



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 06:53 am GMT

PDB ID : 4L29
Title : Structure of wtMHC class I with NY-ESO1 double mutant
Authors : Halabelian, L.; Giorgetti, S.; Bellotti, V.; Bolognesi, M.; Ricagno, S.
Deposited on : 2013-06-04
Resolution : 3.09 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

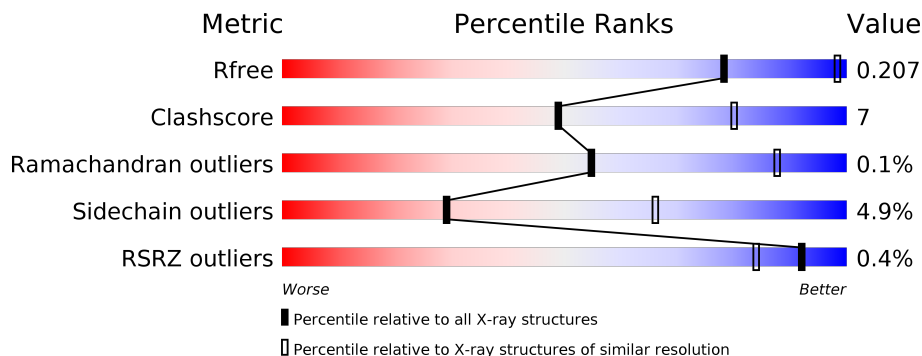
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.09 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.

























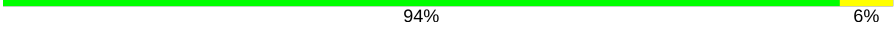

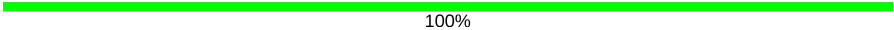
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1001 (3.12-3.08)
Clashscore	112137	1099 (3.12-3.08)
Ramachandran outliers	110173	1057 (3.12-3.08)
Sidechain outliers	110143	1057 (3.12-3.08)
RSRZ outliers	101464	1006 (3.12-3.08)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	276	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between;"> 72% 27% </div> </div>
1	C	276	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="display: flex; justify-content: space-between;"> 82% 17% </div> </div>
1	E	276	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="display: flex; justify-content: space-between;"> 83% 16% </div> </div>
1	G	276	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="display: flex; justify-content: space-between;"> 85% 14% </div> </div>
1	I	276	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="display: flex; justify-content: space-between;"> 83% 14% </div> </div>
1	K	276	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="display: flex; justify-content: space-between;"> 80% 19% </div> </div>

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Mol	Chain	Length	Quality of chain
1	M	276	 83% 16% .
1	O	276	 84% 15% .
1	Q	276	 90% 9% .
1	S	276	 82% 17% .
1	U	276	 80% 19% .
1	W	276	 82% 17% .
1	Y	276	 76% 22% .
1	a	276	 96% .
2	B	100	 68% 31% .
2	D	100	 86% 14% .
2	F	100	 84% 15% .
2	H	100	 82% 18% .
2	J	100	 85% 15% .
2	L	100	 91% 8% .
2	N	100	 84% 15% .
2	P	100	 82% 18% .
2	R	100	 78% 20% .
2	T	100	 81% 17% .
2	V	100	 82% 17% .
2	X	100	 83% 17% .
2	Z	100	 88% 10% .
2	b	100	 94% 6% .
3	c	9	 89% 11% .
3	e	9	 100% .
3	f	9	 89% 11% .

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Mol	Chain	Length	Quality of chain
3	g	9	 100%
3	h	9	 89% 11%
3	i	9	 100%
3	j	9	 89% 11%
3	k	9	 100%
3	l	9	 100%
3	m	9	 89% 11%
3	n	9	 89% 11%
3	o	9	 100%
3	p	9	 100%
3	q	9	 100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GOL	O	302	-	-	-	X
5	GOL	i	101	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 44627 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 HLA class I histocompatibility antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	C	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	E	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	G	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	I	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	K	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	M	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	O	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	Q	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	S	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	U	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	W	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	Y	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	a	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	D	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	F	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	H	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	J	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	L	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	N	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	P	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	R	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	T	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	V	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	X	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	Z	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	b	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	INITIATING METHIONINE	UNP P61769
D	0	MET	-	INITIATING METHIONINE	UNP P61769
F	0	MET	-	INITIATING METHIONINE	UNP P61769
H	0	MET	-	INITIATING METHIONINE	UNP P61769
J	0	MET	-	INITIATING METHIONINE	UNP P61769
L	0	MET	-	INITIATING METHIONINE	UNP P61769
N	0	MET	-	INITIATING METHIONINE	UNP P61769
P	0	MET	-	INITIATING METHIONINE	UNP P61769
R	0	MET	-	INITIATING METHIONINE	UNP P61769
T	0	MET	-	INITIATING METHIONINE	UNP P61769
V	0	MET	-	INITIATING METHIONINE	UNP P61769
X	0	MET	-	INITIATING METHIONINE	UNP P61769

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Chain	Residue	Modelled	Actual	Comment	Reference
Z	0	MET	-	INITIATING METHIONINE	UNP P61769
b	0	MET	-	INITIATING METHIONINE	UNP P61769

- Molecule 3 is a protein called NY-ESO1 DOUBLE MUTANT (1Y, 9V).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	m	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	i	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	k	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	f	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	l	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	h	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	e	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	n	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	p	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	o	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	c	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	g	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	q	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			
3	j	9	Total	C	N	O	S	0	0	0
			82	57	11	13	1			

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	P	1	Total	Cl	0	0
			1	1		
4	K	1	Total	Cl	0	0
			1	1		

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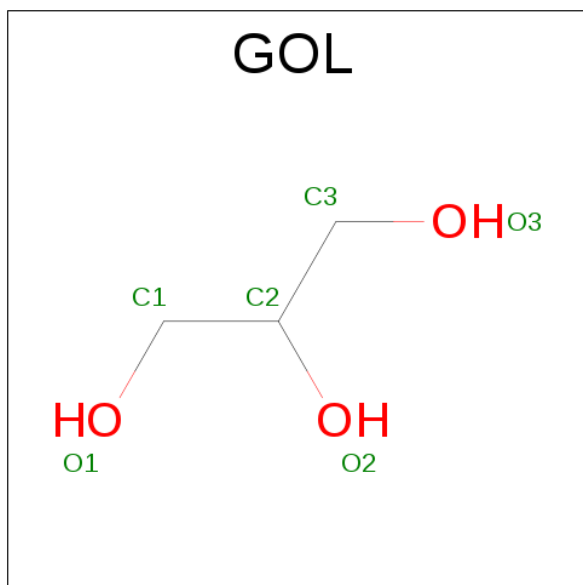
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Cl 1	0	0
4	W	1	Total 1	Cl 1	0	0
4	N	2	Total 2	Cl 2	0	0
4	S	1	Total 1	Cl 1	0	0
4	J	1	Total 1	Cl 1	0	0
4	E	1	Total 1	Cl 1	0	0
4	b	1	Total 1	Cl 1	0	0
4	V	1	Total 1	Cl 1	0	0
4	A	1	Total 1	Cl 1	0	0
4	R	1	Total 1	Cl 1	0	0
4	D	1	Total 1	Cl 1	0	0
4	I	3	Total 3	Cl 3	0	0
4	Z	1	Total 1	Cl 1	0	0
4	a	1	Total 1	Cl 1	0	0
4	U	1	Total 1	Cl 1	0	0
4	L	1	Total 1	Cl 1	0	0
4	G	1	Total 1	Cl 1	0	0
4	Q	2	Total 2	Cl 2	0	0
4	H	1	Total 1	Cl 1	0	0
4	C	1	Total 1	Cl 1	0	0
4	T	1	Total 1	Cl 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	O	1	Total	Cl	0	0
			1	1		
4	Y	1	Total	Cl	0	0
			1	1		
4	F	1	Total	Cl	0	0
			1	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	i	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		
5	f	1	Total	C	O	0	0
			6	3	3		
5	l	1	Total	C	O	0	0
			6	3	3		
5	K	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	h	1	Total C O 6 3 3	0	0
5	e	1	Total C O 6 3 3	0	0
5	O	1	Total C O 6 3 3	0	0
5	O	1	Total C O 6 3 3	0	0
5	Q	1	Total C O 6 3 3	0	0
5	S	1	Total C O 6 3 3	0	0
5	o	1	Total C O 6 3 3	0	0
5	c	1	Total C O 6 3 3	0	0
5	g	1	Total C O 6 3 3	0	0
5	Y	1	Total C O 6 3 3	0	0
5	q	1	Total C O 6 3 3	0	0
5	a	1	Total C O 6 3 3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	2	Total O 2 2	0	0
6	C	4	Total O 4 4	0	0
6	D	2	Total O 2 2	0	0
6	E	3	Total O 3 3	0	0
6	F	2	Total O 2 2	0	0
6	G	7	Total O 7 7	0	0
6	H	1	Total O 1 1	0	0

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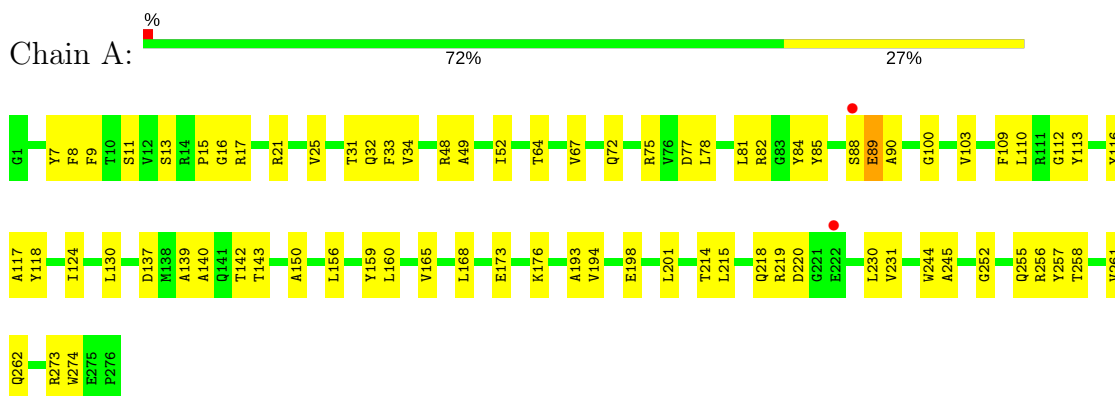
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	I	3	Total 3	O 3	0	0
6	J	2	Total 2	O 2	0	0
6	K	3	Total 3	O 3	0	0
6	L	3	Total 3	O 3	0	0
6	M	1	Total 1	O 1	0	0
6	P	3	Total 3	O 3	0	0
6	Q	2	Total 2	O 2	0	0
6	R	3	Total 3	O 3	0	0
6	S	1	Total 1	O 1	0	0
6	T	1	Total 1	O 1	0	0
6	U	3	Total 3	O 3	0	0
6	V	2	Total 2	O 2	0	0
6	X	3	Total 3	O 3	0	0
6	Y	2	Total 2	O 2	0	0
6	Z	2	Total 2	O 2	0	0

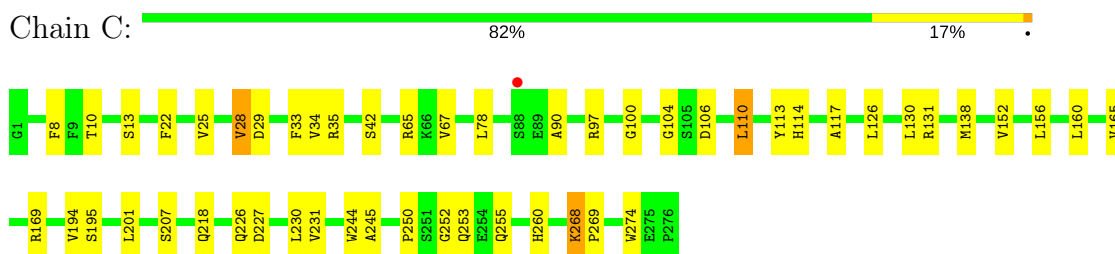
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

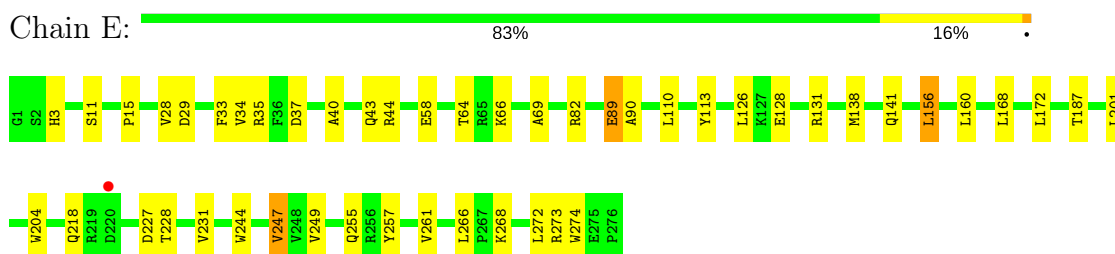
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



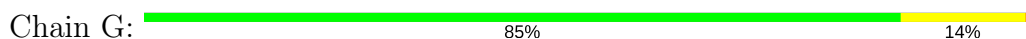
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain

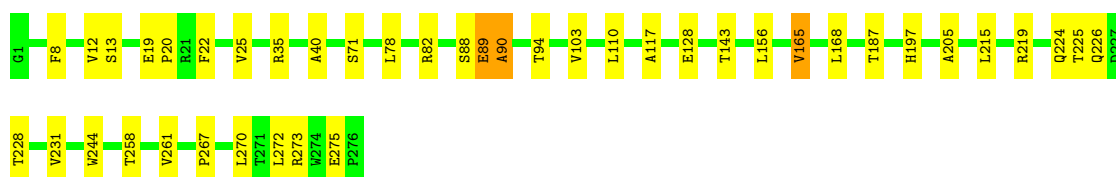


- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain

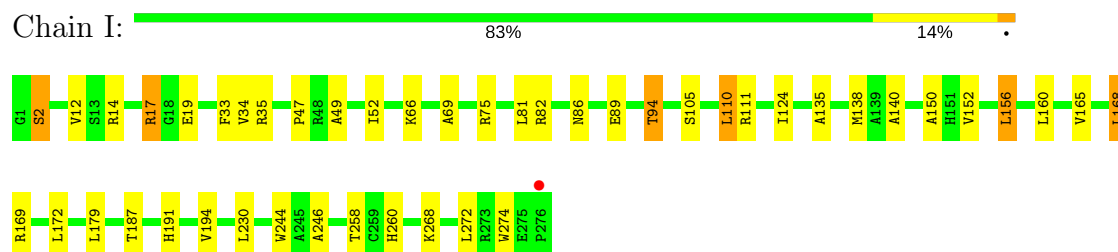


- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain

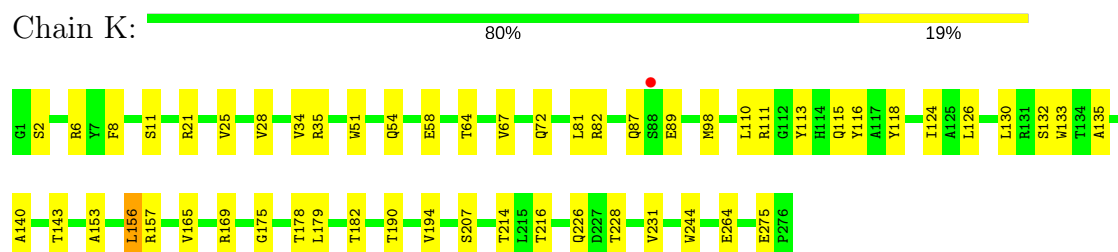




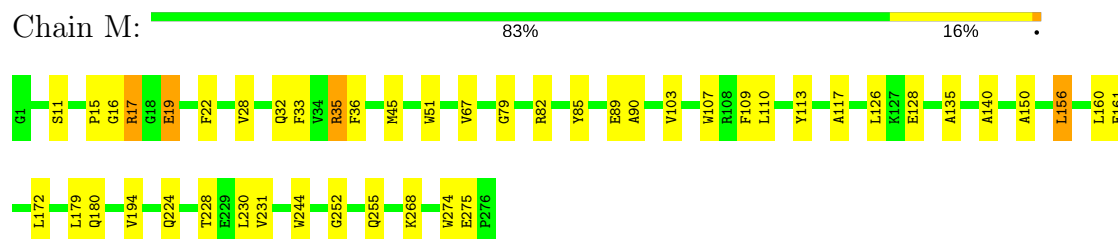
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



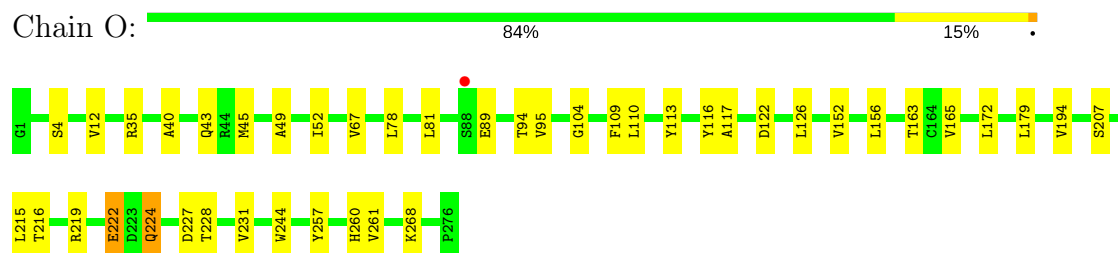
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



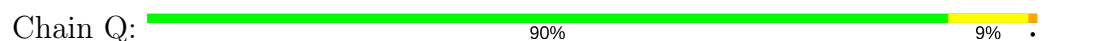
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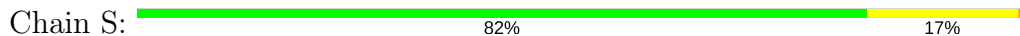


- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain





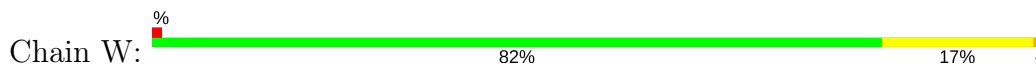
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



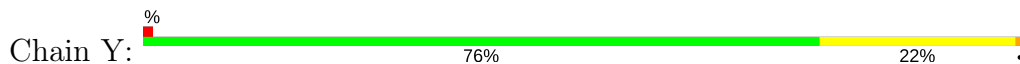
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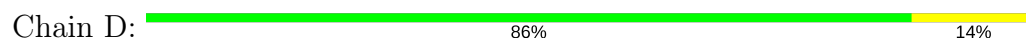




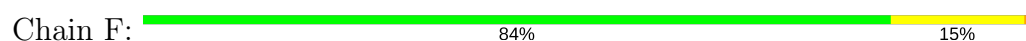
- Molecule 2: Beta-2-microglobulin



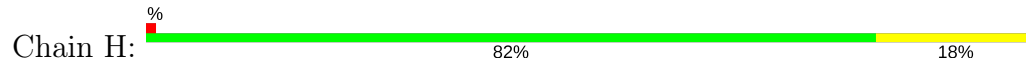
- Molecule 2: Beta-2-microglobulin



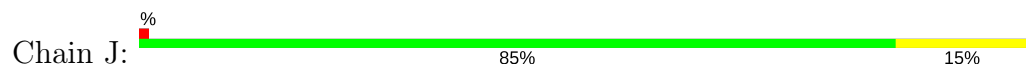
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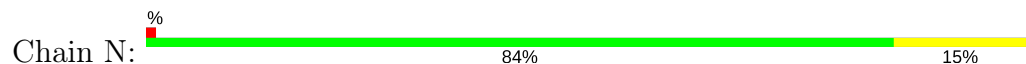
- Molecule 2: Beta-2-microglobulin

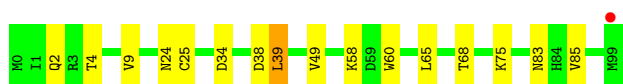


- Molecule 2: Beta-2-microglobulin

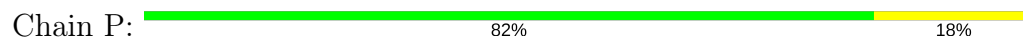


- Molecule 2: Beta-2-microglobulin

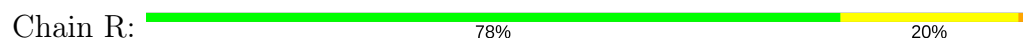




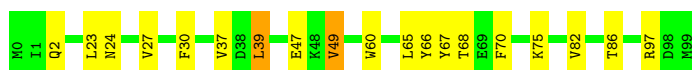
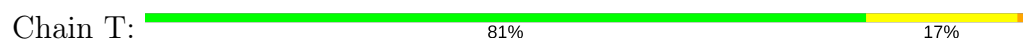
- Molecule 2: Beta-2-microglobulin



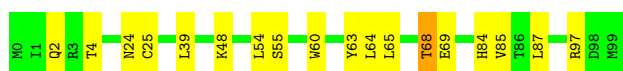
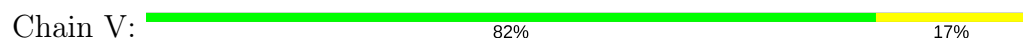
- Molecule 2: Beta-2-microglobulin



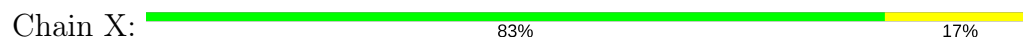
- Molecule 2: Beta-2-microglobulin



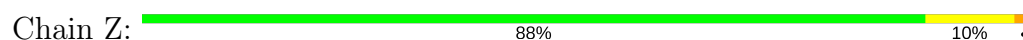
- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin




- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain m:  89% 11%



- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain i:  100%


There are no outlier residues recorded for this chain.

- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain k:  100%

There are no outlier residues recorded for this chain.

- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain f:  89% 11%



- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain l:  100%

There are no outlier residues recorded for this chain.

- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain h:  89% 11%




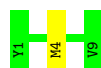
- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain e:  100%

There are no outlier residues recorded for this chain.

- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain n:  89% 11%



- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain p:  100%


There are no outlier residues recorded for this chain.

- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain o:  100%

There are no outlier residues recorded for this chain.

- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain c:  89% 11%



- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain g:  100%


There are no outlier residues recorded for this chain.

- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain q:  100%

There are no outlier residues recorded for this chain.

- Molecule 3: NY-ESO1 DOUBLE MUTANT (1Y, 9V)

Chain j:  89% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.57Å 313.36Å 314.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	53.80 – 3.09 53.88 – 3.09	Depositor EDS
% Data completeness (in resolution range)	98.8 (53.80-3.09) 99.1 (53.88-3.09)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 3.07Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.171 , 0.201 0.185 , 0.207	Depositor DCC
R_{free} test set	9163 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	65.6	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for -h,l,k	Xtriage
Reported twinning fraction	0.671 for H, K, L 0.329 for -H, L, K	Depositor
Outliers	1 of 182843 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	44627	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.49	0/2320	0.64	0/3149
1	C	0.58	1/2320 (0.0%)	0.65	0/3149
1	E	0.49	0/2320	0.64	0/3149
1	G	0.52	1/2320 (0.0%)	0.61	0/3149
1	I	0.51	0/2320	0.61	0/3149
1	K	0.49	0/2320	0.63	0/3149
1	M	0.48	0/2320	0.62	0/3149
1	O	0.46	0/2320	0.60	0/3149
1	Q	0.52	0/2320	0.64	0/3149
1	S	0.47	0/2320	0.61	0/3149
1	U	0.47	0/2320	0.58	0/3149
1	W	0.52	1/2320 (0.0%)	0.61	0/3149
1	Y	0.47	0/2320	0.60	0/3149
1	a	0.44	0/2320	0.59	0/3149
2	B	0.47	0/860	0.59	0/1162
2	D	0.53	0/860	0.63	0/1162
2	F	0.50	0/860	0.59	0/1162
2	H	0.48	0/860	0.60	0/1162
2	J	0.50	0/860	0.57	0/1162
2	L	0.50	0/860	0.60	0/1162
2	N	0.50	0/860	0.61	0/1162
2	P	0.46	0/860	0.58	0/1162
2	R	0.47	0/860	0.56	0/1162
2	T	0.56	1/860 (0.1%)	0.63	0/1162
2	V	0.49	0/860	0.58	0/1162
2	X	0.48	0/860	0.59	0/1162
2	Z	0.45	0/860	0.58	0/1162
2	b	0.46	0/860	0.56	0/1162
3	c	0.51	0/84	0.55	0/113
3	e	0.43	0/84	0.58	0/113
3	f	0.45	0/84	0.50	0/113
3	g	0.46	0/84	0.51	0/113

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
3	h	0.48	0/84	0.58	0/113
3	i	0.46	0/84	0.62	0/113
3	j	0.46	0/84	0.53	0/113
3	k	0.51	0/84	0.59	0/113
3	l	0.42	0/84	0.59	0/113
3	m	0.46	0/84	0.69	0/113
3	n	0.45	0/84	0.68	0/113
3	o	0.45	0/84	0.48	0/113
3	p	0.48	0/84	0.57	0/113
3	q	0.43	0/84	0.54	0/113
All	All	0.49	4/45696 (0.0%)	0.61	0/61936

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	T	49	VAL	CB-CG1	-5.63	1.41	1.52
1	C	28	VAL	CB-CG1	-5.37	1.41	1.52
1	G	90	ALA	CA-CB	-5.34	1.41	1.52
1	W	28	VAL	CA-CB	-5.01	1.44	1.54

There are no bond angle outliers.

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	2254	0	2103	66	0
1	C	2254	0	2103	41	0
1	E	2254	0	2103	43	0
1	G	2254	0	2103	25	0
1	I	2254	0	2103	28	0
1	K	2254	0	2103	26	0
1	M	2254	0	2103	37	0
1	O	2254	0	2103	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Q	2254	0	2103	19	0
1	S	2254	0	2103	32	0
1	U	2254	0	2103	35	0
1	W	2254	0	2103	31	0
1	Y	2254	0	2103	46	0
1	a	2254	0	2103	0	0
2	B	837	0	803	22	0
2	D	837	0	803	10	0
2	F	837	0	803	7	0
2	H	837	0	803	7	0
2	J	837	0	803	14	0
2	L	837	0	803	4	0
2	N	837	0	803	11	0
2	P	837	0	803	13	0
2	R	837	0	803	13	0
2	T	837	0	803	15	0
2	V	837	0	803	19	0
2	X	837	0	803	10	0
2	Z	837	0	803	5	0
2	b	837	0	803	0	0
3	c	82	0	87	0	0
3	e	82	0	87	0	0
3	f	82	0	87	0	0
3	g	82	0	87	0	0
3	h	82	0	87	0	0
3	i	82	0	87	0	0
3	j	82	0	87	0	0
3	k	82	0	87	0	0
3	l	82	0	87	0	0
3	m	82	0	87	0	0
3	n	82	0	87	0	0
3	o	82	0	87	0	0
3	p	82	0	87	0	0
3	q	82	0	87	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	3	0	0	0	0
4	J	1	0	0	0	0
4	K	1	0	0	0	0
4	L	1	0	0	0	0
4	N	2	0	0	0	0
4	O	1	0	0	0	0
4	P	1	0	0	0	0
4	Q	2	0	0	0	0
4	R	1	0	0	0	0
4	S	1	0	0	0	0
4	T	1	0	0	0	0
4	U	1	0	0	0	0
4	V	1	0	0	0	0
4	W	1	0	0	0	0
4	Y	1	0	0	0	0
4	Z	1	0	0	0	0
4	a	1	0	0	0	0
4	b	1	0	0	0	0
5	A	6	0	8	0	0
5	C	12	0	16	1	0
5	E	6	0	8	0	0
5	K	6	0	8	0	0
5	O	12	0	16	1	0
5	Q	6	0	8	0	0
5	S	6	0	8	0	0
5	Y	6	0	8	0	0
5	a	6	0	8	0	0
5	c	6	0	8	0	0
5	e	6	0	8	0	0
5	f	6	0	8	0	0
5	g	6	0	8	0	0
5	h	6	0	8	0	0
5	i	6	0	8	0	0
5	l	6	0	8	0	0
5	o	6	0	8	0	0
5	q	6	0	8	0	0
6	A	2	0	0	0	0
6	C	4	0	0	0	0
6	D	2	0	0	0	0
6	E	3	0	0	1	0
6	F	2	0	0	0	0
6	G	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	H	1	0	0	0	0
6	I	3	0	0	0	0
6	J	2	0	0	1	0
6	K	3	0	0	0	0
6	L	3	0	0	0	0
6	M	1	0	0	0	0
6	P	3	0	0	0	0
6	Q	2	0	0	0	0
6	R	3	0	0	0	0
6	S	1	0	0	0	0
6	T	1	0	0	0	0
6	U	3	0	0	0	0
6	V	2	0	0	0	0
6	X	3	0	0	0	0
6	Y	2	0	0	0	0
6	Z	2	0	0	0	0
All	All	44627	0	42062	528	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 7.

All (528) 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)
1:K:126:LEU:HD22	1:K:156:LEU:HD13	1.47	0.97
1:A:90:ALA:HB1	2:N:85:VAL:HG22	1.45	0.97
1:A:231:VAL:HG11	1:A:244:TRP:CE2	2.05	0.91
1:U:126:LEU:HD22	1:U:156:LEU:HD13	1.50	0.90
1:C:90:ALA:HB3	2:V:2:GLN:HE22	1.35	0.88
2:L:39:LEU:HD23	2:L:68:THR:HG22	1.55	0.88
1:M:16:GLY:O	1:M:17:ARG:HG2	1.78	0.83
1:C:90:ALA:HB1	2:V:85:VAL:HG13	1.61	0.83
1:E:90:ALA:HB1	2:J:85:VAL:HG22	1.61	0.83
2:X:7:ILE:CD1	2:X:82:VAL:HG21	2.10	0.81
1:U:215:LEU:HD22	1:U:261:VAL:HG22	1.62	0.80
1:Y:126:LEU:HD22	1:Y:156:LEU:HD13	1.63	0.80
1:C:90:ALA:CB	2:V:2:GLN:HE22	1.96	0.79
1:W:28:VAL:HG11	1:W:179:LEU:HD13	1.63	0.79
1:A:156:LEU:CD2	1:A:160:LEU:HD11	4.04	0.78
1:E:90:ALA:HB3	2:J:2:GLN:OE1	1.83	0.78
1:C:90:ALA:HB3	2:V:2:GLN:NE2	1.99	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:82:ARG:HD3	1:I:89:GLU:HA	1.66	0.77
1:U:126:LEU:CD2	1:U:156:LEU:HD13	2.14	0.77
1:S:65:ARG:HD3	1:W:190:THR:HG22	1.67	0.76
1:U:215:LEU:CD2	1:U:261:VAL:HG22	2.15	0.76
1:W:51:TRP:CZ2	1:W:179:LEU:HD11	2.21	0.76
2:T:47:GLU:O	2:T:47:GLU:HG3	1.84	0.76
1:C:230:LEU:HD23	1:C:245:ALA:HB2	1.67	0.75
2:B:11:SER:OG	2:B:13:HIS:O	2.56	0.75
1:G:215:LEU:CD2	1:G:261:VAL:HG22	2.16	0.75
1:G:224:GLN:O	1:G:228:THR:HG23	1.86	0.74
1:W:135:ALA:HB1	1:W:140:ALA:HB3	1.69	0.73
1:W:201:LEU:HD22	1:W:274:TRP:CZ2	2.22	0.73
2:X:7:ILE:HD13	2:X:82:VAL:HG21	1.71	0.72
1:M:255:GLN:N	1:M:255:GLN:OE1	2.21	0.72
1:C:90:ALA:HB1	2:V:85:VAL:CG1	2.20	0.71
1:K:98:MET:CE	1:K:115:GLN:OE1	2.39	0.71
1:A:150:ALA:HB1	1:M:252:GLY:HA3	61.73	0.70
1:A:82:ARG:HD2	1:A:89:GLU:HA	1.71	0.70
2:B:2:GLN:OE1	1:M:90:ALA:HB3	1.91	0.70
2:T:39:LEU:HD22	2:T:49:VAL:CG1	2.21	0.69
1:S:218:GLN:OE1	1:S:260:HIS:NE2	2.24	0.69
2:B:54:LEU:HA	2:B:64:LEU:HD21	2.63	0.69
1:M:231:VAL:CG1	1:M:244:TRP:CZ2	2.76	0.69
1:Q:28:VAL:HG11	1:Q:179:LEU:HD13	1.76	0.68
1:C:90:ALA:CB	2:V:2:GLN:NE2	2.55	0.68
1:E:126:LEU:HD22	1:E:156:LEU:HD13	1.74	0.68
1:I:135:ALA:HB1	1:I:140:ALA:HB3	1.76	0.68
1:K:98:MET:HE2	1:K:115:GLN:OE1	1.95	0.67
1:Y:231:VAL:HG13	1:Y:244:TRP:CZ2	2.30	0.67
1:A:231:VAL:HG11	1:A:244:TRP:CZ2	2.30	0.66
2:D:24:ASN:HB3	2:D:65:LEU:HD11	1.77	0.66
1:C:201:LEU:HD22	1:C:274:TRP:CZ2	2.31	0.66
1:E:228:THR:HG22	1:E:247:VAL:HG23	1.78	0.66
1:K:64:THR:O	1:K:67:VAL:HG12	1.96	0.65
1:M:156:LEU:HD22	1:M:160:LEU:CD1	2.25	0.65
1:C:65:ARG:HD3	1:K:190:THR:HG22	1.79	0.65
1:Y:156:LEU:HD22	1:Y:160:LEU:HD11	1.79	0.65
1:E:90:ALA:CB	2:J:85:VAL:HG22	2.27	0.65
1:Y:231:VAL:CG1	1:Y:244:TRP:CZ2	2.80	0.65
1:A:32:GLN:HE21	1:A:48:ARG:HG3	2.88	0.64
1:I:12:VAL:HG22	1:I:94:THR:HG23	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:40:ALA:O	1:O:43:GLN:NE2	2.30	0.64
2:X:2:GLN:OE1	1:Y:90:ALA:HB3	1.97	0.64
1:Q:117:ALA:HB2	2:R:60:TRP:CE2	2.33	0.64
1:M:231:VAL:HG13	1:M:244:TRP:CZ2	2.33	0.64
2:T:47:GLU:O	2:T:47:GLU:CG	2.45	0.64
1:E:44:ARG:HA	1:E:64:THR:HG23	1.80	0.64
1:Q:266:LEU:HD13	1:Q:270:LEU:HG	1.80	0.63
1:S:163:THR:HG22	1:S:167:TRP:CD1	2.33	0.63
1:M:156:LEU:HD22	1:M:160:LEU:HD11	1.81	0.63
2:V:24:ASN:HB3	2:V:65:LEU:HD11	1.80	0.63
1:Y:126:LEU:HD13	1:Y:133:TRP:CH2	2.34	0.63
2:P:2:GLN:HB3	2:P:86:THR:HG22	1.80	0.62
2:X:45:ARG:NH1	2:X:47:GLU:OE1	2.30	0.62
1:W:28:VAL:O	1:W:28:VAL:HG12	1.97	0.62
1:C:230:LEU:CD2	1:C:245:ALA:HB2	2.30	0.62
1:O:224:GLN:O	1:O:228:THR:HG23	1.98	0.62
1:A:231:VAL:CG1	1:A:244:TRP:CE2	2.82	0.62
1:I:2:SER:OG	1:I:110:LEU:HD12	1.98	0.62
1:Y:270:LEU:HD22	1:Y:272:LEU:CD2	2.30	0.62
2:T:39:LEU:HD22	2:T:49:VAL:HG11	1.81	0.62
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.35	0.61
1:Y:219:ARG:HG3	1:Y:257:TYR:CZ	2.35	0.61
1:W:49:ALA:O	1:W:52:ILE:HG22	2.00	0.61
1:A:33:PHE:CD1	1:A:34:VAL:HG13	2.35	0.61
1:M:126:LEU:HD22	1:M:156:LEU:HD13	1.82	0.61
2:X:7:ILE:HD11	2:X:82:VAL:HG21	1.80	0.61
2:V:84:HIS:N	2:V:87:LEU:HD12	2.16	0.60
2:V:54:LEU:HA	2:V:64:LEU:HD21	1.83	0.60
2:B:46:ILE:O	2:B:49:VAL:HG23	2.42	0.60
1:C:126:LEU:HD23	1:C:130:LEU:HD22	1.83	0.60
2:D:2:GLN:HE22	1:U:88:SER:CB	2.13	0.60
1:Y:49:ALA:O	1:Y:52:ILE:HG22	2.01	0.60
1:A:90:ALA:CB	2:N:85:VAL:HG22	2.27	0.60
1:S:103:VAL:CG1	1:S:165:VAL:HG13	2.32	0.60
1:W:135:ALA:HB1	1:W:140:ALA:CB	2.31	0.60
2:N:25:CYS:HB2	2:N:39:LEU:HD11	1.83	0.59
1:W:231:VAL:HG11	1:W:244:TRP:CZ2	2.37	0.59
2:H:9:VAL:HG22	1:K:143:THR:HG21	70.92	0.59
1:K:116:TYR:HB3	1:K:124:ILE:HG22	1.83	0.59
1:E:231:VAL:CG1	1:E:244:TRP:CZ2	2.86	0.59
1:M:231:VAL:HG11	1:M:244:TRP:CZ2	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:U:126:LEU:HD22	1:U:156:LEU:CD1	2.28	0.59
1:G:88:SER:C	1:G:90:ALA:H	2.07	0.59
1:A:32:GLN:NE2	1:A:48:ARG:HG3	2.89	0.58
1:C:117:ALA:HB2	2:D:60:TRP:CE2	2.38	0.58
1:E:15:PRO:HB2	2:J:85:VAL:HG23	1.85	0.58
2:N:9:VAL:HG11	1:O:81:LEU:HD21	71.05	0.58
2:T:39:LEU:HD22	2:T:49:VAL:HG13	1.85	0.58
1:U:82:ARG:HD2	1:U:89:GLU:HA	1.86	0.58
1:A:231:VAL:CG1	1:A:244:TRP:CZ2	2.87	0.57
1:A:143:THR:HG21	2:J:9:VAL:HG23	43.06	0.57
1:A:252:GLY:HA3	1:Y:150:ALA:HB1	153.72	0.57
1:Y:270:LEU:HD22	1:Y:272:LEU:HD23	1.86	0.57
1:A:49:ALA:O	1:A:52:ILE:HG22	2.14	0.57
1:E:37:ASP:HB3	1:E:40:ALA:HB2	1.86	0.57
1:A:252:GLY:O	1:A:255:GLN:NE2	2.38	0.56
1:C:165:VAL:HG12	1:C:169:ARG:NH1	2.19	0.56
1:C:230:LEU:HD23	1:C:245:ALA:CB	2.35	0.56
1:C:231:VAL:CG1	1:C:244:TRP:CZ2	2.87	0.56
1:E:138:MET:N	1:M:128:GLU:OE2	2.35	0.56
1:A:150:ALA:CB	1:M:252:GLY:HA3	61.85	0.56
1:Y:156:LEU:HD22	1:Y:160:LEU:CD1	2.35	0.56
1:A:215:LEU:HD22	1:A:261:VAL:HG22	1.86	0.56
1:W:201:LEU:HD13	1:W:274:TRP:NE1	2.20	0.56
1:Q:13:SER:HB3	1:Q:78:LEU:HD13	1.86	0.56
1:K:231:VAL:CG1	1:K:244:TRP:CZ2	2.88	0.56
1:Y:8:PHE:HB2	1:Y:25:VAL:HG23	1.88	0.56
1:M:172:LEU:O	1:M:180:GLN:NE2	2.35	0.56
1:U:28:VAL:HG11	1:U:179:LEU:HD13	1.87	0.56
1:A:13:SER:HB3	1:A:78:LEU:HD13	1.88	0.56
1:C:152:VAL:HG13	5:C:302:GOL:H11	1.88	0.56
1:S:12:VAL:HG22	1:S:94:THR:CG2	2.36	0.56
2:T:37:VAL:HG13	2:T:82:VAL:HG22	1.87	0.56
1:G:82:ARG:HD2	1:G:89:GLU:HA	1.87	0.55
2:P:24:ASN:HB3	2:P:65:LEU:HD11	1.86	0.55
1:S:12:VAL:HG22	1:S:94:THR:HG23	1.87	0.55
1:M:11:SER:HG	1:M:22:PHE:HD1	1.53	0.55
1:Q:231:VAL:HG13	1:Q:244:TRP:CZ2	2.41	0.55
1:G:88:SER:O	1:G:90:ALA:N	2.40	0.55
1:A:252:GLY:O	1:U:150:ALA:HB1	2.06	0.55
1:S:224:GLN:O	1:S:228:THR:HG23	2.07	0.55
1:C:231:VAL:HG13	1:C:244:TRP:CZ2	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:ALA:HB1	2:N:85:VAL:CG2	2.31	0.54
1:G:215:LEU:HD23	1:G:261:VAL:HG22	1.87	0.54
1:M:28:VAL:HG23	1:M:33:PHE:CD2	2.42	0.54
1:Q:117:ALA:HB2	2:R:60:TRP:CZ2	2.42	0.54
1:Y:72:GLN:O	1:Y:76:VAL:HG23	2.07	0.54
1:A:116:TYR:HB3	1:A:124:ILE:HG22	2.10	0.54
1:C:126:LEU:CD2	1:C:130:LEU:HD22	2.38	0.54
2:P:39:LEU:HD23	2:P:49:VAL:CG2	2.36	0.54
1:I:14:ARG:HB3	1:I:17:ARG:HB3	1.90	0.54
1:O:231:VAL:HG13	1:O:244:TRP:CZ2	2.43	0.54
1:W:231:VAL:CG1	1:W:244:TRP:CZ2	2.91	0.54
1:A:90:ALA:HB3	2:N:2:GLN:OE1	2.08	0.54
1:A:109:PHE:HB2	1:A:165:VAL:HG21	1.89	0.54
1:I:168:LEU:HD12	1:I:168:LEU:O	2.08	0.54
1:I:110:LEU:HD13	1:I:111:ARG:NH2	2.23	0.53
1:K:11:SER:HA	1:K:21:ARG:O	2.07	0.53
2:D:84:HIS:CE1	2:D:86:THR:HG23	2.42	0.53
2:D:85:VAL:HG22	1:U:90:ALA:HA	1.90	0.53
1:G:226:GLN:CD	1:G:226:GLN:H	2.12	0.53
1:G:88:SER:C	1:G:90:ALA:N	2.60	0.53
1:I:258:THR:OG1	1:I:260:HIS:NE2	2.42	0.53
2:R:7:ILE:CD1	2:R:82:VAL:HG21	2.38	0.53
1:U:110:LEU:HD23	1:U:110:LEU:O	2.09	0.53
1:E:231:VAL:HG13	1:E:244:TRP:CZ2	2.43	0.53
1:G:35:ARG:NH2	1:G:40:ALA:HB2	2.24	0.53
1:I:165:VAL:HG12	1:I:169:ARG:CZ	2.39	0.53
1:M:117:ALA:HB2	2:N:60:TRP:CE2	2.44	0.53
1:C:104:GLY:CA	1:C:110:LEU:HD12	2.39	0.53
1:C:90:ALA:CB	2:V:85:VAL:HG13	2.36	0.53
1:M:16:GLY:C	1:M:17:ARG:CG	2.77	0.53
1:Q:156:LEU:HD22	1:Q:160:LEU:HG	1.91	0.53
1:O:45:MET:HE1	1:O:67:VAL:HB	1.90	0.53
1:I:49:ALA:O	1:I:52:ILE:HG22	2.09	0.52
1:C:250:PRO:HB2	1:C:253:GLN:HE21	1.75	0.52
1:K:130:LEU:HB3	1:K:157:ARG:HG3	1.92	0.52
1:E:33:PHE:CD1	1:E:34:VAL:HG13	2.43	0.52
1:A:84:TYR:HB3	1:A:139:ALA:HB1	2.80	0.52
1:I:187:THR:HB	1:I:272:LEU:HD11	1.90	0.52
2:F:9:VAL:HG13	1:G:143:THR:OG1	109.32	0.52
1:G:219:ARG:HB3	1:G:224:GLN:HE21	1.75	0.52
1:S:231:VAL:HG13	1:S:244:TRP:CE2	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:117:ALA:HB2	2:Z:60:TRP:CE2	2.45	0.52
1:S:231:VAL:CG1	1:S:244:TRP:CE2	2.93	0.52
1:Y:33:PHE:CD1	1:Y:34:VAL:HG13	2.45	0.52
2:B:85:VAL:HG23	1:M:15:PRO:HB2	1.92	0.52
1:U:169:ARG:HA	1:U:172:LEU:HD12	1.92	0.52
1:A:255:GLN:CD	1:A:255:GLN:H	2.14	0.51
2:X:1:ILE:HD12	2:X:1:ILE:C	2.31	0.51
1:A:159:TYR:HB2	1:S:276:PRO:HB2	1.92	0.51
1:A:72:GLN:HE22	1:A:75:ARG:HH21	1.58	0.51
1:G:215:LEU:HD22	1:G:261:VAL:HG22	1.88	0.51
1:K:81:LEU:HD13	1:K:118:TYR:CD1	2.45	0.51
1:O:152:VAL:HG13	5:O:302:GOL:H11	1.93	0.51
1:O:219:ARG:O	1:O:222:GLU:HG3	2.10	0.51
1:Q:253:GLN:HA	1:Q:255:GLN:NE2	2.26	0.51
1:A:11:SER:HA	1:A:21:ARG:O	2.10	0.51
1:A:193:ALA:O	1:U:72:GLN:NE2	2.40	0.51
2:T:23:LEU:HB2	2:T:70:PHE:CE1	2.46	0.51
1:C:33:PHE:CD1	1:C:34:VAL:HG13	2.45	0.51
1:K:231:VAL:HG11	1:K:244:TRP:CZ2	2.46	0.51
1:U:25:VAL:HG21	2:V:55:SER:OG	2.11	0.51
1:G:13:SER:HB3	1:G:78:LEU:HD13	1.92	0.51
1:E:261:VAL:HG12	1:E:266:LEU:HD12	1.93	0.50
2:P:2:GLN:HB3	2:P:86:THR:CG2	2.40	0.50
2:P:39:LEU:HD23	2:P:49:VAL:HG21	1.93	0.50
1:G:258:THR:HG22	1:G:273:ARG:HA	1.94	0.50
1:M:16:GLY:C	1:M:17:ARG:HG2	2.30	0.50
1:U:117:ALA:HB2	2:V:60:TRP:CE2	2.46	0.50
1:A:230:LEU:HD23	1:A:245:ALA:HB2	2.20	0.50
1:C:201:LEU:HD13	1:C:274:TRP:NE1	2.26	0.50
1:S:117:ALA:HB2	2:T:60:TRP:CE2	2.46	0.50
1:O:122:ASP:OD1	2:P:60:TRP:NE1	2.39	0.50
1:O:12:VAL:HG22	1:O:94:THR:HG23	1.94	0.50
1:K:126:LEU:HD22	1:K:156:LEU:CD1	2.30	0.50
1:A:142:THR:OG1	1:I:111:ARG:NH1	2.45	0.50
1:M:135:ALA:HB1	1:M:140:ALA:CB	2.42	0.50
1:O:172:LEU:HD23	1:O:179:LEU:HD13	1.94	0.50
1:S:231:VAL:HG13	1:S:244:TRP:CZ2	2.47	0.50
1:E:156:LEU:HD22	1:E:160:LEU:HD11	1.93	0.49
1:S:150:ALA:HB1	1:W:252:GLY:HA3	1.94	0.49
1:S:103:VAL:HG11	1:S:165:VAL:HG13	1.93	0.49
1:A:156:LEU:HD22	1:A:160:LEU:HD11	3.10	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:261:VAL:HG11	1:E:266:LEU:HD11	1.94	0.49
1:O:126:LEU:HD22	1:O:156:LEU:HD13	1.93	0.49
1:U:110:LEU:C	1:U:110:LEU:HD23	2.32	0.49
1:W:32:GLN:NE2	1:W:35:ARG:HB3	2.27	0.49
1:A:100:GLY:O	1:A:160:LEU:HD22	2.37	0.49
2:V:25:CYS:HB2	2:V:39:LEU:HD11	1.94	0.49
1:A:214:THR:HG21	1:A:262:GLN:HE21	5.41	0.49
2:B:37:VAL:HB	2:B:66:TYR:CZ	2.55	0.49
1:U:117:ALA:HB2	2:V:60:TRP:CZ2	2.47	0.49
1:W:249:VAL:HG12	1:W:257:TYR:CZ	2.48	0.49
1:I:244:TRP:CZ3	1:I:246:ALA:HB2	2.48	0.49
2:T:24:ASN:HB3	2:T:65:LEU:HD11	1.94	0.49
1:A:156:LEU:CD2	1:A:160:LEU:CD1	4.68	0.49
1:E:89:GLU:HG3	1:E:89:GLU:O	2.13	0.49
1:M:231:VAL:HG11	1:M:244:TRP:CE2	2.48	0.49
1:U:208:PHE:CZ	1:U:213:ILE:HG21	2.48	0.49
1:C:90:ALA:HB3	2:V:2:GLN:OE1	2.13	0.48
1:E:201:LEU:HD22	1:E:274:TRP:CZ2	2.48	0.48
1:I:150:ALA:HB3	1:I:152:VAL:HG23	1.95	0.48
1:S:203:CYS:O	1:S:244:TRP:HA	2.13	0.48
2:J:39:LEU:HD23	2:J:68:THR:HG22	1.95	0.48
2:J:4:THR:HG22	6:J:202:HOH:O	2.13	0.48
1:O:117:ALA:HB2	2:P:60:TRP:CE2	2.48	0.48
1:A:258:THR:HG22	1:A:273:ARG:HA	2.38	0.48
1:A:8:PHE:HB2	1:A:25:VAL:CG2	2.43	0.48
1:A:84:TYR:HB3	1:A:139:ALA:CB	3.20	0.48
1:M:45:MET:HE2	1:M:67:VAL:HB	1.93	0.48
1:M:79:GLY:O	1:M:82:ARG:HB3	2.13	0.48
1:Y:126:LEU:HD13	1:Y:133:TRP:CZ3	2.48	0.48
1:C:90:ALA:HB3	2:V:2:GLN:CD	2.34	0.48
1:A:117:ALA:HB2	2:B:60:TRP:CZ2	2.51	0.48
1:C:231:VAL:HG11	1:C:244:TRP:CZ2	2.48	0.48
1:A:16:GLY:O	2:N:83:ASN:ND2	2.47	0.48
2:D:85:VAL:HG22	1:U:90:ALA:CA	2.44	0.48
1:W:209:TYR:CD1	1:W:210:PRO:HA	2.49	0.47
1:O:78:LEU:CD2	1:O:95:VAL:HG23	2.44	0.47
1:U:21:ARG:NH2	1:U:37:ASP:OD2	2.43	0.47
1:A:88:SER:C	1:A:90:ALA:H	2.18	0.47
1:C:138:MET:HG2	1:G:128:GLU:O	2.15	0.47
1:A:64:THR:O	1:A:67:VAL:HG12	2.22	0.47
1:E:261:VAL:HG12	1:E:266:LEU:CD1	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:51:TRP:O	1:K:54:GLN:HG2	2.14	0.47
1:M:45:MET:CE	1:M:67:VAL:HB	2.44	0.47
1:C:165:VAL:HG12	1:C:169:ARG:CZ	2.45	0.47
1:E:66:LYS:O	1:E:69:ALA:HB3	2.14	0.47
1:G:231:VAL:HG11	1:G:244:TRP:CZ2	2.50	0.47
1:S:231:VAL:CG1	1:S:244:TRP:CZ2	2.97	0.47
1:Y:185:PRO:HA	1:Y:208:PHE:HB3	1.97	0.47
1:K:133:TRP:NE1	1:K:153:ALA:HB2	2.29	0.47
1:Y:231:VAL:HG13	1:Y:244:TRP:CE2	2.49	0.47
1:A:218:GLN:O	1:A:257:TYR:HA	2.53	0.47
1:M:224:GLN:O	1:M:228:THR:HG23	2.13	0.47
1:U:51:TRP:CZ2	1:U:179:LEU:HD11	2.49	0.47
2:X:5:PRO:HB3	2:X:30:PHE:HB3	1.96	0.47
1:Y:153:ALA:HB3	1:Y:154:GLU:OE1	2.15	0.47
1:I:165:VAL:HG12	1:I:169:ARG:NH1	2.30	0.47
1:C:90:ALA:CB	2:V:2:GLN:OE1	2.62	0.47
2:Z:98:ASP:OD1	2:Z:98:ASP:N	2.48	0.47
1:C:252:GLY:HA3	1:M:150:ALA:HB1	1.97	0.47
1:I:47:PRO:HB3	1:I:52:ILE:HG23	1.97	0.47
1:S:253:GLN:OE1	1:S:256:ARG:NH1	2.48	0.47
2:X:2:GLN:OE1	1:Y:90:ALA:CB	2.62	0.47
1:K:82:ARG:HG3	1:K:87:GLN:HB2	1.97	0.47
1:Y:109:PHE:HD2	1:Y:165:VAL:HG21	1.80	0.47
1:Y:184:ALA:HB2	1:Y:265:GLY:O	2.15	0.47
1:Y:234:ARG:NH2	2:Z:99:MET:OXT	2.48	0.47
1:M:36:PHE:CD1	1:M:67:VAL:HG11	2.50	0.46
1:S:163:THR:HG22	1:S:167:TRP:HD1	1.78	0.46
1:A:112:GLY:O	1:A:130:LEU:HD21	2.48	0.46
1:S:126:LEU:HD12	1:S:132:SER:O	2.16	0.46
1:S:187:THR:HB	1:S:272:LEU:HD11	1.96	0.46
2:X:35:ILE:O	2:X:35:ILE:HG23	2.15	0.46
1:A:201:LEU:HD22	1:A:274:TRP:CZ2	2.51	0.46
1:C:100:GLY:O	1:C:160:LEU:HD22	2.15	0.46
1:C:218:GLN:OE1	1:C:260:HIS:NE2	2.48	0.46
1:E:187:THR:HB	1:E:272:LEU:HD11	1.98	0.46
1:I:33:PHE:CD1	1:I:34:VAL:HG13	2.51	0.46
1:Y:231:VAL:CG1	1:Y:244:TRP:CE2	2.98	0.46
1:Y:81:LEU:HD13	1:Y:118:TYR:CD1	2.51	0.46
1:G:103:VAL:HG13	1:G:168:LEU:HD23	1.97	0.46
1:C:28:VAL:O	1:C:29:ASP:HB2	2.16	0.46
1:M:51:TRP:CZ2	1:M:179:LEU:HD11	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:THR:HG22	1:A:32:GLN:O	2.16	0.46
2:D:85:VAL:HG13	1:U:90:ALA:HB1	1.98	0.46
1:O:215:LEU:HD22	1:O:261:VAL:HG22	1.98	0.46
1:S:12:VAL:HG13	1:S:94:THR:HG23	1.98	0.46
2:T:27:VAL:HG23	2:T:30:PHE:CE1	2.50	0.46
1:W:230:LEU:HD23	1:W:245:ALA:HB2	1.97	0.46
1:W:13:SER:HB3	1:W:78:LEU:HD22	1.98	0.45
1:Y:7:TYR:HB3	1:Y:9:PHE:CE2	2.51	0.45
1:U:182:THR:HG21	1:U:264:GLU:HG2	1.97	0.45
1:K:165:VAL:HG12	1:K:169:ARG:CZ	2.46	0.45
1:E:231:VAL:HG11	1:E:244:TRP:CE2	2.51	0.45
1:I:172:LEU:HD23	1:I:179:LEU:HD13	1.97	0.45
2:N:49:VAL:HG22	2:N:68:THR:HB	1.99	0.45
2:X:54:LEU:HD11	2:X:62:PHE:HB3	1.98	0.45
1:Y:111:ARG:HG2	1:Y:112:GLY:N	2.31	0.45
1:A:143:THR:HG21	2:J:9:VAL:CG2	43.61	0.45
1:E:44:ARG:HD2	2:P:16:GLU:HG3	1.98	0.45
2:R:32:PRO:HD2	2:R:85:VAL:HG11	1.99	0.45
1:S:33:PHE:O	1:S:52:ILE:HG21	2.17	0.45
1:G:8:PHE:HB2	1:G:25:VAL:HG23	1.99	0.45
2:T:39:LEU:HD23	2:T:68:THR:HG22	1.99	0.45
1:E:231:VAL:CG1	1:E:244:TRP:CE2	3.00	0.45
1:O:231:VAL:CG1	1:O:244:TRP:CZ2	3.00	0.45
2:R:59:ASP:O	2:R:60:TRP:HB2	2.17	0.45
1:U:33:PHE:CD1	1:U:34:VAL:HG13	2.52	0.45
1:A:137:ASP:H	1:A:140:ALA:HB3	1.89	0.45
1:I:124:ILE:O	1:I:124:ILE:HG23	2.15	0.45
1:I:66:LYS:O	1:I:69:ALA:HB3	2.17	0.45
2:B:54:LEU:CA	2:B:64:LEU:HD21	3.22	0.45
2:L:25:CYS:HB2	2:L:39:LEU:HD11	1.99	0.45
2:R:5:PRO:HB3	2:R:30:PHE:HB3	1.99	0.45
1:Y:165:VAL:HG12	1:Y:169:ARG:NH1	2.32	0.45
2:H:5:PRO:HB3	2:H:30:PHE:HB3	1.98	0.45
1:E:40:ALA:HB3	6:E:401:HOH:O	2.16	0.44
1:M:82:ARG:HD3	1:M:89:GLU:HA	1.99	0.44
1:O:49:ALA:O	1:O:52:ILE:HG22	2.17	0.44
1:A:194:VAL:HG22	1:A:198:GLU:O	2.17	0.44
2:H:40:LEU:HD23	2:H:45:ARG:HA	1.99	0.44
1:A:77:ASP:CG	2:J:9:VAL:HG12	44.22	0.44
1:Y:98:MET:C	1:Y:98:MET:SD	2.96	0.44
1:A:219:ARG:O	1:A:219:ARG:HG3	4.45	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:23:LEU:O	2:B:67:TYR:HA	2.17	0.44
1:E:231:VAL:HG11	1:E:244:TRP:CZ2	2.52	0.44
2:F:54:LEU:HD12	2:F:64:LEU:HG	1.99	0.44
1:M:103:VAL:HB	1:M:107:TRP:HA	1.99	0.44
1:Y:152:VAL:O	1:Y:155:GLN:N	2.50	0.44
1:E:90:ALA:CA	2:J:85:VAL:HG22	2.48	0.44
2:H:34:ASP:O	2:H:84:HIS:HD2	2.00	0.44
1:O:45:MET:CE	1:O:67:VAL:HB	2.46	0.44
1:Q:32:GLN:NE2	1:Q:35:ARG:HB3	2.33	0.44
1:Y:74:HIS:CE1	1:Y:97:ARG:HE	2.35	0.44
1:U:13:SER:HB3	1:U:78:LEU:HD13	2.00	0.44
1:A:81:LEU:HD13	1:A:118:TYR:CD1	2.53	0.44
1:E:28:VAL:HG23	1:E:33:PHE:CD2	2.52	0.44
1:G:12:VAL:HG22	1:G:94:THR:HG23	1.99	0.44
1:I:138:MET:HB3	1:I:138:MET:HE3	1.86	0.44
1:Q:21:ARG:HD3	1:Q:39:ASP:OD2	2.18	0.44
1:A:201:LEU:HD13	1:A:274:TRP:NE1	2.42	0.44
1:E:261:VAL:CG1	1:E:266:LEU:CD1	2.96	0.44
1:G:103:VAL:HG12	1:G:165:VAL:CG1	2.48	0.44
1:S:196:ASP:OD2	1:S:197:HIS:N	2.50	0.44
1:K:214:THR:HG22	1:K:216:THR:HG23	2.00	0.43
1:S:109:PHE:HZ	1:S:130:LEU:HD11	1.83	0.43
1:Y:95:VAL:HG13	1:Y:116:TYR:CE2	2.52	0.43
1:A:255:GLN:HA	1:A:274:TRP:HB2	2.30	0.43
2:B:39:LEU:O	2:B:46:ILE:HD12	3.44	0.43
1:E:3:HIS:HA	1:E:29:ASP:OD1	2.18	0.43
1:G:187:THR:HB	1:G:272:LEU:HD11	2.00	0.43
1:K:98:MET:HE1	1:K:115:GLN:OE1	2.16	0.43
1:Q:28:VAL:CG1	1:Q:179:LEU:HD13	2.46	0.43
1:Q:231:VAL:CG1	1:Q:244:TRP:CZ2	3.01	0.43
2:R:39:LEU:HD23	2:R:68:THR:HG22	2.00	0.43
1:Y:51:TRP:CE2	1:Y:179:LEU:HD11	2.53	0.43
1:A:252:GLY:HA3	1:U:150:ALA:CB	2.48	0.43
1:E:82:ARG:HD3	1:E:89:GLU:HA	1.99	0.43
1:U:226:GLN:HE21	1:U:226:GLN:HA	1.83	0.43
1:K:8:PHE:HB2	1:K:25:VAL:CG2	2.48	0.43
1:M:231:VAL:CG1	1:M:244:TRP:CE2	3.02	0.43
1:S:85:TYR:OH	1:S:137:ASP:OD2	2.35	0.43
2:T:2:GLN:HB3	2:T:86:THR:HG22	1.99	0.43
1:W:81:LEU:HD13	1:W:118:TYR:CD1	2.54	0.43
1:Y:219:ARG:HG3	1:Y:257:TYR:CE2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:40:LEU:HD23	2:R:45:ARG:HA	2.00	0.43
1:Y:25:VAL:HG12	1:Y:32:GLN:NE2	2.33	0.43
1:A:103:VAL:HG13	1:A:168:LEU:HD23	2.00	0.43
1:C:8:PHE:HB2	1:C:25:VAL:HG22	2.00	0.43
1:E:138:MET:SD	1:M:110:LEU:HD22	2.59	0.43
1:E:247:VAL:HG13	1:E:249:VAL:HG13	2.00	0.43
2:P:5:PRO:HB3	2:P:30:PHE:HB3	2.00	0.43
1:S:103:VAL:HG12	1:S:165:VAL:HG13	2.01	0.43
2:B:5:PRO:HB3	2:B:30:PHE:HB3	2.00	0.43
2:R:7:ILE:HD13	2:R:82:VAL:HG21	1.99	0.43
1:A:215:LEU:CD2	1:A:261:VAL:HG22	2.48	0.43
2:F:51:HIS:HB3	2:F:66:TYR:CD2	2.54	0.43
1:U:82:ARG:CD	1:U:89:GLU:HA	2.49	0.43
2:V:68:THR:HG23	2:V:69:GLU:O	2.19	0.43
1:A:219:ARG:O	1:A:220:ASP:C	2.93	0.43
1:I:191:HIS:HB2	1:I:274:TRP:CH2	2.54	0.43
1:I:244:TRP:HZ3	1:I:246:ALA:HB2	1.83	0.43
1:Q:12:VAL:O	1:Q:21:ARG:N	2.52	0.43
1:U:252:GLY:O	1:U:255:GLN:NE2	2.51	0.43
1:E:43:GLN:HE21	2:P:74:GLU:HG2	1.84	0.43
1:W:203:CYS:O	1:W:244:TRP:HB2	2.19	0.43
1:A:7:TYR:HB3	1:A:9:PHE:CE2	2.55	0.42
1:A:88:SER:O	1:A:90:ALA:N	2.52	0.42
2:B:73:THR:O	2:B:97:ARG:NH2	2.97	0.42
1:G:117:ALA:HB2	2:H:60:TRP:CE2	2.54	0.42
2:P:2:GLN:HG3	2:P:32:PRO:HD3	2.01	0.42
1:Y:270:LEU:HD22	1:Y:272:LEU:HD21	1.99	0.42
1:A:173:GLU:O	1:A:176:LYS:HB2	2.41	0.42
1:C:97:ARG:HH11	1:C:114:HIS:CE1	2.36	0.42
2:H:23:LEU:O	2:H:67:TYR:HA	2.19	0.42
1:M:109:PHE:CD2	1:M:161:GLU:HA	2.54	0.42
1:M:32:GLN:NE2	1:M:35:ARG:HB3	2.34	0.42
2:T:37:VAL:HG21	2:T:66:TYR:CE1	2.54	0.42
1:W:216:THR:HG21	1:W:223:ASP:OD2	2.19	0.42
1:C:268:LYS:HG2	1:C:269:PRO:HD2	2.00	0.42
1:E:89:GLU:CG	1:E:89:GLU:O	2.65	0.42
1:I:81:LEU:HD21	2:L:9:VAL:HG11	104.64	0.42
2:J:24:ASN:HB3	2:J:65:LEU:HD11	2.01	0.42
1:K:133:TRP:HE1	1:K:153:ALA:HB2	1.84	0.42
1:K:175:GLY:O	1:K:179:LEU:N	2.50	0.42
2:B:10:TYR:N	2:B:10:TYR:CD1	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:21:ASN:HB3	2:F:70:PHE:CE1	2.53	0.42
2:F:7:ILE:HB	2:F:93:VAL:HG21	2.01	0.42
2:H:21:ASN:OD1	2:H:22:PHE:N	2.48	0.42
1:O:78:LEU:HD23	1:O:95:VAL:HG23	2.02	0.42
1:U:109:PHE:CD2	1:U:161:GLU:HA	2.54	0.42
1:W:30:ASP:HB2	1:W:209:TYR:HE1	1.84	0.42
2:B:24:ASN:HB3	2:B:65:LEU:HD11	2.37	0.42
1:U:194:VAL:HG23	1:U:195:SER:N	2.34	0.42
2:F:98:ASP:OD1	2:F:98:ASP:N	2.53	0.42
1:U:236:ALA:HB2	1:U:242:GLN:HE21	1.84	0.42
2:Z:5:PRO:HB3	2:Z:30:PHE:HB3	2.02	0.42
1:A:15:PRO:HG2	2:N:85:VAL:HG23	2.01	0.42
1:I:165:VAL:CG1	1:I:169:ARG:CZ	2.97	0.42
1:O:104:GLY:HA2	1:O:110:LEU:HD12	2.01	0.42
2:T:23:LEU:O	2:T:67:TYR:HA	2.19	0.42
1:S:235:PRO:HG2	2:T:65:LEU:HD22	2.01	0.42
1:W:64:THR:O	1:W:67:VAL:HG12	2.19	0.42
1:C:10:THR:HG21	2:D:54:LEU:HD23	2.00	0.42
1:M:255:GLN:HA	1:M:274:TRP:HB2	2.02	0.42
1:O:12:VAL:CG2	1:O:94:THR:HG23	2.50	0.42
1:W:28:VAL:O	1:W:29:ASP:HB2	2.20	0.42
1:Y:172:LEU:O	1:Y:180:GLN:NE2	2.52	0.42
1:Y:59:TYR:O	1:Y:63:GLU:HG2	2.20	0.42
2:D:33:SER:HB2	2:D:54:LEU:HD21	2.02	0.42
1:E:28:VAL:HG23	1:E:33:PHE:CE2	2.55	0.42
1:K:135:ALA:HB1	1:K:140:ALA:CB	2.50	0.42
1:I:156:LEU:HD22	1:I:160:LEU:CD1	2.50	0.41
1:O:109:PHE:HD2	1:O:165:VAL:HG21	1.84	0.41
1:O:219:ARG:HG3	1:O:257:TYR:CZ	2.55	0.41
1:A:220:ASP:OD2	1:A:256:ARG:NH2	2.47	0.41
1:C:13:SER:HB3	1:C:78:LEU:HD13	2.02	0.41
1:E:168:LEU:HD12	1:E:172:LEU:HG	2.03	0.41
1:E:218:GLN:O	1:E:257:TYR:HA	2.21	0.41
2:P:15:ALA:HB2	2:P:95:TRP:HZ2	1.84	0.41
2:R:2:GLN:HB3	2:R:86:THR:HG22	2.01	0.41
1:S:238:ASP:OD1	1:S:240:THR:OG1	2.36	0.41
1:W:230:LEU:CD2	1:W:245:ALA:HB2	2.50	0.41
1:Y:116:TYR:HB3	1:Y:124:ILE:HG22	2.01	0.41
1:W:160:LEU:O	1:W:165:VAL:HG23	2.19	0.41
1:W:33:PHE:CD1	1:W:34:VAL:HG13	2.55	0.41
1:G:19:GLU:HB3	1:G:20:PRO:HD2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:24:ASN:HB3	2:N:65:LEU:HD11	2.02	0.41
1:Q:137:ASP:C	1:Q:137:ASP:OD1	2.58	0.41
2:B:54:LEU:HD12	2:B:64:LEU:HG	2.51	0.41
1:E:90:ALA:HB1	2:J:85:VAL:CG2	2.41	0.41
2:J:23:LEU:O	2:J:67:TYR:HA	2.20	0.41
2:J:31:HIS:ND1	2:J:32:PRO:HA	2.35	0.41
1:M:19:GLU:CA	1:M:19:GLU:OE1	2.69	0.41
1:W:133:TRP:HB2	1:W:144:LYS:HZ3	1.86	0.41
1:Y:51:TRP:CZ2	1:Y:179:LEU:HD11	2.54	0.41
2:B:23:LEU:HB3	2:B:68:THR:HG22	2.02	0.41
1:C:22:PHE:HE2	1:C:67:VAL:HG22	1.85	0.41
1:W:231:VAL:HG11	1:W:244:TRP:CE2	2.55	0.41
1:A:85:TYR:CE2	1:A:139:ALA:HB3	3.64	0.41
1:E:187:THR:HA	1:E:204:TRP:O	2.20	0.41
1:E:249:VAL:HG12	1:E:257:TYR:CZ	2.56	0.41
1:E:201:LEU:HD13	1:E:274:TRP:NE1	2.35	0.41
1:Q:235:PRO:HG2	2:R:65:LEU:HD22	2.02	0.41
1:S:231:VAL:HG11	1:S:244:TRP:CE2	2.56	0.41
1:W:126:LEU:HD22	1:W:156:LEU:HD12	2.03	0.41
1:Y:52:ILE:HD12	1:Y:52:ILE:HA	1.95	0.41
2:Z:81:ARG:HB2	2:Z:92:ILE:HD13	2.03	0.41
1:I:138:MET:HG2	1:M:85:TYR:CE1	2.56	0.41
1:S:112:GLY:HA3	1:S:160:LEU:HD13	2.03	0.41
1:Y:45:MET:HE1	1:Y:67:VAL:HB	2.02	0.41
1:Y:95:VAL:HG11	1:Y:116:TYR:OH	2.21	0.41
2:B:37:VAL:HG11	2:B:66:TYR:CD1	2.67	0.41
2:B:54:LEU:HD13	2:B:64:LEU:HD11	2.79	0.41
1:E:37:ASP:HB3	1:E:40:ALA:CB	2.50	0.41
2:P:85:VAL:HG13	2:P:86:THR:N	2.36	0.41
2:R:3:ARG:HB3	2:R:29:GLY:O	2.21	0.41
1:S:150:ALA:HB3	1:S:152:VAL:HG23	2.03	0.41
1:G:22:PHE:CD2	1:G:71:SER:HB3	2.55	0.41
1:Q:201:LEU:HD13	1:Q:274:TRP:NE1	2.36	0.41
1:U:70:HIS:NE2	1:U:99:TYR:OH	2.43	0.41
1:W:126:LEU:HD22	1:W:156:LEU:CD1	2.51	0.41
1:Y:28:VAL:HG11	1:Y:179:LEU:HD13	2.03	0.41
1:E:44:ARG:HA	1:E:64:THR:CG2	2.49	0.41
1:K:82:ARG:CD	1:K:89:GLU:HA	2.51	0.41
2:L:24:ASN:HB3	2:L:65:LEU:HD11	2.03	0.41
2:B:84:HIS:CE1	2:B:86:THR:HG23	2.90	0.40
1:C:117:ALA:HB2	2:D:60:TRP:CZ2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:126:LEU:HD22	1:Q:156:LEU:HD13	2.03	0.40
1:U:175:GLY:O	1:U:179:LEU:N	2.51	0.40
2:B:7:ILE:HG12	2:B:27:VAL:HG12	2.03	0.40
1:Q:74:HIS:HD2	1:Q:95:VAL:HG11	1.86	0.40
1:W:167:TRP:CE3	1:W:170:ARG:HD3	2.56	0.40
2:B:13:HIS:HB2	2:B:21:ASN:HD21	2.79	0.40
1:G:205:ALA:HB2	1:G:215:LEU:HD21	2.03	0.40
1:K:182:THR:HG21	1:K:264:GLU:HG2	2.03	0.40
2:R:24:ASN:HB3	2:R:65:LEU:HD11	2.04	0.40
2:F:40:LEU:HA	2:F:44:GLU:O	2.21	0.40
1:K:28:VAL:HG11	1:K:179:LEU:HD13	2.03	0.40
1:S:178:THR:OG1	1:S:179:LEU:N	2.53	0.40
1:C:90:ALA:CB	2:V:2:GLN:CD	2.88	0.40
1:I:19:GLU:HG2	1:I:75:ARG:HH22	1.87	0.40
1:O:95:VAL:HG13	1:O:116:TYR:CE2	2.57	0.40
1:O:216:THR:O	1:O:260:HIS:N	2.55	0.40
1:Q:78:LEU:CD2	1:Q:95:VAL:HG23	2.52	0.40
1:U:55:GLU:OE2	1:U:170:ARG:NE	2.53	0.40
1:Y:64:THR:O	1:Y:67:VAL:HG12	2.21	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	274/276 (99%)	263 (96%)	10 (4%)	1 (0%)	38	75
1	C	274/276 (99%)	262 (96%)	12 (4%)	0	100	100
1	E	274/276 (99%)	262 (96%)	12 (4%)	0	100	100
1	G	274/276 (99%)	264 (96%)	9 (3%)	1 (0%)	38	75
1	I	274/276 (99%)	268 (98%)	6 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	K	274/276 (99%)	264 (96%)	10 (4%)	0	100	100
1	M	274/276 (99%)	267 (97%)	7 (3%)	0	100	100
1	O	274/276 (99%)	265 (97%)	9 (3%)	0	100	100
1	Q	274/276 (99%)	267 (97%)	7 (3%)	0	100	100
1	S	274/276 (99%)	264 (96%)	10 (4%)	0	100	100
1	U	274/276 (99%)	265 (97%)	9 (3%)	0	100	100
1	W	274/276 (99%)	266 (97%)	8 (3%)	0	100	100
1	Y	274/276 (99%)	258 (94%)	16 (6%)	0	100	100
1	a	274/276 (99%)	260 (95%)	14 (5%)	0	100	100
2	B	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
2	D	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
2	F	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
2	H	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
2	J	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
2	L	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
2	N	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
2	P	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
2	R	98/100 (98%)	92 (94%)	6 (6%)	0	100	100
2	T	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
2	V	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
2	X	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
2	Z	98/100 (98%)	97 (99%)	1 (1%)	0	100	100
2	b	98/100 (98%)	94 (96%)	3 (3%)	1 (1%)	18	57
3	c	7/9 (78%)	7 (100%)	0	0	100	100
3	e	7/9 (78%)	7 (100%)	0	0	100	100
3	f	7/9 (78%)	7 (100%)	0	0	100	100
3	g	7/9 (78%)	7 (100%)	0	0	100	100
3	h	7/9 (78%)	7 (100%)	0	0	100	100
3	i	7/9 (78%)	7 (100%)	0	0	100	100
3	j	7/9 (78%)	7 (100%)	0	0	100	100
3	k	7/9 (78%)	7 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	l	7/9 (78%)	7 (100%)	0	0	100	100
3	m	7/9 (78%)	7 (100%)	0	0	100	100
3	n	7/9 (78%)	7 (100%)	0	0	100	100
3	o	7/9 (78%)	7 (100%)	0	0	100	100
3	p	7/9 (78%)	7 (100%)	0	0	100	100
3	q	7/9 (78%)	7 (100%)	0	0	100	100
All	All	5306/5390 (98%)	5121 (96%)	182 (3%)	3 (0%)	55	88

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	GLU
1	G	267	PRO
2	b	35	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain 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 sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	232/232 (100%)	229 (99%)	3 (1%)	73	91
1	C	232/232 (100%)	218 (94%)	14 (6%)	22	58
1	E	232/232 (100%)	217 (94%)	15 (6%)	20	55
1	G	232/232 (100%)	224 (97%)	8 (3%)	42	77
1	I	232/232 (100%)	220 (95%)	12 (5%)	27	63
1	K	232/232 (100%)	215 (93%)	17 (7%)	16	50
1	M	232/232 (100%)	223 (96%)	9 (4%)	37	73
1	O	232/232 (100%)	221 (95%)	11 (5%)	30	67
1	Q	232/232 (100%)	228 (98%)	4 (2%)	66	88
1	S	232/232 (100%)	219 (94%)	13 (6%)	25	61
1	U	232/232 (100%)	219 (94%)	13 (6%)	25	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	W	232/232 (100%)	221 (95%)	11 (5%)	30	67
1	Y	232/232 (100%)	218 (94%)	14 (6%)	22	58
1	a	232/232 (100%)	220 (95%)	12 (5%)	27	63
2	B	95/95 (100%)	89 (94%)	6 (6%)	21	56
2	D	95/95 (100%)	90 (95%)	5 (5%)	26	63
2	F	95/95 (100%)	90 (95%)	5 (5%)	26	63
2	H	95/95 (100%)	89 (94%)	6 (6%)	21	56
2	J	95/95 (100%)	92 (97%)	3 (3%)	44	78
2	L	95/95 (100%)	91 (96%)	4 (4%)	34	71
2	N	95/95 (100%)	89 (94%)	6 (6%)	21	56
2	P	95/95 (100%)	92 (97%)	3 (3%)	44	78
2	R	95/95 (100%)	89 (94%)	6 (6%)	21	56
2	T	95/95 (100%)	92 (97%)	3 (3%)	44	78
2	V	95/95 (100%)	90 (95%)	5 (5%)	26	63
2	X	95/95 (100%)	89 (94%)	6 (6%)	21	56
2	Z	95/95 (100%)	88 (93%)	7 (7%)	16	49
2	b	95/95 (100%)	90 (95%)	5 (5%)	26	63
3	c	9/9 (100%)	8 (89%)	1 (11%)	7	29
3	e	9/9 (100%)	9 (100%)	0	100	100
3	f	9/9 (100%)	8 (89%)	1 (11%)	7	29
3	g	9/9 (100%)	9 (100%)	0	100	100
3	h	9/9 (100%)	8 (89%)	1 (11%)	7	29
3	i	9/9 (100%)	9 (100%)	0	100	100
3	j	9/9 (100%)	8 (89%)	1 (11%)	7	29
3	k	9/9 (100%)	9 (100%)	0	100	100
3	l	9/9 (100%)	9 (100%)	0	100	100
3	m	9/9 (100%)	8 (89%)	1 (11%)	7	29
3	n	9/9 (100%)	8 (89%)	1 (11%)	7	29
3	o	9/9 (100%)	9 (100%)	0	100	100
3	p	9/9 (100%)	9 (100%)	0	100	100
3	q	9/9 (100%)	9 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	4704/4704 (100%)	4472 (95%)	232 (5%)	29 66

All (232) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
1	A	110	LEU
1	A	113	TYR
2	B	4	THR
2	B	57	SER
2	B	70	PHE
2	B	77	GLU
2	B	89	GLN
2	B	97	ARG
3	m	9	VAL
1	C	35	ARG
1	C	42	SER
1	C	106	ASP
1	C	110	LEU
1	C	113	TYR
1	C	131	ARG
1	C	156	LEU
1	C	194	VAL
1	C	195	SER
1	C	207	SER
1	C	226	GLN
1	C	227	ASP
1	C	255	GLN
1	C	268	LYS
2	D	39	LEU
2	D	45	ARG
2	D	70	PHE
2	D	97	ARG
2	D	98	ASP
1	E	11	SER
1	E	35	ARG
1	E	58	GLU
1	E	89	GLU
1	E	110	LEU
1	E	113	TYR
1	E	128	GLU
1	E	131	ARG

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Mol	Chain	Res	Type
1	E	141	GLN
1	E	156	LEU
1	E	227	ASP
1	E	247	VAL
1	E	255	GLN
1	E	268	LYS
1	E	273	ARG
2	F	36	GLU
2	F	39	LEU
2	F	70	PHE
2	F	85	VAL
2	F	89	GLN
1	G	89	GLU
1	G	110	LEU
1	G	156	LEU
1	G	165	VAL
1	G	197	HIS
1	G	225	THR
1	G	270	LEU
1	G	275	GLU
2	H	4	THR
2	H	47	GLU
2	H	70	PHE
2	H	75	LYS
2	H	85	VAL
2	H	89	GLN
3	f	7	THR
1	I	2	SER
1	I	17	ARG
1	I	35	ARG
1	I	86	ASN
1	I	94	THR
1	I	105	SER
1	I	110	LEU
1	I	156	LEU
1	I	168	LEU
1	I	194	VAL
1	I	230	LEU
1	I	268	LYS
2	J	47	GLU
2	J	49	VAL
2	J	70	PHE

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Mol	Chain	Res	Type
1	K	2	SER
1	K	6	ARG
1	K	34	VAL
1	K	35	ARG
1	K	58	GLU
1	K	72	GLN
1	K	110	LEU
1	K	111	ARG
1	K	113	TYR
1	K	132	SER
1	K	156	LEU
1	K	178	THR
1	K	194	VAL
1	K	207	SER
1	K	226	GLN
1	K	228	THR
1	K	275	GLU
2	L	39	LEU
2	L	44	GLU
2	L	51	HIS
2	L	70	PHE
3	h	9	VAL
1	M	17	ARG
1	M	19	GLU
1	M	35	ARG
1	M	113	TYR
1	M	156	LEU
1	M	194	VAL
1	M	230	LEU
1	M	268	LYS
1	M	275	GLU
2	N	4	THR
2	N	34	ASP
2	N	38	ASP
2	N	39	LEU
2	N	58	LYS
2	N	75	LYS
1	O	4	SER
1	O	35	ARG
1	O	89	GLU
1	O	113	TYR
1	O	163	THR

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Mol	Chain	Res	Type
1	O	194	VAL
1	O	207	SER
1	O	222	GLU
1	O	224	GLN
1	O	227	ASP
1	O	268	LYS
2	P	4	THR
2	P	57	SER
2	P	70	PHE
3	n	4	MET
1	Q	35	ARG
1	Q	110	LEU
1	Q	156	LEU
1	Q	181	ARG
2	R	0	MET
2	R	2	GLN
2	R	3	ARG
2	R	48	LYS
2	R	50	GLU
2	R	77	GLU
1	S	25	VAL
1	S	34	VAL
1	S	35	ARG
1	S	67	VAL
1	S	87	GLN
1	S	94	THR
1	S	113	TYR
1	S	122	ASP
1	S	142	THR
1	S	163	THR
1	S	178	THR
1	S	194	VAL
1	S	268	LYS
2	T	39	LEU
2	T	75	LYS
2	T	97	ARG
1	U	2	SER
1	U	19	GLU
1	U	35	ARG
1	U	67	VAL
1	U	86	ASN
1	U	113	TYR

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Mol	Chain	Res	Type
1	U	127	LYS
1	U	132	SER
1	U	156	LEU
1	U	163	THR
1	U	178	THR
1	U	207	SER
1	U	230	LEU
2	V	4	THR
2	V	48	LYS
2	V	63	TYR
2	V	68	THR
2	V	97	ARG
3	c	9	VAL
1	W	35	ARG
1	W	113	TYR
1	W	115	GLN
1	W	122	ASP
1	W	156	LEU
1	W	163	THR
1	W	168	LEU
1	W	216	THR
1	W	225	THR
1	W	251	SER
1	W	268	LYS
2	X	39	LEU
2	X	48	LYS
2	X	58	LYS
2	X	70	PHE
2	X	76	ASP
2	X	98	ASP
1	Y	25	VAL
1	Y	35	ARG
1	Y	122	ASP
1	Y	128	GLU
1	Y	129	ASP
1	Y	138	MET
1	Y	154	GLU
1	Y	156	LEU
1	Y	178	THR
1	Y	194	VAL
1	Y	216	THR
1	Y	227	ASP

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Mol	Chain	Res	Type
1	Y	230	LEU
1	Y	270	LEU
2	Z	4	THR
2	Z	34	ASP
2	Z	39	LEU
2	Z	70	PHE
2	Z	75	LYS
2	Z	98	ASP
2	Z	99	MET
1	a	35	ARG
1	a	67	VAL
1	a	89	GLU
1	a	110	LEU
1	a	113	TYR
1	a	121	LYS
1	a	176	LYS
1	a	219	ARG
1	a	226	GLN
1	a	227	ASP
1	a	262	GLN
1	a	270	LEU
2	b	4	THR
2	b	11	SER
2	b	39	LEU
2	b	70	PHE
2	b	75	LYS
3	j	7	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	115	GLN
2	D	2	GLN
1	E	43	GLN
1	E	255	GLN
1	G	224	GLN
1	I	155	GLN
1	M	155	GLN
1	S	32	GLN
1	S	155	GLN
1	U	226	GLN

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Mol	Chain	Res	Type
1	U	242	GLN
1	U	255	GLN
1	Y	32	GLN
1	a	32	GLN
1	a	155	GLN
1	a	262	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 50 ligands modelled in this entry, 30 are monoatomic - leaving 20 for Mogul analysis.

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	GOL	A	302	-	5,5,5	0.60	0	5,5,5	0.74	0
5	GOL	C	302	-	5,5,5	0.51	0	5,5,5	0.35	0
5	GOL	C	303	-	5,5,5	0.34	0	5,5,5	0.22	0
5	GOL	E	302	-	5,5,5	0.53	0	5,5,5	0.78	0
5	GOL	K	302	-	5,5,5	0.33	0	5,5,5	0.62	0
5	GOL	O	302	-	5,5,5	0.40	0	5,5,5	0.26	0
5	GOL	O	303	-	5,5,5	0.37	0	5,5,5	0.28	0
5	GOL	Q	303	-	5,5,5	0.58	0	5,5,5	0.46	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GOL	S	302	-	5,5,5	0.37	0	5,5,5	0.54	0
5	GOL	Y	302	-	5,5,5	0.47	0	5,5,5	0.27	0
5	GOL	a	302	-	5,5,5	0.46	0	5,5,5	0.49	0
5	GOL	c	101	-	5,5,5	0.39	0	5,5,5	0.42	0
5	GOL	e	101	-	5,5,5	0.49	0	5,5,5	0.38	0
5	GOL	f	101	-	5,5,5	0.54	0	5,5,5	0.83	0
5	GOL	g	101	-	5,5,5	0.49	0	5,5,5	0.27	0
5	GOL	h	101	-	5,5,5	0.49	0	5,5,5	0.51	0
5	GOL	i	101	-	5,5,5	0.49	0	5,5,5	0.49	0
5	GOL	l	101	-	5,5,5	0.22	0	5,5,5	0.37	0
5	GOL	o	101	-	5,5,5	0.44	0	5,5,5	0.29	0
5	GOL	q	101	-	5,5,5	0.40	0	5,5,5	0.29	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	GOL	A	302	-	-	0/4/4/4	0/0/0/0
5	GOL	C	302	-	-	0/4/4/4	0/0/0/0
5	GOL	C	303	-	-	0/4/4/4	0/0/0/0
5	GOL	E	302	-	-	0/4/4/4	0/0/0/0
5	GOL	K	302	-	-	0/4/4/4	0/0/0/0
5	GOL	O	302	-	-	0/4/4/4	0/0/0/0
5	GOL	O	303	-	-	0/4/4/4	0/0/0/0
5	GOL	Q	303	-	-	0/4/4/4	0/0/0/0
5	GOL	S	302	-	-	0/4/4/4	0/0/0/0
5	GOL	Y	302	-	-	0/4/4/4	0/0/0/0
5	GOL	a	302	-	-	0/4/4/4	0/0/0/0
5	GOL	c	101	-	-	0/4/4/4	0/0/0/0
5	GOL	e	101	-	-	0/4/4/4	0/0/0/0
5	GOL	f	101	-	-	0/4/4/4	0/0/0/0
5	GOL	g	101	-	-	0/4/4/4	0/0/0/0
5	GOL	h	101	-	-	0/4/4/4	0/0/0/0
5	GOL	i	101	-	-	0/4/4/4	0/0/0/0
5	GOL	l	101	-	-	0/4/4/4	0/0/0/0
5	GOL	o	101	-	-	0/4/4/4	0/0/0/0
5	GOL	q	101	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	302	GOL	1	0
5	O	302	GOL	1	0

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	276/276 (100%)	0.10	2 (0%) 87 75	38, 51, 71, 87	2 (0%)
1	C	276/276 (100%)	0.02	1 (0%) 92 84	34, 47, 68, 82	2 (0%)
1	E	276/276 (100%)	0.05	1 (0%) 92 84	34, 49, 66, 81	1 (0%)
1	G	276/276 (100%)	-0.06	0 100 100	36, 50, 69, 80	1 (0%)
1	I	276/276 (100%)	0.00	1 (0%) 92 84	33, 49, 71, 83	1 (0%)
1	K	276/276 (100%)	0.04	1 (0%) 92 84	33, 51, 65, 72	2 (0%)
1	M	276/276 (100%)	0.01	0 100 100	40, 53, 70, 83	1 (0%)
1	O	276/276 (100%)	0.11	1 (0%) 92 84	39, 54, 69, 78	2 (0%)
1	Q	276/276 (100%)	0.03	1 (0%) 92 84	35, 48, 63, 80	1 (0%)
1	S	276/276 (100%)	0.13	1 (0%) 92 84	41, 58, 74, 82	2 (0%)
1	U	276/276 (100%)	0.03	0 100 100	41, 56, 72, 102	2 (0%)
1	W	276/276 (100%)	0.27	3 (1%) 80 65	41, 60, 74, 87	2 (0%)
1	Y	276/276 (100%)	0.17	3 (1%) 80 65	39, 58, 73, 78	2 (0%)
1	a	276/276 (100%)	0.34	2 (0%) 87 75	45, 63, 78, 98	2 (0%)
2	B	100/100 (100%)	0.04	0 100 100	38, 53, 71, 93	0
2	D	100/100 (100%)	0.07	0 100 100	34, 50, 72, 89	0
2	F	100/100 (100%)	0.07	0 100 100	34, 52, 69, 83	0
2	H	100/100 (100%)	0.09	1 (1%) 82 67	39, 57, 73, 113	0
2	J	100/100 (100%)	0.06	1 (1%) 82 67	38, 53, 70, 85	0
2	L	100/100 (100%)	0.07	0 100 100	38, 53, 70, 86	0
2	N	100/100 (100%)	-0.02	1 (1%) 82 67	42, 55, 73, 155	0
2	P	100/100 (100%)	-0.01	0 100 100	41, 56, 79, 116	0
2	R	100/100 (100%)	-0.03	0 100 100	37, 52, 69, 80	0
2	T	100/100 (100%)	0.07	0 100 100	43, 60, 79, 95	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
2	V	100/100 (100%)	0.05	0 100 100	42, 57, 71, 84	0
2	X	100/100 (100%)	0.00	0 100 100	45, 57, 73, 84	0
2	Z	100/100 (100%)	0.06	0 100 100	44, 58, 75, 83	0
2	b	100/100 (100%)	0.08	0 100 100	48, 61, 75, 85	0
3	c	9/9 (100%)	-0.10	0 100 100	58, 58, 60, 62	0
3	e	9/9 (100%)	-0.02	0 100 100	52, 55, 56, 57	0
3	f	9/9 (100%)	0.13	0 100 100	42, 44, 50, 50	0
3	g	9/9 (100%)	0.20	0 100 100	50, 52, 56, 56	0
3	h	9/9 (100%)	-0.01	0 100 100	47, 49, 53, 54	0
3	i	9/9 (100%)	-0.08	0 100 100	40, 41, 43, 45	0
3	j	9/9 (100%)	0.22	0 100 100	56, 61, 66, 66	0
3	k	9/9 (100%)	-0.14	0 100 100	39, 43, 48, 49	0
3	l	9/9 (100%)	0.18	0 100 100	39, 42, 48, 48	0
3	m	9/9 (100%)	-0.06	0 100 100	42, 47, 49, 51	0
3	n	9/9 (100%)	0.10	0 100 100	53, 55, 58, 58	0
3	o	9/9 (100%)	0.32	0 100 100	55, 61, 64, 64	0
3	p	9/9 (100%)	0.01	0 100 100	38, 42, 45, 46	0
3	q	9/9 (100%)	0.34	0 100 100	67, 71, 73, 73	0
All	All	5390/5390 (100%)	0.08	20 (0%) 92 84	33, 54, 73, 155	23 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	88	SER	3.4
1	W	1	GLY	3.4
1	Q	225	THR	3.3
1	O	88	SER	3.1
1	I	276	PRO	3.0
2	H	98	ASP	2.6
1	a	276	PRO	2.4
1	Y	89	GLU	2.3
2	N	99	MET	2.3
1	E	220	ASP	2.3
1	A	222	GLU	2.3
1	W	252	GLY	2.3
1	C	88	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	88	SER	2.2
1	Y	180	GLN	2.2
1	W	276	PRO	2.1
1	S	160	LEU	2.1
1	Y	1	GLY	2.1
1	a	225	THR	2.0
2	J	99	MET	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	GOL	O	302	6/6	0.94	0.38	2.92	53,54,55,56	0
5	GOL	i	101	6/6	0.91	0.30	2.21	68,68,69,69	0
5	GOL	A	302	6/6	0.96	0.30	1.34	25,28,28,29	0
5	GOL	o	101	6/6	0.93	0.34	1.05	58,58,58,58	0
5	GOL	c	101	6/6	0.91	0.30	0.96	60,61,62,63	0
5	GOL	E	302	6/6	0.95	0.27	0.93	38,39,43,45	0
5	GOL	S	302	6/6	0.78	0.23	0.73	53,55,55,55	0
5	GOL	g	101	6/6	0.93	0.34	0.64	56,56,57,57	0
5	GOL	h	101	6/6	0.90	0.28	0.51	50,51,52,52	0
5	GOL	l	101	6/6	0.91	0.28	0.29	35,38,39,41	0
5	GOL	C	302	6/6	0.97	0.24	0.01	37,39,40,41	0
5	GOL	e	101	6/6	0.94	0.25	-0.01	51,53,54,55	0
5	GOL	q	101	6/6	0.94	0.29	-0.35	61,61,62,62	0
5	GOL	f	101	6/6	0.95	0.22	-0.90	35,37,37,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	GOL	Q	303	6/6	0.98	0.21	-1.04	36,37,38,39	0
5	GOL	a	302	6/6	0.84	0.22	-1.19	57,58,59,59	0
5	GOL	K	302	6/6	0.88	0.19	-	50,52,54,54	0
5	GOL	C	303	6/6	0.81	0.29	-	55,57,58,58	0
5	GOL	O	303	6/6	0.80	0.34	-	79,80,81,82	0
4	CL	H	101	1/1	0.84	0.12	-	83,83,83,83	0
4	CL	P	101	1/1	0.93	0.11	-	70,70,70,70	0
4	CL	N	101	1/1	0.96	0.10	-	65,65,65,65	0
4	CL	A	301	1/1	0.96	0.21	-	66,66,66,66	0
4	CL	E	301	1/1	0.96	0.13	-	44,44,44,44	0
4	CL	F	101	1/1	0.96	0.10	-	68,68,68,68	0
5	GOL	Y	302	6/6	0.86	0.20	-	58,59,59,60	0
4	CL	R	101	1/1	0.95	0.15	-	72,72,72,72	0
4	CL	I	302	1/1	0.98	0.25	-	52,52,52,52	0
4	CL	V	101	1/1	0.83	0.15	-	69,69,69,69	0
4	CL	I	301	1/1	0.88	0.23	-	57,57,57,57	0
4	CL	I	303	1/1	0.97	0.24	-	50,50,50,50	0
4	CL	Q	302	1/1	0.97	0.15	-	57,57,57,57	0
4	CL	b	101	1/1	0.90	0.09	-	73,73,73,73	0
4	CL	G	301	1/1	0.95	0.20	-	54,54,54,54	0
4	CL	W	301	1/1	0.97	0.25	-	59,59,59,59	0
4	CL	J	101	1/1	0.72	0.32	-	67,67,67,67	0
4	CL	K	301	1/1	0.87	0.27	-	69,69,69,69	0
4	CL	O	301	1/1	0.90	0.21	-	59,59,59,59	0
4	CL	L	101	1/1	0.86	0.12	-	74,74,74,74	0
4	CL	a	301	1/1	0.80	0.13	-	67,67,67,67	0
4	CL	S	301	1/1	0.94	0.13	-	52,52,52,52	0
4	CL	U	301	1/1	0.97	0.23	-	64,64,64,64	0
4	CL	T	101	1/1	0.80	0.15	-	63,63,63,63	0
4	CL	D	101	1/1	0.84	0.13	-	76,76,76,76	0
4	CL	Z	101	1/1	0.73	0.15	-	59,59,59,59	0
4	CL	Y	301	1/1	0.87	0.23	-	60,60,60,60	0
4	CL	N	102	1/1	0.91	0.19	-	50,50,50,50	0
4	CL	Q	301	1/1	0.88	0.32	-	46,46,46,46	0
4	CL	B	101	1/1	0.89	0.25	-	64,64,64,64	0
4	CL	C	301	1/1	0.90	0.23	-	60,60,60,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.