A:169:LEU:HD23	1:A:192:VAL:HG21	1.91	0.52
3:O:107:LEU:HD21	3:O:180:TRP:CD2	2.44	0.52
1:M:84:MET:HB2	1:M:87:LEU:HD21	1.91	0.52
4:P:1076:ASP:HA	4:P:1080:ARG:HB2	1.92	0.52
1:A:33:TYR:HE2	4:D:1067:ILE:HD11	1.75	0.51
1:A:136:PRO:HG3	1:A:148:LEU:HB3	1.92	0.51
3:O:14:ASN:HB2	4:P:1011:PHE:HB3	1.91	0.51
1:M:33:TYR:CD2	1:M:99:ARG:NH2	2.79	0.51
4:L:1010:GLN:HG3	4:L:1031:ILE:HB	1.93	0.51
3:K:99:VAL:HG11	3:K:182:PRO:HB3	1.92	0.51
1:E:178:ALA:HB2	1:E:188:LEU:HD23	1.91	0.51
2:B:147:VAL:HG21	2:B:176:LEU:HD22	1.93	0.51
4:D:1129:VAL:HG22	4:D:1175:VAL:HG22	1.91	0.51
3:O:48:LEU:HD12	4:P:1153:TRP:CE2	2.46	0.51
1:A:178:ALA:HB2	1:A:188:LEU:HD23	1.92	0.51
3:G:48:LEU:HD12	4:H:1153:TRP:CD2	2.46	0.51
3:C:33:GLU:HB2	3:C:140:LEU:HD11	1.92	0.50
2:J:21:ILE:HD12	2:J:73:LEU:HD23	1.92	0.50
4:P:1018:THR:HB	4:P:1023:ARG:HB2	1.92	0.50
2:J:163:SER:OG	2:J:177:SER:OG	2.30	0.50
1:M:49:MET:SD	1:M:95:TYR:HE2	2.35	0.50
3:C:14:ASN:HB2	4:D:1011:PHE:HB3	1.93	0.50
4:D:1131:TRP:CD1	4:D:1159:VAL:HG12	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:VAL:HG11	1:A:87:LEU:HD12	1.94	0.50
2:N:126:LEU:HD13	2:N:131:ALA:HB2	1.94	0.49
1:I:92:THR:HG23	1:I:120:THR:HA	1.95	0.49
2:B:48:ILE:HA	2:B:54:LEU:HA	1.94	0.49
2:F:121:PRO:HG3	2:F:133:VAL:HG22	1.94	0.49
1:I:28:THR:O	1:I:33:TYR:OH	2.29	0.49
3:G:48:LEU:HD12	4:H:1153:TRP:CD1	2.48	0.49
2:J:161:GLN:HB2	2:J:179:THR:HB	1.94	0.48
4:D:1101:ILE:HD11	4:D:1173:CYS:HB2	1.95	0.48
2:N:21:ILE:HD12	2:N:73:LEU:HD23	1.94	0.48
4:D:1076:ASP:HA	4:D:1080:ARG:HB2	1.96	0.48
1:E:174:HIS:CD2	2:F:138:ASN:HD21	2.32	0.48
1:A:41:ALA:HB3	1:A:44:GLN:HB2	1.94	0.47
3:G:72:LEU:HD13	4:H:1009:TYR:HB2	1.95	0.47
4:H:4:GLN:OE1	4:H:1028:SER:OG	2.31	0.47
1:A:178:ALA:HA	1:A:188:LEU:HB3	1.97	0.47
4:H:4:GLN:HG2	4:H:1078:VAL:HG11	1.95	0.47
1:E:34:TRP:HB3	1:E:51:CYS:SG	2.54	0.47
1:E:158:GLU:HG2	1:E:159:PRO:CA	2.45	0.47
1:I:34:TRP:HB3	1:I:51:CYS:SG	2.55	0.47
4:L:1116:VAL:HG12	4:L:1160:MET:HG3	1.97	0.47
2:B:21:ILE:HD12	2:B:73:LEU:HD23	1.97	0.47
3:K:68:LEU:HD13	4:L:1009:TYR:HD2	1.80	0.47
2:B:48:ILE:HG12	2:B:54:LEU:HB3	1.96	0.46
1:E:178:ALA:HB1	1:E:186:TYR:HB3	1.96	0.46
3:O:120:ASN:HB2	3:O:168:GLU:HB2	1.96	0.46
3:G:114:ILE:HD13	3:G:119:VAL:HG11	1.97	0.46
2:F:55:ALA:HB3	2:F:58:ILE:HD13	1.97	0.46
3:G:48:LEU:HD12	4:H:1153:TRP:CE2	2.50	0.46
4:H:1116:VAL:HG12	4:H:1160:MET:HB3	1.96	0.46
3:O:48:LEU:HD13	3:O:51:LEU:HD13	1.97	0.46
4:P:1129:VAL:HG22	4:P:1175:VAL:HG22	1.98	0.46
4:L:1114:LEU:HD11	4:L:1160:MET:HB3	1.97	0.46
3:C:145:HIS:CE1	4:D:1031:ILE:CD1	2.99	0.46
1:I:160:VAL:CG1	1:I:210:HIS:ND1	2.79	0.46
1:M:49:MET:SD	1:M:95:TYR:CE2	3.09	0.46
4:H:1117:CYS:SG	4:H:1131:TRP:CZ2	3.08	0.45
1:A:150:CYS:SG	1:A:164:TRP:CH2	3.09	0.45
1:E:160:VAL:CG1	1:E:210:HIS:ND1	2.80	0.45
4:H:1103:PRO:HA	4:H:1115:LEU:HA	1.98	0.45
2:B:145:ALA:HB2	2:B:199:HIS:HD2	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:123:TRP:CD1	3:G:134:VAL:CG2	2.99	0.45
1:E:37:TRP:CD2	1:E:82:LEU:HD12	2.52	0.45
3:O:44:THR:HG21	3:O:56:PHE:HB3	1.99	0.45
1:I:154:ASP:HA	1:I:185:LEU:CB	2.47	0.45
1:I:20:LEU:HB2	1:I:82:LEU:HB3	1.99	0.45
1:E:12:VAL:HG11	1:E:87:LEU:HD12	2.00	0.44
2:F:94:ILE:HD13	2:F:94:ILE:H	1.80	0.44
1:A:92:THR:HG23	1:A:120:THR:HA	1.98	0.44
4:D:7:LEU:HD21	4:D:1067:ILE:HG21	2.00	0.44
2:N:35:TRP:HB2	2:N:48:ILE:HG22	1.99	0.44
1:A:84:MET:HB3	1:A:87:LEU:HD21	1.99	0.44
4:D:1132:PHE:HB2	4:D:1172:THR:HG23	1.98	0.44
1:E:179:VAL:HG11	2:F:161:GLN:HB3	1.98	0.44
3:O:114:ILE:HD13	3:O:119:VAL:HG11	2.00	0.44
3:C:136:GLU:HA	3:C:151:SER:HA	1.99	0.44
1:I:35:MET:HB3	1:I:80:LEU:HD22	1.98	0.44
1:A:84:MET:HE3	1:A:87:LEU:HD11	1.99	0.44
2:F:151:VAL:CG1	2:F:190:HIS:CD2	3.00	0.44
3:G:124:LEU:HB2	3:G:164:ASP:HB2	1.99	0.44
1:I:37:TRP:HE1	1:I:80:LEU:HG	1.82	0.44
2:F:7:SER:HG	2:F:22:THR:HG1	1.61	0.43
2:F:151:VAL:O	2:F:152:ASP:OD1	2.36	0.43
4:P:1149:ARG:NH2	4:P:1155:PHE:HZ	2.15	0.43
4:P:2:PHE:N	4:P:1082:ASN:HD21	2.12	0.43
1:E:177:PRO:HG2	2:F:163:SER:HB2	2.00	0.43
2:F:33:LEU:HB3	2:F:51:VAL:HG22	2.00	0.43
3:K:72:LEU:HD13	4:L:1009:TYR:HB2	1.98	0.43
3:G:110:LEU:HG	3:G:150:ILE:HG12	2.00	0.43
1:M:3:GLN:HG3	1:M:25:SER:HB3	2.00	0.43
1:M:149:GLY:HA3	1:M:191:VAL:HA	2.00	0.43
2:B:94:ILE:HD13	2:B:94:ILE:H	1.83	0.43
3:C:124:LEU:HD11	3:C:166:LYS:HD2	2.01	0.43
4:D:1037:ILE:HD12	4:D:1038:VAL:HG23	1.99	0.43
2:F:137:LEU:HD13	2:F:176:LEU:HB3	2.01	0.43
1:I:178:ALA:HB1	1:I:186:TYR:HB3	1.99	0.43
1:E:34:TRP:HB2	1:E:100:ASP:O	2.19	0.42
4:D:7:LEU:HD23	4:D:1061:TRP:CD1	2.54	0.42
3:G:48:LEU:HD23	3:G:49:PRO:HD2	2.01	0.42
4:H:1131:TRP:HE1	4:H:1159:VAL:CG1	2.30	0.42
4:H:1084:GLN:HA	4:H:1087:LEU:HD12	2.00	0.42
1:I:154:ASP:HA	1:I:185:LEU:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:1120:THR:HG22	4:D:1156:GLN:HG3	2.02	0.42
1:E:92:THR:HG23	1:E:120:THR:HA	2.00	0.42
1:M:12:VAL:HG23	1:M:121:VAL:HG22	2.01	0.42
4:P:1148:ILE:HB	4:P:1156:GLN:HG3	2.00	0.42
1:E:158:GLU:CG	1:E:159:PRO:HA	2.49	0.42
4:H:1098:THR:OG1	4:H:1120:THR:OG1	2.33	0.42
1:I:151:LEU:O	1:I:188:LEU:O	2.37	0.42
2:J:134:VAL:HG12	2:J:179:THR:HA	2.01	0.42
1:E:84:MET:HB3	1:E:87:LEU:HD21	2.02	0.41
3:G:124:LEU:HD12	3:G:177:LEU:HD11	2.02	0.41
1:M:6:GLU:OE2	1:M:97:CYS:SG	2.78	0.41
1:E:6:GLU:OE2	1:E:97:CYS:SG	2.78	0.41
2:F:30:TYR:HD2	2:F:92:HIS:NE2	2.18	0.41
2:N:118:ILE:HB	2:N:208:LYS:HB3	2.01	0.41
2:B:129:GLY:HA2	2:B:184:LYS:HB2	2.01	0.41
3:G:104:PRO:HB3	3:G:156:LEU:HG	2.01	0.41
3:C:91:VAL:HB	3:C:178:LYS:HD3	2.02	0.41
1:E:174:HIS:CD2	2:F:138:ASN:ND2	2.89	0.41
3:K:24:GLN:HE21	3:K:24:GLN:HB3	1.71	0.41
1:E:191:VAL:HG21	2:F:136:LEU:HD11	2.03	0.41
3:K:88:VAL:HA	3:K:89:PRO:HD3	1.97	0.41
1:M:34:TRP:HB2	1:M:100:ASP:O	2.21	0.41
4:L:1050:VAL:HG22	4:L:1051:THR:HG23	2.03	0.40
1:M:92:THR:HG23	1:M:120:THR:HA	2.03	0.40
3:C:72:LEU:HD13	4:D:1009:TYR:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	102/228 (45%)	98 (96%)	4 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	213/228 (93%)	193 (91%)	19 (9%)	1 (0%)	24	53
1	I	213/228 (93%)	192 (90%)	18 (8%)	3 (1%)	9	27
1	M	200/228 (88%)	190 (95%)	10 (5%)	0	100	100
2	B	174/215 (81%)	165 (95%)	7 (4%)	2 (1%)	11	34
2	F	210/215 (98%)	186 (89%)	19 (9%)	5 (2%)	4	16
2	J	207/215 (96%)	188 (91%)	16 (8%)	3 (1%)	9	27
2	N	210/215 (98%)	191 (91%)	18 (9%)	1 (0%)	24	53
3	C	177/189 (94%)	170 (96%)	7 (4%)	0	100	100
3	G	177/189 (94%)	166 (94%)	10 (6%)	1 (1%)	21	48
3	K	177/189 (94%)	171 (97%)	4 (2%)	2 (1%)	11	34
3	O	176/189 (93%)	168 (96%)	8 (4%)	0	100	100
4	D	57/226 (25%)	53 (93%)	4 (7%)	0	100	100
4	H	170/226 (75%)	156 (92%)	12 (7%)	2 (1%)	10	31
4	L	171/226 (76%)	157 (92%)	11 (6%)	3 (2%)	6	22
4	P	178/226 (79%)	169 (95%)	9 (5%)	0	100	100
All	All	2812/3432 (82%)	2613 (93%)	176 (6%)	23 (1%)	16	41

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	141	TYR
2	B	139	ASN
2	F	158	GLY
4	H	1121	ASP
2	J	139	ASN
4	L	1121	ASP
2	F	30	TYR
4	H	1133	ARG
1	I	86	SER
1	I	152	VAL
3	K	12	GLY
3	K	139	PHE
2	B	144	GLU
1	I	221	VAL
2	J	144	GLU
2	N	139	ASN
2	F	183	SER

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Mol	Chain	Res	Type
2	J	30	TYR
4	L	1033	ASN
2	F	152	ASP
3	G	12	GLY
1	E	136	PRO
4	L	1124	PRO

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	G	1000	-	14,14,15	0.24	0	17,19,21	0.43	0
5	NAG	K	1000	-	14,14,15	0.25	0	17,19,21	0.40	0
5	NAG	O	1000	-	14,14,15	0.41	0	17,19,21	0.31	0
5	NAG	C	1000	-	14,14,15	0.24	0	17,19,21	0.45	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	G	1000	-	-	1/6/23/26	0/1/1/1
5	NAG	K	1000	-	-	0/6/23/26	0/1/1/1
5	NAG	O	1000	-	-	0/6/23/26	0/1/1/1
5	NAG	C	1000	-	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	G	1000	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	219/228 (96%)	0.71	8 (3%) 45 36	24, 30, 41, 50	0
1	E	217/228 (95%)	0.92	16 (7%) 20 15	20, 34, 56, 66	0
1	I	217/228 (95%)	0.99	25 (11%) 9 7	31, 38, 50, 56	0
1	M	206/228 (90%)	1.36	54 (26%) 1 1	26, 37, 86, 93	0
2	B	214/215 (99%)	0.69	9 (4%) 40 32	23, 36, 48, 55	0
2	F	212/215 (98%)	1.26	42 (19%) 3 2	26, 42, 60, 69	0
2	J	209/215 (97%)	0.96	14 (6%) 24 17	30, 39, 47, 51	0
2	N	212/215 (98%)	1.39	49 (23%) 2 1	34, 51, 73, 82	0
3	C	179/189 (94%)	0.54	5 (2%) 55 44	15, 31, 40, 45	0
3	G	179/189 (94%)	0.97	16 (8%) 15 11	23, 36, 50, 52	0
3	K	179/189 (94%)	0.92	16 (8%) 15 11	26, 40, 53, 58	0
3	O	178/189 (94%)	0.84	9 (5%) 33 25	28, 39, 54, 64	0
4	D	181/226 (80%)	0.96	15 (8%) 17 12	27, 32, 43, 50	0
4	H	178/226 (78%)	1.00	15 (8%) 17 12	26, 35, 46, 53	0
4	L	179/226 (79%)	0.98	22 (12%) 8 6	29, 41, 55, 58	0
4	P	186/226 (82%)	0.94	13 (6%) 22 16	28, 38, 50, 56	0
All	All	3145/3432 (91%)	0.97	328 (10%) 11 8	15, 37, 67, 93	0

All (328) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	151	LEU	6.5
1	M	200	GLY	5.1
1	E	201	THR	5.1
2	N	207	THR	5.1
4	D	1136	GLN	4.6

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Mol	Chain	Res	Type	RSRZ
2	F	153	ASN	4.6
2	N	159	ASN	4.5
1	M	133	PRO	4.4
1	M	199	LEU	4.2
4	L	9	TYR	4.1
2	F	129	GLY	4.1
1	M	191	VAL	4.1
3	O	110	LEU	4.0
1	M	131	VAL	3.9
4	H	1101	ILE	3.9
1	M	187	SER	3.9
2	N	128	SER	3.8
2	N	116	VAL	3.8
2	F	207	THR	3.8
2	N	147	VAL	3.7
3	C	99	VAL	3.7
4	D	1135	ASP	3.6
3	G	164	ASP	3.6
1	M	135	ALA	3.5
2	N	118	ILE	3.5
1	M	162	VAL	3.5
4	H	1162	GLU	3.5
2	N	193	TYR	3.5
2	N	132	SER	3.4
2	N	139	ASN	3.4
1	A	144	GLY	3.4
1	M	204	TYR	3.4
1	M	147	ALA	3.4
1	M	176	PHE	3.4
1	M	129	PRO	3.4
1	M	216	LYS	3.3
2	F	132	SER	3.3
2	J	164	VAL	3.2
1	M	145	THR	3.2
2	N	149	TRP	3.2
1	I	151	LEU	3.2
3	O	5	ALA	3.2
4	P	1187	GLU	3.2
3	G	98	PRO	3.2
4	D	1141	GLY	3.1
1	M	136	PRO	3.1
1	I	148	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	I	191	VAL	3.1
4	P	6	GLU	3.1
2	B	195	CYS	3.1
1	M	134	LEU	3.1
2	F	140	PHE	3.0
1	M	198	SER	3.0
2	N	119	PHE	3.0
2	F	34	ALA	3.0
1	M	152	VAL	3.0
1	M	168	ALA	3.0
1	M	179	VAL	3.0
4	L	1050	VAL	3.0
2	F	212	ARG	2.9
4	L	1084	GLN	2.9
1	M	132	PHE	2.9
1	M	194	VAL	2.9
2	F	138	ASN	2.9
3	O	174	LYS	2.9
2	F	159	ASN	2.9
1	M	146	ALA	2.9
2	N	178	SER	2.9
1	M	45	GLY	2.9
2	N	208	LYS	2.9
1	I	75	ASN	2.9
4	P	1038	VAL	2.9
1	I	193	THR	2.9
2	J	181	THR	2.9
4	L	1045	GLY	2.8
1	E	26	GLY	2.8
4	D	1187	GLU	2.8
2	N	131	ALA	2.8
1	I	84	MET	2.8
1	I	192	VAL	2.8
2	F	141	TYR	2.8
4	D	1171	TYR	2.8
1	M	169	LEU	2.8
1	M	127	LYS	2.8
1	M	164	TRP	2.7
1	M	180	LEU	2.7
2	F	136	LEU	2.7
2	F	176	LEU	2.7
4	D	9	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
4	D	1083	TYR	2.7
4	L	1037	ILE	2.7
1	I	171	SER	2.7
1	A	222	GLU	2.7
2	N	151	VAL	2.7
2	N	206	VAL	2.7
1	A	220	LYS	2.7
2	N	150	LYS	2.7
1	M	154	ASP	2.7
1	E	132	PHE	2.7
2	N	135	CYS	2.7
1	M	195	PRO	2.7
2	F	204	SER	2.7
2	F	147	VAL	2.7
3	G	51	LEU	2.7
1	M	150	CYS	2.7
4	D	1060	TYR	2.7
2	F	211	ASN	2.7
2	N	173	THR	2.7
3	K	21	PRO	2.7
1	E	194	VAL	2.7
1	I	221	VAL	2.7
1	M	163	SER	2.7
3	K	158	SER	2.7
1	A	207	ASN	2.6
2	N	176	LEU	2.6
4	P	1031	ILE	2.6
4	P	1164	THR	2.6
2	J	197	VAL	2.6
4	H	1158	LEU	2.6
3	G	125	SER	2.6
4	L	1085	LEU	2.6
4	P	1132	PHE	2.6
3	K	173	ASP	2.6
3	G	180	TRP	2.6
1	E	107	TYR	2.6
1	E	52	VAL	2.6
1	I	173	VAL	2.6
3	K	92	THR	2.6
2	N	200	GLN	2.6
2	F	71	PHE	2.6
3	C	51	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
4	D	1089	THR	2.5
1	M	210	HIS	2.5
2	F	203	SER	2.5
1	M	149	GLY	2.5
3	K	75	LEU	2.5
4	D	7	LEU	2.5
2	F	194	ALA	2.5
2	F	88	CYS	2.5
1	M	159	PRO	2.5
2	N	120	PRO	2.5
3	G	160	GLU	2.5
4	L	1172	THR	2.5
4	P	1133	ARG	2.5
1	I	150	CYS	2.5
1	I	219	LYS	2.5
3	K	171	GLY	2.5
1	E	204	TYR	2.5
1	M	156	PHE	2.5
2	N	161	GLN	2.5
1	I	28	THR	2.5
2	N	179	THR	2.5
1	I	199	LEU	2.4
3	O	112	ASP	2.4
2	F	151	VAL	2.4
2	N	197	VAL	2.4
3	O	160	GLU	2.4
4	H	1017	PHE	2.4
1	M	148	LEU	2.4
1	M	188	LEU	2.4
3	G	96	LYS	2.4
2	J	200	GLN	2.4
1	I	161	THR	2.4
2	J	78	LEU	2.4
2	N	107	ILE	2.4
1	E	51	CYS	2.4
2	N	114	PRO	2.4
1	I	172	GLY	2.4
1	M	130	SER	2.4
2	F	7	SER	2.4
4	L	1187	GLU	2.4
4	D	1057	ALA	2.4
3	C	163	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
4	H	1139	THR	2.4
2	F	202	LEU	2.4
3	G	140	LEU	2.4
2	N	138	ASN	2.4
4	L	8	PRO	2.4
1	E	69	PHE	2.4
2	B	214	GLU	2.4
2	N	166	GLU	2.4
2	F	25	ALA	2.4
2	N	175	SER	2.4
2	F	50	TYR	2.4
3	G	155	LEU	2.4
2	F	116	VAL	2.3
2	N	180	LEU	2.3
4	H	1171	TYR	2.3
4	D	1095	VAL	2.3
2	F	168	ASP	2.3
4	H	4	GLN	2.3
2	F	175	SER	2.3
2	F	134	VAL	2.3
1	I	74	ASP	2.3
3	O	173	ASP	2.3
2	F	161	GLN	2.3
2	B	204	SER	2.3
3	K	146	SER	2.3
4	H	1054	GLY	2.3
1	E	29	PHE	2.3
1	M	207	ASN	2.3
4	P	1134	ASN	2.3
1	E	49	MET	2.3
2	B	96	GLU	2.3
1	A	123	SER	2.3
2	F	114	PRO	2.3
4	P	3	PRO	2.3
1	M	29	PHE	2.2
1	M	165	ASN	2.2
1	M	209	ASN	2.2
2	N	73	LEU	2.2
2	N	148	GLN	2.2
1	M	206	CYS	2.2
4	L	1051	THR	2.2
1	A	107	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
2	N	67	SER	2.2
4	L	1044	VAL	2.2
3	K	163	TYR	2.2
3	O	74	SER	2.2
1	M	128	GLY	2.2
2	J	176	LEU	2.2
2	N	189	LYS	2.2
2	F	152	ASP	2.2
1	M	215	THR	2.2
2	F	197	VAL	2.2
3	O	88	VAL	2.2
4	P	1021	THR	2.2
1	M	186	TYR	2.2
2	J	95	SER	2.2
4	L	1003	SER	2.2
4	L	1182	SER	2.2
2	J	119	PHE	2.2
2	F	137	LEU	2.2
2	F	180	LEU	2.2
3	K	65	ILE	2.2
2	N	164	VAL	2.2
1	M	203	THR	2.2
2	F	46	LEU	2.2
2	N	182	LEU	2.2
1	M	173	VAL	2.2
2	N	177	SER	2.2
3	C	84	ALA	2.2
4	L	1138	GLU	2.2
4	L	1133	ARG	2.1
1	E	14	PRO	2.1
2	N	133	VAL	2.1
3	G	91	VAL	2.1
4	H	5	PRO	2.1
4	H	1078	VAL	2.1
2	B	99	PHE	2.1
3	G	15	LEU	2.1
4	H	1085	LEU	2.1
2	J	143	ARG	2.1
3	K	108	ILE	2.1
4	D	1174	HIS	2.1
2	F	192	VAL	2.1
4	P	1175	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	I	92	THR	2.1
1	M	185	LEU	2.1
3	K	140	LEU	2.1
3	K	172	LEU	2.1
2	F	187	TYR	2.1
4	L	1074	ALA	2.1
1	A	30	SER	2.1
4	L	1010	GLN	2.1
3	K	126	ASN	2.1
4	H	1178	PRO	2.1
1	I	26	GLY	2.1
2	B	126	LEU	2.1
2	N	123	ASP	2.1
2	N	181	THR	2.1
3	G	78	ARG	2.1
1	M	89	ALA	2.1
4	D	1148	ILE	2.1
4	H	1083	TYR	2.1
3	G	93	VAL	2.1
3	G	119	VAL	2.1
3	K	42	LYS	2.1
3	O	13	VAL	2.1
4	H	1170	VAL	2.1
4	L	1142	VAL	2.1
1	E	75	ASN	2.1
1	E	148	LEU	2.1
3	G	177	LEU	2.1
2	B	201	GLY	2.1
2	F	210	PHE	2.1
2	J	66	GLY	2.1
1	M	193	THR	2.1
4	P	1139	THR	2.1
1	I	89	ALA	2.1
2	N	174	TYR	2.1
2	F	184	LYS	2.0
2	J	209	SER	2.0
2	J	88	CYS	2.0
3	K	37	VAL	2.0
1	E	169	LEU	2.0
2	J	28	ASN	2.0
4	L	7	LEU	2.0
3	G	60	PHE	2.0

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Mol	Chain	Res	Type	RSRZ
2	F	179	THR	2.0
4	L	1018	THR	2.0
1	I	124	ALA	2.0
2	J	60	ALA	2.0
2	N	93	ASP	2.0
4	L	1127	ILE	2.0
1	E	81	TYR	2.0
1	I	152	VAL	2.0
1	I	189	SER	2.0
2	F	133	VAL	2.0
2	N	167	GLN	2.0
4	L	1063	SER	2.0
4	P	1116	VAL	2.0
1	M	212	PRO	2.0
2	N	121	PRO	2.0
4	H	1	PRO	2.0
1	A	20	LEU	2.0
2	N	202	LEU	2.0
3	C	110	LEU	2.0
1	I	27	PHE	2.0
2	N	210	PHE	2.0
3	K	133	GLY	2.0
2	B	198	THR	2.0
2	F	130	THR	2.0
2	N	185	ALA	2.0
2	B	39	LYS	2.0
2	N	162	GLU	2.0
4	D	1137	GLU	2.0
1	I	162	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	G	1000	14/15	0.59	0.18	62,62,62,62	0
5	NAG	O	1000	14/15	0.63	0.18	64,64,64,64	0
5	NAG	C	1000	14/15	0.78	0.21	52,52,52,52	0
5	NAG	K	1000	14/15	0.81	0.14	57,57,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.