



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 02:48 PM GMT

PDB ID : 3TBW  
Title : CRYSTAL STRUCTURE OF THE MURINE CLASS I MAJOR HISTO-COMPATIBILITY COMPLEX H-2DB IN COMPLEX WITH THE LCMV-DERIVED GP33 ALTERED PEPTIDE ligand (A2G, V3P, Y4S)  
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Deposited on : 2011-08-08  
Resolution : 2.15 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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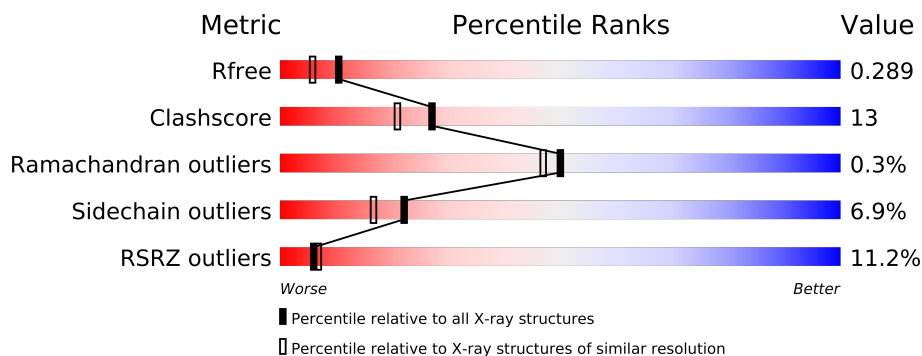
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.15 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))
$R_{free}$	66092	1094 (2.18-2.14)
Clashscore	79885	1299 (2.18-2.14)
Ramachandran outliers	78287	1272 (2.18-2.14)
Sidechain outliers	78261	1272 (2.18-2.14)
RSRZ outliers	66119	1094 (2.18-2.14)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	338	
1	C	338	
1	E	338	
1	G	338	
2	B	99	
2	D	99	
2	F	99	
2	H	99	
3	I	9	
3	J	9	
3	K	9	
3	L	9	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13054 atoms, of which 0 are hydrogen and 0 are deuterium.

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 H-2 class I histocompatibility antigen, D-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	14	0	0
			2248	1420	398	421	9			
1	C	274	Total	C	N	O	S	14	0	0
			2248	1420	398	421	9			
1	E	274	Total	C	N	O	S	30	0	0
			2248	1418	398	423	9			
1	G	267	Total	C	N	O	S	30	0	0
			2196	1387	390	410	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			820	524	138	151	7			
2	D	99	Total	C	N	O	S	0	0	0
			820	524	138	151	7			
2	F	99	Total	C	N	O	S	0	0	0
			820	524	138	151	7			
2	H	99	Total	C	N	O	S	0	0	0
			820	524	138	151	7			

- Molecule 3 is a protein called GLYCOPROTEIN GPC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	9	Total	C	N	O	S	0	0	0
			65	41	11	12	1			
3	J	9	Total	C	N	O	S	0	0	0
			65	41	11	12	1			
3	K	9	Total	C	N	O	S	0	0	0
			65	41	11	12	1			
3	L	9	Total	C	N	O	S	0	0	0
			65	41	11	12	1			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	2	GLY	ALA	ENGINEERED MUTATION	UNP P07399
I	3	PRO	VAL	ENGINEERED MUTATION	UNP P07399
I	4	SER	TYR	ENGINEERED MUTATION	UNP P07399
I	9	MET	CYS	ENGINEERED MUTATION	UNP P07399
J	2	GLY	ALA	ENGINEERED MUTATION	UNP P07399
J	3	PRO	VAL	ENGINEERED MUTATION	UNP P07399
J	4	SER	TYR	ENGINEERED MUTATION	UNP P07399
J	9	MET	CYS	ENGINEERED MUTATION	UNP P07399
K	2	GLY	ALA	ENGINEERED MUTATION	UNP P07399
K	3	PRO	VAL	ENGINEERED MUTATION	UNP P07399
K	4	SER	TYR	ENGINEERED MUTATION	UNP P07399
K	9	MET	CYS	ENGINEERED MUTATION	UNP P07399
L	2	GLY	ALA	ENGINEERED MUTATION	UNP P07399
L	3	PRO	VAL	ENGINEERED MUTATION	UNP P07399
L	4	SER	TYR	ENGINEERED MUTATION	UNP P07399
L	9	MET	CYS	ENGINEERED MUTATION	UNP P07399

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	124	Total O 124 124	0	0
4	B	51	Total O 51 51	0	0
4	C	87	Total O 87 87	0	0
4	D	52	Total O 52 52	0	0
4	E	84	Total O 84 84	0	0
4	F	31	Total O 31 31	0	0
4	G	83	Total O 83 83	0	0
4	H	45	Total O 45 45	0	0
4	I	8	Total O 8 8	0	0
4	J	2	Total O 2 2	0	0
4	K	2	Total O 2 2	0	0

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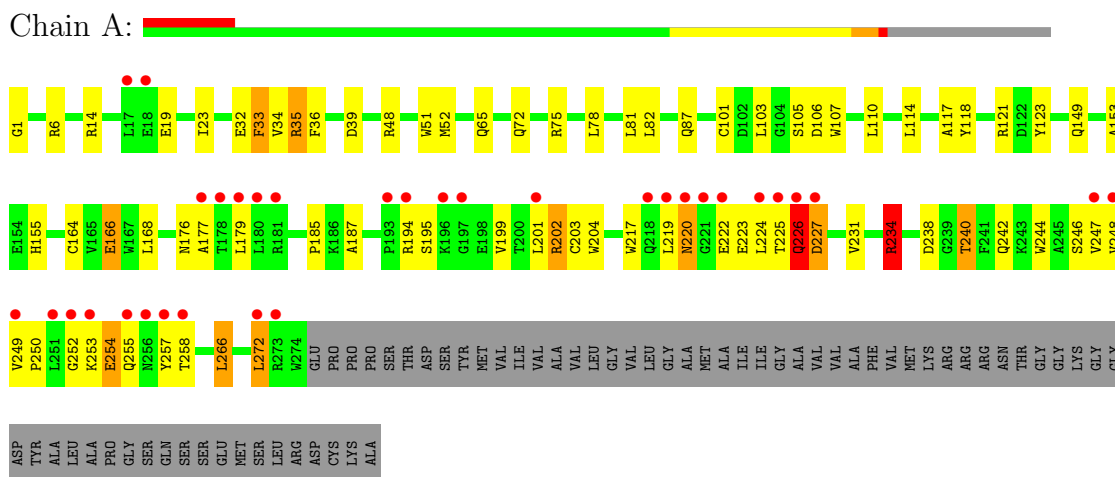
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	5	Total	O	0	0
			5	5		

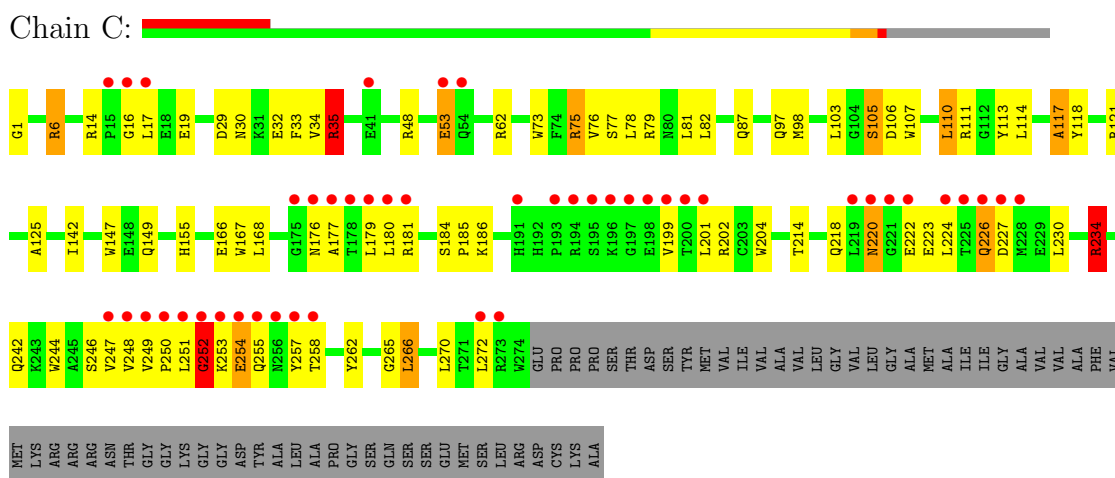
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

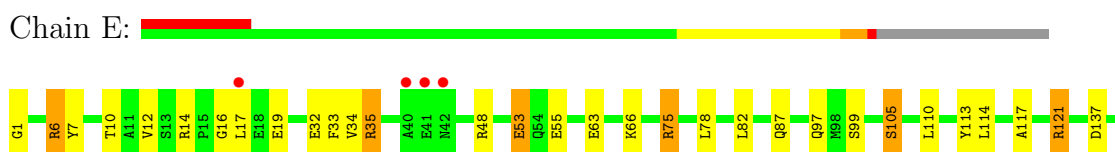
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain

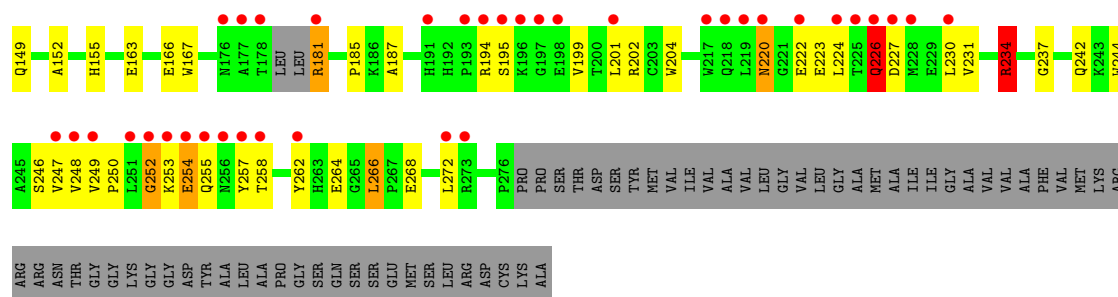


- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



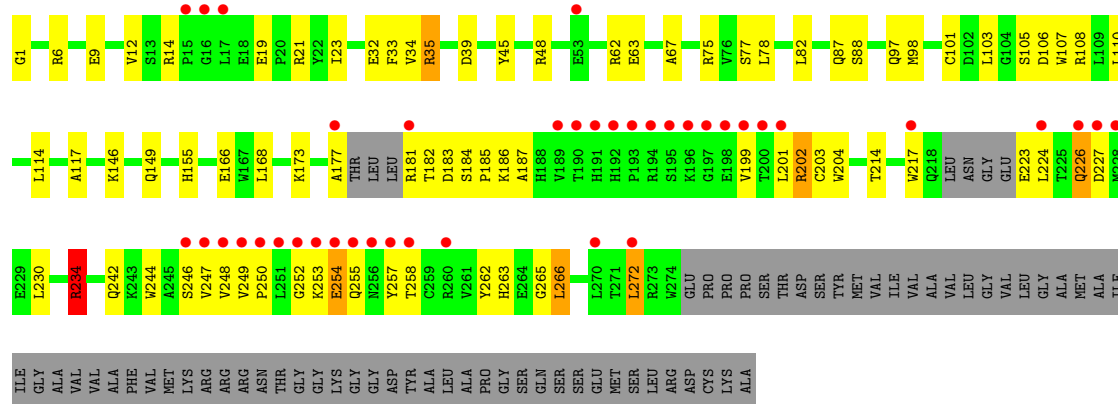
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain





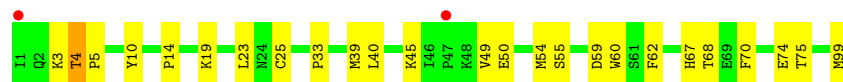
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain

Chain G:



- Molecule 2: Beta-2-microglobulin

Chain B:



- Molecule 2: Beta-2-microglobulin

Chain D:



- Molecule 2: Beta-2-microglobulin

Chain F:



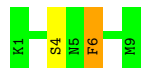
- Molecule 2: Beta-2-microglobulin

Chain H:



- Molecule 3: GLYCOPROTEIN GPC

Chain I: 



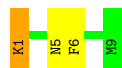
- Molecule 3: GLYCOPROTEIN GPC

Chain J: 



- Molecule 3: GLYCOPROTEIN GPC

Chain K: 



- Molecule 3: GLYCOPROTEIN GPC

Chain L: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.40Å 124.28Å 99.89Å 90.00° 103.23° 90.00°	Depositor
Resolution (Å)	47.48 – 2.15 47.48 – 2.15	Depositor EDS
% Data completeness (in resolution range)	92.6 (47.48-2.15) 98.4 (47.48-2.15)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.24 (at 2.16Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, $R_{free}$	0.232 , 0.283 0.242 , 0.289	Depositor DCC
$R_{free}$ test set	5818 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	35.4	Xtriage
Anisotropy	0.422	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 45.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	3 of 116037 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	13054	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 58.42 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.0489e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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	1.00	4/2314 (0.2%)	1.05	10/3142 (0.3%)
1	C	0.92	3/2314 (0.1%)	0.90	10/3142 (0.3%)
1	E	0.93	2/2314 (0.1%)	0.93	12/3141 (0.4%)
1	G	0.91	2/2260 (0.1%)	0.89	6/3065 (0.2%)
2	B	1.01	1/846 (0.1%)	0.89	1/1148 (0.1%)
2	D	1.02	3/846 (0.4%)	0.89	0/1148
2	F	0.93	1/846 (0.1%)	0.87	1/1148 (0.1%)
2	H	0.91	0/846	0.89	1/1148 (0.1%)
3	I	1.96	2/66 (3.0%)	1.16	0/87
3	J	1.34	0/66	1.17	0/87
3	K	1.37	0/66	1.20	0/87
3	L	1.35	0/66	1.27	1/87 (1.1%)
All	All	0.96	18/12850 (0.1%)	0.94	42/17430 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	1
1	E	0	2
1	G	0	1
All	All	0	5

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	4	SER	CA-CB	9.23	1.66	1.52
2	B	25	CYS	CB-SG	-7.80	1.69	1.82
1	E	152	ALA	CA-CB	7.47	1.68	1.52
1	A	164	CYS	CB-SG	7.12	1.94	1.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	7	TYR	CD2-CE2	6.79	1.49	1.39
1	C	117	ALA	CA-CB	6.67	1.66	1.52
2	D	10	TYR	CD1-CE1	6.33	1.48	1.39
1	G	101	CYS	CB-SG	6.07	1.92	1.82
1	A	33	PHE	CE1-CZ	5.84	1.48	1.37
1	C	76	VAL	CB-CG2	5.70	1.64	1.52
2	D	93	VAL	CB-CG1	-5.65	1.41	1.52
2	F	60	TRP	CE3-CZ3	5.47	1.47	1.38
1	A	123	TYR	CE1-CZ	5.45	1.45	1.38
1	A	101	CYS	CB-SG	5.40	1.91	1.82
1	G	9	GLU	CG-CD	5.25	1.59	1.51
1	C	125	ALA	CA-CB	5.05	1.63	1.52
3	I	6	PHE	CB-CG	5.02	1.59	1.51
2	D	56	PHE	CD1-CE1	5.01	1.49	1.39

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	ARG	NE-CZ-NH1	19.05	129.82	120.30
1	A	234	ARG	NE-CZ-NH2	-15.87	112.36	120.30
1	A	35	ARG	NE-CZ-NH2	-15.47	112.56	120.30
1	A	35	ARG	NE-CZ-NH1	12.80	126.70	120.30
1	E	234	ARG	NE-CZ-NH2	-12.28	114.16	120.30
1	E	234	ARG	NE-CZ-NH1	11.26	125.93	120.30
1	G	35	ARG	NE-CZ-NH2	-10.81	114.90	120.30
1	A	6	ARG	NE-CZ-NH2	-10.49	115.06	120.30
1	E	121	ARG	NE-CZ-NH2	-9.75	115.42	120.30
1	C	35	ARG	NE-CZ-NH2	-8.68	115.96	120.30
1	C	234	ARG	NE-CZ-NH1	8.24	124.42	120.30
1	C	35	ARG	NE-CZ-NH1	7.91	124.25	120.30
1	G	234	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	E	121	ARG	NE-CZ-NH1	7.57	124.09	120.30
1	G	234	ARG	NE-CZ-NH2	-7.47	116.57	120.30
1	A	121	ARG	NE-CZ-NH2	-7.32	116.64	120.30
1	E	35	ARG	NE-CZ-NH1	7.11	123.85	120.30
1	A	220	ASN	N-CA-C	6.58	128.75	111.00
1	E	53	GLU	CB-CA-C	-6.51	97.38	110.40
1	A	121	ARG	NE-CZ-NH1	6.13	123.37	120.30
1	G	39	ASP	CB-CA-C	-6.07	98.25	110.40
2	H	97	ARG	NE-CZ-NH2	-6.03	117.29	120.30
1	E	6	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	C	75	ARG	NE-CZ-NH1	5.82	123.21	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	35	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	C	6	ARG	NE-CZ-NH2	-5.67	117.46	120.30
1	C	234	ARG	NE-CZ-NH2	-5.67	117.46	120.30
1	C	121	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	E	137	ASP	CB-CG-OD2	-5.57	113.29	118.30
1	G	6	ARG	NE-CZ-NH2	-5.55	117.52	120.30
3	L	9	MET	CG-SD-CE	5.49	108.99	100.20
2	B	59	ASP	CB-CG-OD1	5.40	123.16	118.30
1	A	39	ASP	CB-CG-OD1	5.38	123.14	118.30
1	C	252	GLY	N-CA-C	5.33	126.42	113.10
1	E	35	ARG	NE-CZ-NH2	-5.23	117.69	120.30
1	A	6	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	C	53	GLU	CB-CA-C	-5.21	99.99	110.40
2	F	59	ASP	CB-CG-OD1	5.17	122.95	118.30
1	E	220	ASN	N-CA-C	5.07	124.68	111.00
1	E	252	GLY	N-CA-C	5.03	125.69	113.10
1	C	220	ASN	N-CA-C	5.00	124.51	111.00
1	E	75	ARG	NE-CZ-NH2	-5.00	117.80	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	252	GLY	Peptide
1	C	252	GLY	Peptide
1	E	252	GLY	Peptide
1	E	53	GLU	Peptide
1	G	252	GLY	Peptide

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2248	0	2123	57	0
1	C	2248	0	2123	65	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2248	0	2113	57	0
1	G	2196	0	2066	74	0
2	B	820	0	796	20	0
2	D	820	0	796	23	0
2	F	820	0	796	23	0
2	H	820	0	796	33	0
3	I	65	0	66	1	0
3	J	65	0	66	9	0
3	K	65	0	66	8	0
3	L	65	0	66	7	0
4	A	124	0	0	11	0
4	B	51	0	0	2	0
4	C	87	0	0	5	0
4	D	52	0	0	5	0
4	E	84	0	0	4	0
4	F	31	0	0	3	0
4	G	83	0	0	9	0
4	H	45	0	0	3	0
4	I	8	0	0	0	0
4	J	2	0	0	0	0
4	K	2	0	0	0	0
4	L	5	0	0	0	0
All	All	13054	0	11873	315	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (315) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:39:MET:HE2	2:H:49:VAL:HG13	1.41	1.02
2:D:39:MET:SD	4:D:190:HOH:O	2.18	1.01
2:F:39:MET:HE2	2:F:49:VAL:HG13	1.46	0.97
1:G:184:SER:HA	4:G:389:HOH:O	1.73	0.89
2:D:39:MET:HE2	2:D:49:VAL:HG13	1.56	0.87
1:C:167:TRP:CE2	3:J:1:LYS:HD2	2.10	0.86
2:B:39:MET:HE2	2:B:49:VAL:HG13	1.59	0.83
1:C:186:LYS:HE3	1:G:177:ALA:HB2	1.64	0.80
1:G:146:LYS:NZ	4:G:450:HOH:O	2.15	0.79
1:C:97:GLN:HE22	3:J:5:ASN:HD21	1.33	0.76
1:G:185:PRO:HD2	1:G:266:LEU:HD13	1.68	0.76
1:A:234:ARG:HD3	2:B:10:TYR:CE2	2.22	0.74
2:D:39:MET:CE	2:D:68:THR:HB	2.18	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:234:ARG:HD3	2:H:10:TYR:CE2	2.24	0.73
1:G:202:ARG:HD2	1:G:244:TRP:CD2	2.24	0.72
1:E:226:GLN:CG	1:E:226:GLN:O	2.38	0.72
1:C:185:PRO:HD2	1:C:266:LEU:HD13	1.72	0.71
2:F:39:MET:HE1	2:F:68:THR:HG22	1.73	0.71
1:C:202:ARG:HD2	1:C:244:TRP:CD2	2.26	0.71
2:H:39:MET:HE2	2:H:49:VAL:CG1	2.19	0.70
1:G:185:PRO:HD3	4:G:389:HOH:O	1.91	0.69
1:A:19:GLU:OE2	1:A:75:ARG:HD2	1.93	0.69
1:C:155:HIS:HB3	3:J:6:PHE:CZ	2.28	0.69
1:G:155:HIS:HB3	3:L:6:PHE:CZ	2.28	0.69
1:A:185:PRO:HD2	1:A:266:LEU:HD13	1.74	0.68
1:E:55:GLU:OE2	4:E:364:HOH:O	2.12	0.68
1:A:223:GLU:HB3	4:A:556:HOH:O	1.94	0.67
1:C:234:ARG:HD3	2:D:10:TYR:CE2	2.29	0.67
1:C:186:LYS:HE3	1:G:177:ALA:CB	2.25	0.67
1:C:19:GLU:OE2	1:C:75:ARG:HD2	1.94	0.66
2:H:39:MET:CE	2:H:68:THR:HB	2.26	0.66
1:A:32:GLU:OE2	1:A:48:ARG:HD2	1.95	0.66
1:G:32:GLU:OE2	1:G:48:ARG:HD2	1.96	0.66
1:E:202:ARG:HD2	1:E:244:TRP:CD2	2.31	0.66
1:A:202:ARG:HD2	1:A:244:TRP:CD2	2.31	0.65
1:G:19:GLU:OE2	1:G:75:ARG:HD2	1.96	0.65
1:G:201:LEU:O	1:G:246:SER:HA	1.96	0.65
2:D:39:MET:HE1	2:D:68:THR:HB	1.79	0.65
1:E:97:GLN:HE22	3:K:5:ASN:HD21	1.44	0.65
1:E:234:ARG:HD3	2:F:10:TYR:CE2	2.31	0.65
2:B:39:MET:HE1	2:B:68:THR:HG22	1.80	0.64
1:G:146:LYS:CD	4:G:450:HOH:O	2.45	0.64
2:H:91:LYS:HD3	4:H:522:HOH:O	1.97	0.64
2:H:9:VAL:CG2	2:H:93:VAL:HG22	2.28	0.64
1:C:53:GLU:O	2:H:19:LYS:CE	2.46	0.63
1:A:226:GLN:CG	1:A:226:GLN:O	2.41	0.63
1:C:201:LEU:O	1:C:246:SER:HA	1.99	0.63
2:F:1:ILE:N	4:F:311:HOH:O	2.32	0.62
1:A:155:HIS:HB3	3:I:6:PHE:CZ	2.34	0.62
2:D:39:MET:HE1	2:D:68:THR:CB	2.30	0.62
1:C:53:GLU:O	2:H:19:LYS:HD3	2.00	0.62
1:C:167:TRP:CZ2	3:J:1:LYS:HD2	2.35	0.62
2:H:9:VAL:HG21	2:H:93:VAL:HG22	1.82	0.61
1:C:53:GLU:O	2:H:19:LYS:CD	2.47	0.61
1:A:234:ARG:HD2	1:A:242:GLN:HB2	1.82	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:146:LYS:HD2	4:G:450:HOH:O	2.01	0.61
1:E:185:PRO:HD2	1:E:266:LEU:HD13	1.83	0.61
1:C:62:ARG:HG2	4:C:401:HOH:O	2.01	0.60
2:F:40:LEU:HD23	2:F:45:LYS:HA	1.83	0.60
2:F:39:MET:HE1	2:F:68:THR:CG2	2.31	0.60
1:A:194:ARG:HG2	4:A:508:HOH:O	2.02	0.60
1:C:167:TRP:CE2	3:J:1:LYS:CD	2.82	0.59
1:C:77:SER:HB3	3:J:9:MET:HG2	1.84	0.59
1:G:103:LEU:HD11	1:G:168:LEU:HD23	1.84	0.59
2:H:39:MET:HE1	2:H:68:THR:HB	1.83	0.59
1:E:19:GLU:OE2	1:E:75:ARG:HD2	2.03	0.58
1:C:79:ARG:NH1	4:C:570:HOH:O	2.36	0.58
1:G:12:VAL:HG21	2:H:33:PRO:HG2	1.86	0.58
1:G:263:HIS:CE1	4:G:389:HOH:O	2.57	0.58
1:E:226:GLN:CD	1:E:226:GLN:O	2.42	0.58
1:G:88:SER:O	4:G:374:HOH:O	2.17	0.57
1:C:222:GLU:OE2	1:C:222:GLU:HA	2.04	0.57
1:E:163:GLU:OE1	3:K:1:LYS:HD3	2.05	0.57
1:C:53:GLU:O	2:H:19:LYS:NZ	2.38	0.56
1:A:201:LEU:O	1:A:246:SER:HA	2.04	0.56
2:F:39:MET:CE	2:F:68:THR:CG2	2.84	0.56
1:C:177:ALA:N	1:G:186:LYS:HZ1	2.04	0.56
2:F:39:MET:CE	2:F:68:THR:HG22	2.35	0.56
1:E:32:GLU:OE2	1:E:48:ARG:HD2	2.06	0.56
2:B:3:LYS:HE3	4:B:118:HOH:O	2.06	0.56
1:G:32:GLU:OE2	1:G:35:ARG:HD2	2.05	0.56
1:E:201:LEU:O	1:E:246:SER:HA	2.07	0.55
2:B:40:LEU:HD23	2:B:45:LYS:HA	1.89	0.55
2:B:39:MET:HE1	2:B:68:THR:CG2	2.36	0.55
1:C:110:LEU:HD13	4:C:575:HOH:O	2.06	0.55
2:D:39:MET:HE1	2:D:68:THR:HG22	1.88	0.55
1:E:262:TYR:CG	1:G:108:ARG:HD3	2.42	0.55
1:E:97:GLN:NE2	3:K:5:ASN:HD21	2.04	0.55
1:A:222:GLU:HA	1:A:222:GLU:OE2	2.07	0.55
1:E:155:HIS:HB3	3:K:6:PHE:CZ	2.41	0.55
2:F:39:MET:CE	2:F:68:THR:HB	2.37	0.55
1:E:223:GLU:C	1:E:224:LEU:HD22	2.27	0.55
2:H:39:MET:HE1	2:H:68:THR:CB	2.37	0.55
2:D:39:MET:HE1	2:D:68:THR:CG2	2.36	0.55
1:E:234:ARG:HD3	2:F:10:TYR:CZ	2.42	0.54
1:A:227:ASP:O	4:A:381:HOH:O	2.18	0.54
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.42	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:39:MET:CE	2:B:68:THR:HB	2.38	0.54
1:C:249:VAL:HG22	1:C:257:TYR:CZ	2.42	0.54
1:C:176:ASN:HB3	1:G:186:LYS:HE2	1.88	0.54
2:H:40:LEU:HD23	2:H:45:LYS:HA	1.90	0.54
2:D:19:LYS:HE3	4:D:309:HOH:O	2.08	0.54
1:E:222:GLU:HA	1:E:222:GLU:OE2	2.07	0.54
1:C:33:PHE:CD2	1:C:34:VAL:HG13	2.43	0.53
1:E:262:TYR:CD2	1:G:108:ARG:HG2	2.43	0.53
1:A:226:GLN:CD	1:A:226:GLN:O	2.47	0.53
1:E:66:LYS:NZ	3:K:1:LYS:HG2	2.23	0.53
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.43	0.53
1:E:226:GLN:HG2	1:E:226:GLN:O	2.09	0.53
1:A:234:ARG:HD3	2:B:10:TYR:CZ	2.43	0.52
1:G:1:GLY:N	1:G:105:SER:HB3	2.24	0.52
1:E:1:GLY:N	1:E:105:SER:HB3	2.24	0.52
1:A:202:ARG:HG2	1:A:204:TRP:NE1	2.25	0.52
1:C:234:ARG:HD2	1:C:242:GLN:HB2	1.91	0.52
1:G:201:LEU:HD12	1:G:249:VAL:HG21	1.90	0.52
1:E:234:ARG:HD2	1:E:242:GLN:HB2	1.91	0.52
1:G:223:GLU:C	1:G:224:LEU:HD22	2.30	0.52
1:G:234:ARG:HD3	2:H:10:TYR:CZ	2.45	0.52
1:A:272:LEU:HD22	4:A:383:HOH:O	2.10	0.52
1:C:1:GLY:N	1:C:105:SER:HB3	2.25	0.51
1:G:247:VAL:HG22	1:G:248:VAL:N	2.25	0.51
1:G:117:ALA:HB2	2:H:60:TRP:CE2	2.45	0.51
1:G:249:VAL:HG22	1:G:257:TYR:CZ	2.46	0.51
1:C:14:ARG:HB3	1:C:14:ARG:NH1	2.25	0.51
1:A:223:GLU:C	1:A:224:LEU:HD22	2.31	0.51
1:E:264:GLU:HG2	4:E:351:HOH:O	2.10	0.51
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.46	0.51
2:D:33:PRO:HG3	2:D:62:PHE:CZ	2.46	0.51
1:A:226:GLN:HG2	1:A:226:GLN:O	2.11	0.50
1:A:195:SER:N	4:A:508:HOH:O	2.43	0.50
1:E:201:LEU:HD12	1:E:249:VAL:HG21	1.94	0.50
2:H:39:MET:CE	2:H:68:THR:CG2	2.89	0.50
1:E:163:GLU:OE1	3:K:1:LYS:CD	2.58	0.50
1:C:32:GLU:OE2	1:C:35:ARG:HD2	2.11	0.50
1:A:244:TRP:HZ2	2:B:99:MET:HE2	1.77	0.50
1:C:218:GLN:NE2	4:C:355:HOH:O	2.45	0.50
1:A:72:GLN:HG2	4:A:389:HOH:O	2.12	0.50
2:B:39:MET:CE	2:B:68:THR:CG2	2.90	0.50
1:E:262:TYR:CD2	1:G:108:ARG:HD3	2.46	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:244:TRP:HZ2	2:F:99:MET:HE2	1.76	0.49
1:C:223:GLU:C	1:C:224:LEU:HD22	2.32	0.49
1:C:226:GLN:CG	1:C:226:GLN:O	2.60	0.49
1:G:77:SER:HB3	3:L:9:MET:HG2	1.95	0.49
1:C:97:GLN:NE2	3:J:5:ASN:HD21	2.04	0.49
1:G:14:ARG:HB3	1:G:14:ARG:NH1	2.26	0.49
2:H:40:LEU:CD1	4:H:446:HOH:O	2.59	0.49
1:C:32:GLU:OE2	1:C:48:ARG:HD2	2.12	0.49
2:F:39:MET:HE1	2:F:68:THR:CB	2.42	0.49
1:E:82:LEU:HD12	1:E:87:GLN:HB2	1.95	0.49
2:H:39:MET:HE1	2:H:67:HIS:C	2.33	0.49
1:C:106:ASP:O	1:C:107:TRP:HB2	2.11	0.48
1:A:14:ARG:NH1	1:A:14:ARG:HB3	2.28	0.48
2:F:21:ASN:OD1	2:F:22:ILE:N	2.44	0.48
2:H:51:MET:HB3	4:H:405:HOH:O	2.13	0.48
1:C:81:LEU:HD13	1:C:118:TYR:CD1	2.48	0.48
1:G:183:ASP:O	4:G:389:HOH:O	2.20	0.48
1:E:99:SER:HB2	4:E:354:HOH:O	2.14	0.48
2:F:91:LYS:NZ	4:F:552:HOH:O	2.47	0.48
1:G:202:ARG:HD3	1:G:244:TRP:CE3	2.49	0.48
1:G:199:VAL:HG21	1:G:254:GLU:OE1	2.13	0.47
1:C:250:PRO:HB2	1:C:253:LYS:HB2	1.96	0.47
1:G:234:ARG:HD2	1:G:242:GLN:HB2	1.95	0.47
1:C:255:GLN:HA	1:C:255:GLN:OE1	2.14	0.47
1:C:98:MET:HG3	2:D:60:TRP:CZ3	2.49	0.47
2:D:23:LEU:O	2:D:67:HIS:HA	2.15	0.47
1:A:223:GLU:CG	4:A:556:HOH:O	2.62	0.47
1:C:82:LEU:HD12	1:C:87:GLN:HB2	1.96	0.47
1:G:249:VAL:HG13	1:G:257:TYR:CE2	2.49	0.47
1:E:249:VAL:HG22	1:E:257:TYR:CZ	2.50	0.47
1:G:82:LEU:HD12	1:G:87:GLN:HB2	1.97	0.47
1:A:166:GLU:OE2	4:A:404:HOH:O	2.20	0.47
2:B:39:MET:CE	2:B:68:THR:HG22	2.43	0.47
1:A:231:VAL:HG13	1:A:244:TRP:CZ2	2.50	0.47
1:C:234:ARG:HD3	2:D:10:TYR:CZ	2.49	0.46
1:E:14:ARG:NH1	1:E:14:ARG:HB3	2.30	0.46
1:E:202:ARG:HG2	1:E:204:TRP:NE1	2.30	0.46
2:D:21:ASN:OD1	2:D:22:ILE:N	2.43	0.46
1:A:247:VAL:HG22	1:A:248:VAL:N	2.31	0.46
2:B:23:LEU:O	2:B:67:HIS:HA	2.16	0.46
1:G:226:GLN:CG	1:G:226:GLN:O	2.62	0.46
2:H:79:ALA:HB2	2:H:94:TYR:CD1	2.50	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:39:MET:CE	2:H:68:THR:HG22	2.45	0.46
1:C:177:ALA:HB2	1:G:186:LYS:HE3	1.97	0.46
1:C:184:SER:OG	1:C:265:GLY:O	2.30	0.46
1:A:187:ALA:HA	1:A:204:TRP:O	2.16	0.46
1:G:249:VAL:CG1	1:G:254:GLU:HA	2.45	0.46
1:G:202:ARG:CD	1:G:244:TRP:CD2	2.98	0.46
2:D:40:LEU:HD23	2:D:45:LYS:HA	1.96	0.46
1:G:202:ARG:HG2	1:G:204:TRP:NE1	2.30	0.46
1:C:201:LEU:HD12	1:C:249:VAL:HG21	1.97	0.46
2:D:39:MET:CE	2:D:68:THR:CB	2.90	0.46
2:B:39:MET:HE1	2:B:68:THR:CB	2.45	0.46
2:D:39:MET:CE	2:D:68:THR:CG2	2.94	0.45
2:B:4:THR:HA	2:B:5:PRO:HD3	1.84	0.45
1:E:250:PRO:HB2	1:E:253:LYS:HB2	1.98	0.45
2:B:33:PRO:HG3	2:B:62:PHE:CZ	2.51	0.45
2:H:64:ILE:CG2	2:H:65:LEU:N	2.78	0.45
1:E:117:ALA:HB2	2:F:60:TRP:CE2	2.52	0.45
1:A:103:LEU:HD11	1:A:168:LEU:HD23	1.97	0.45
2:B:14:PRO:HD3	4:B:157:HOH:O	2.17	0.45
1:G:185:PRO:HD2	1:G:266:LEU:CD1	2.43	0.45
1:C:247:VAL:HG22	1:C:248:VAL:N	2.32	0.45
1:C:176:ASN:HB3	1:G:186:LYS:CE	2.46	0.45
1:E:231:VAL:HG13	1:E:244:TRP:CZ2	2.52	0.45
1:A:1:GLY:N	1:A:105:SER:HB3	2.30	0.45
1:E:199:VAL:HG21	1:E:254:GLU:OE1	2.17	0.45
2:F:9:VAL:CG2	2:F:93:VAL:HG22	2.46	0.45
1:A:202:ARG:HD3	1:A:244:TRP:CE3	2.52	0.45
1:G:106:ASP:O	1:G:107:TRP:HB2	2.16	0.45
1:A:155:HIS:CD2	4:A:466:HOH:O	2.69	0.45
1:A:82:LEU:HD12	1:A:87:GLN:HB2	1.99	0.45
1:G:63:GLU:OE2	3:L:1:LYS:HD3	2.17	0.44
1:E:234:ARG:HD2	1:E:242:GLN:OE1	2.17	0.44
1:A:51:TRP:CZ3	1:A:52:MET:SD	3.10	0.44
1:C:214:THR:HB	1:C:262:TYR:HB2	1.99	0.44
2:H:9:VAL:HG23	2:H:93:VAL:HG22	1.97	0.44
1:G:103:LEU:CD1	1:G:168:LEU:HD23	2.48	0.44
1:E:10:THR:HG22	1:E:12:VAL:HG23	1.99	0.44
1:C:103:LEU:HD11	1:C:168:LEU:HD23	1.99	0.44
1:A:201:LEU:HD12	1:A:249:VAL:HG21	1.99	0.44
2:D:33:PRO:HG3	2:D:62:PHE:CE2	2.53	0.44
1:C:251:LEU:HD23	1:C:252:GLY:N	2.33	0.44
1:E:247:VAL:HG22	1:E:248:VAL:N	2.31	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.52	0.44
4:C:401:HOH:O	3:J:1:LYS:CE	2.66	0.44
1:G:181:ARG:HG2	1:G:182:THR:N	2.32	0.44
1:E:167:TRP:CZ2	3:K:1:LYS:HE3	2.53	0.44
1:E:63:GLU:OE2	3:K:1:LYS:HG3	2.17	0.44
1:E:268:GLU:OE1	1:G:173:LYS:HE2	2.18	0.44
1:E:255:GLN:HA	1:E:255:GLN:OE1	2.17	0.44
1:G:250:PRO:HB2	1:G:253:LYS:HB2	2.00	0.44
1:G:184:SER:OG	1:G:265:GLY:O	2.30	0.43
1:G:255:GLN:HA	1:G:255:GLN:OE1	2.18	0.43
1:E:6:ARG:NH2	1:E:113:TYR:CE1	2.86	0.43
1:E:262:TYR:CD1	1:G:108:ARG:HD3	2.53	0.43
1:C:73:TRP:CD1	3:J:8:THR:HG22	2.53	0.43
2:F:4:THR:HA	2:F:5:PRO:HD3	1.83	0.43
1:G:23:ILE:HD12	2:H:54:MET:SD	2.58	0.43
1:A:65:GLN:NE2	4:A:421:HOH:O	2.51	0.43
1:G:214:THR:HB	1:G:262:TYR:HB2	2.00	0.43
2:H:39:MET:HE1	2:H:68:THR:HG22	2.01	0.43
1:G:155:HIS:HB3	3:L:6:PHE:CE1	2.53	0.43
1:A:36:PHE:C	1:A:36:PHE:CD1	2.92	0.43
1:A:250:PRO:HB2	1:A:253:LYS:HB2	2.00	0.43
2:H:39:MET:HE3	2:H:68:THR:CG2	2.49	0.43
1:G:249:VAL:HG11	1:G:254:GLU:HA	2.00	0.43
1:C:199:VAL:HG21	1:C:254:GLU:OE1	2.18	0.43
1:G:77:SER:HB3	3:L:9:MET:CG	2.48	0.43
1:E:16:GLY:O	1:E:17:LEU:C	2.57	0.43
2:B:39:MET:HE1	2:B:68:THR:HB	2.00	0.43
1:C:202:ARG:HG2	1:C:204:TRP:NE1	2.33	0.43
1:E:32:GLU:OE2	1:E:35:ARG:HD2	2.18	0.43
1:G:97:GLN:NE2	3:L:5:ASN:HD21	2.16	0.43
2:H:21:ASN:OD1	2:H:22:ILE:N	2.47	0.43
2:F:64:ILE:CG2	2:F:65:LEU:N	2.82	0.43
1:A:199:VAL:HG21	1:A:254:GLU:OE1	2.19	0.43
1:A:249:VAL:HG13	1:A:257:TYR:CE2	2.54	0.43
1:C:29:ASP:O	1:C:30:ASN:HB2	2.19	0.43
1:A:225:THR:HG22	1:A:225:THR:O	2.19	0.43
1:E:121:ARG:NH1	2:F:1:ILE:HB	2.34	0.42
2:H:39:MET:CE	2:H:68:THR:CB	2.94	0.42
1:A:249:VAL:HG22	1:A:257:TYR:CE2	2.54	0.42
1:G:203:CYS:HB2	1:G:217:TRP:CZ2	2.54	0.42
1:E:244:TRP:CZ2	2:F:99:MET:CE	3.02	0.42
2:D:45:LYS:NZ	4:D:318:HOH:O	2.52	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:177:ALA:HB3	1:E:181:ARG:HD3	2.00	0.42
1:C:249:VAL:HG13	1:C:257:TYR:CE2	2.54	0.42
1:G:187:ALA:O	1:G:272:LEU:HD11	2.19	0.42
1:G:33:PHE:CD2	1:G:34:VAL:HG13	2.54	0.42
1:G:97:GLN:HE22	3:L:5:ASN:HD21	1.67	0.42
1:E:33:PHE:CD2	1:E:34:VAL:HG13	2.55	0.42
1:C:176:ASN:OD1	1:C:180:LEU:HD13	2.19	0.42
1:G:98:MET:HG3	2:H:60:TRP:CZ3	2.54	0.42
1:G:181:ARG:HG2	1:G:182:THR:H	1.83	0.42
2:D:77:THR:HG22	4:D:558:HOH:O	2.19	0.42
1:C:249:VAL:CG1	1:C:254:GLU:HA	2.50	0.42
1:A:81:LEU:HD13	1:A:118:TYR:CD1	2.55	0.42
1:A:255:GLN:OE1	1:A:255:GLN:HA	2.18	0.42
1:A:23:ILE:HD12	2:B:54:MET:SD	2.60	0.42
1:A:219:LEU:HG	1:A:219:LEU:O	2.19	0.42
1:C:33:PHE:C	1:C:48:ARG:HB2	2.40	0.42
1:C:1:GLY:H2	1:C:105:SER:HB3	1.84	0.42
2:H:39:MET:HE1	2:H:68:THR:CG2	2.50	0.42
1:C:266:LEU:HD21	1:C:270:LEU:HG	2.02	0.42
2:H:64:ILE:HG22	2:H:65:LEU:N	2.34	0.42
1:A:153:ALA:HB3	4:A:509:HOH:O	2.19	0.42
1:E:99:SER:CB	4:E:354:HOH:O	2.67	0.41
1:E:194:ARG:HG2	1:E:195:SER:H	1.84	0.41
1:G:263:HIS:O	1:G:266:LEU:HB2	2.20	0.41
1:C:117:ALA:HB2	2:D:60:TRP:CE2	2.55	0.41
1:A:106:ASP:O	1:A:107:TRP:HB2	2.20	0.41
2:F:39:MET:HE1	2:F:68:THR:HB	2.00	0.41
1:G:33:PHE:C	1:G:48:ARG:HB2	2.41	0.41
1:G:21:ARG:NE	1:G:23:ILE:HD11	2.36	0.41
1:G:45:TYR:CE2	1:G:67:ALA:HB2	2.56	0.41
2:D:9:VAL:CG2	2:D:93:VAL:HG22	2.50	0.41
1:A:266:LEU:HD12	1:A:266:LEU:HA	1.96	0.41
1:A:254:GLU:H	1:A:254:GLU:HG2	1.76	0.41
2:D:98:ASP:HA	4:D:553:HOH:O	2.21	0.41
1:C:6:ARG:NH2	1:C:113:TYR:CE1	2.89	0.41
1:E:187:ALA:HA	1:E:204:TRP:O	2.20	0.41
1:E:244:TRP:HZ2	2:F:99:MET:CE	2.33	0.41
1:E:266:LEU:HD12	1:E:266:LEU:HA	1.92	0.41
1:A:238:ASP:OD1	1:A:240:THR:OG1	2.37	0.41
1:A:32:GLU:OE2	1:A:35:ARG:HD2	2.21	0.40
1:G:62:ARG:HG3	4:G:358:HOH:O	2.20	0.40
1:C:147:TRP:N	1:C:147:TRP:CD1	2.86	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:247:VAL:CG2	1:G:248:VAL:N	2.84	0.40
1:G:202:ARG:CD	1:G:244:TRP:CE3	3.04	0.40
1:C:142:ILE:H	1:C:142:ILE:HG13	1.70	0.40
2:F:39:MET:HE2	2:F:49:VAL:CG1	2.33	0.40
1:C:111:ARG:HD3	1:C:113:TYR:OH	2.22	0.40
1:E:237:GLY:HA3	4:F:207:HOH:O	2.21	0.40
2:B:39:MET:HE2	2:B:49:VAL:CG1	2.41	0.40
1:C:16:GLY:O	1:C:17:LEU:C	2.60	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	272/338 (80%)	254 (93%)	17 (6%)	1 (0%)	43	39
1	C	272/338 (80%)	254 (93%)	18 (7%)	0	100	100
1	E	270/338 (80%)	253 (94%)	16 (6%)	1 (0%)	43	39
1	G	261/338 (77%)	246 (94%)	15 (6%)	0	100	100
2	B	97/99 (98%)	93 (96%)	4 (4%)	0	100	100
2	D	97/99 (98%)	94 (97%)	2 (2%)	1 (1%)	22	13
2	F	97/99 (98%)	93 (96%)	3 (3%)	1 (1%)	22	13
2	H	97/99 (98%)	93 (96%)	3 (3%)	1 (1%)	22	13
3	I	7/9 (78%)	5 (71%)	2 (29%)	0	100	100
3	J	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	K	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	L	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	1491/1784 (84%)	1403 (94%)	83 (6%)	5 (0%)	50	47

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	226	GLN
1	E	226	GLN
2	H	47	PRO
2	D	47	PRO
2	F	47	PRO

### 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	232/280 (83%)	215 (93%)	17 (7%)	20	15
1	C	232/280 (83%)	214 (92%)	18 (8%)	18	12
1	E	232/280 (83%)	216 (93%)	16 (7%)	22	17
1	G	226/280 (81%)	212 (94%)	14 (6%)	26	20
2	B	94/94 (100%)	87 (93%)	7 (7%)	20	14
2	D	94/94 (100%)	88 (94%)	6 (6%)	25	19
2	F	94/94 (100%)	87 (93%)	7 (7%)	20	14
2	H	94/94 (100%)	88 (94%)	6 (6%)	25	19
3	I	7/7 (100%)	7 (100%)	0	100	100
3	J	7/7 (100%)	7 (100%)	0	100	100
3	K	7/7 (100%)	6 (86%)	1 (14%)	5	2
3	L	7/7 (100%)	7 (100%)	0	100	100
All	All	1326/1524 (87%)	1234 (93%)	92 (7%)	22	17

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	LEU
1	A	110	LEU
1	A	114	LEU
1	A	149	GLN
1	A	166	GLU
1	A	176	ASN
1	A	179	LEU

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Mol	Chain	Res	Type
1	A	202	ARG
1	A	220	ASN
1	A	226	GLN
1	A	227	ASP
1	A	234	ARG
1	A	240	THR
1	A	254	GLU
1	A	258	THR
1	A	266	LEU
1	A	272	LEU
2	B	4	THR
2	B	19	LYS
2	B	50	GLU
2	B	55	SER
2	B	70	PHE
2	B	74	GLU
2	B	75	THR
1	C	35	ARG
1	C	78	LEU
1	C	105	SER
1	C	110	LEU
1	C	114	LEU
1	C	149	GLN
1	C	166	GLU
1	C	179	LEU
1	C	181	ARG
1	C	220	ASN
1	C	226	GLN
1	C	227	ASP
1	C	230	LEU
1	C	234	ARG
1	C	254	GLU
1	C	258	THR
1	C	266	LEU
1	C	272	LEU
2	D	4	THR
2	D	19	LYS
2	D	50	GLU
2	D	70	PHE
2	D	74	GLU
2	D	75	THR
1	E	78	LEU

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Mol	Chain	Res	Type
1	E	105	SER
1	E	110	LEU
1	E	114	LEU
1	E	149	GLN
1	E	166	GLU
1	E	181	ARG
1	E	220	ASN
1	E	226	GLN
1	E	227	ASP
1	E	230	LEU
1	E	234	ARG
1	E	254	GLU
1	E	258	THR
1	E	266	LEU
1	E	272	LEU
2	F	4	THR
2	F	19	LYS
2	F	50	GLU
2	F	70	PHE
2	F	74	GLU
2	F	75	THR
2	F	87	MET
1	G	78	LEU
1	G	110	LEU
1	G	114	LEU
1	G	149	GLN
1	G	166	GLU
1	G	202	ARG
1	G	226	GLN
1	G	227	ASP
1	G	230	LEU
1	G	234	ARG
1	G	254	GLU
1	G	258	THR
1	G	266	LEU
1	G	272	LEU
2	H	4	THR
2	H	19	LYS
2	H	50	GLU
2	H	74	GLU
2	H	75	THR
2	H	87	MET

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Mol	Chain	Res	Type
3	K	1	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	GLN
1	A	220	ASN
2	B	38	GLN
1	C	97	GLN
1	C	220	ASN
1	E	30	ASN
1	E	97	GLN
1	E	220	ASN
2	F	38	GLN
1	G	30	ASN
1	G	97	GLN
1	G	192	HIS
2	H	38	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

### 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	273/338 (80%)	0.63	33 (12%) 5 6	20, 49, 123, 195	0
1	C	273/338 (80%)	0.98	46 (16%) 2 2	22, 51, 130, 169	0
1	E	271/338 (80%)	0.84	41 (15%) 3 3	25, 52, 125, 160	0
1	G	264/338 (78%)	0.96	40 (15%) 3 3	22, 51, 123, 159	0
2	B	99/99 (100%)	0.37	2 (2%) 62 67	26, 46, 74, 100	0
2	D	99/99 (100%)	0.24	1 (1%) 79 83	27, 43, 76, 102	0
2	F	99/99 (100%)	0.60	5 (5%) 27 30	30, 51, 78, 103	0
2	H	99/99 (100%)	0.37	2 (2%) 62 67	28, 48, 79, 102	0
3	I	9/9 (100%)	-0.09	0 100 100	21, 27, 36, 37	0
3	J	9/9 (100%)	0.08	0 100 100	25, 33, 44, 61	0
3	K	9/9 (100%)	0.05	0 100 100	31, 37, 42, 74	0
3	L	9/9 (100%)	-0.01	0 100 100	27, 32, 40, 48	0
All	All	1513/1784 (84%)	0.71	170 (11%) 6 7	20, 49, 118, 195	0

All (170) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	177	ALA	13.9
1	C	179	LEU	11.0
1	E	178	THR	10.8
1	A	178	THR	10.7
1	G	193	PRO	9.3
1	G	201	LEU	9.3
1	A	17	LEU	9.2
1	E	177	ALA	9.1
1	A	179	LEU	8.1
1	C	195	SER	8.1
1	G	195	SER	8.0

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Mol	Chain	Res	Type	RSRZ
1	G	177	ALA	8.0
1	G	199	VAL	7.8
1	G	258	THR	7.5
1	C	177	ALA	7.4
1	E	219	LEU	7.4
1	C	224	LEU	7.1
1	E	17	LEU	7.1
1	G	224	LEU	6.8
1	C	252	GLY	6.6
1	G	255	GLN	6.6
1	G	194	ARG	6.5
1	C	253	LYS	6.5
1	E	197	GLY	6.5
1	A	252	GLY	6.4
1	C	194	ARG	6.1
1	G	257	TYR	6.1
1	E	196	LYS	5.9
1	C	256	ASN	5.9
1	G	197	GLY	5.8
1	G	227	ASP	5.8
1	E	226	GLN	5.6
1	G	251	LEU	5.6
1	E	253	LYS	5.5
1	C	225	THR	5.5
1	A	220	ASN	5.5
1	C	199	VAL	5.4
1	A	226	GLN	5.4
1	E	224	LEU	5.3
1	G	249	VAL	5.2
1	G	256	ASN	5.2
1	G	247	VAL	5.2
1	G	253	LYS	5.1
1	A	273	ARG	4.9
1	E	41	GLU	4.8
1	A	227	ASP	4.7
1	E	40	ALA	4.7
1	G	252	GLY	4.7
1	E	220	ASN	4.7
1	E	227	ASP	4.6
1	G	228	MET	4.6
1	C	53	GLU	4.6
1	C	220	ASN	4.6

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Mol	Chain	Res	Type	RSRZ
1	C	41	GLU	4.5
1	C	219	LEU	4.5
1	C	197	GLY	4.4
1	E	256	ASN	4.4
1	G	250	PRO	4.4
1	G	196	LYS	4.4
1	C	249	VAL	4.3
1	C	196	LYS	4.2
1	C	257	TYR	4.1
1	A	193	PRO	4.1
1	E	273	ARG	4.1
1	C	227	ASP	4.1
1	C	248	VAL	4.0
1	E	195	SER	4.0
1	C	250	PRO	4.0
1	E	257	TYR	4.0
1	C	255	GLN	4.0
1	C	251	LEU	3.9
1	G	272	LEU	3.9
1	G	248	VAL	3.9
1	C	247	VAL	3.9
1	E	255	GLN	3.8
1	A	180	LEU	3.7
1	A	196	LYS	3.7
1	C	16	GLY	3.7
1	E	193	PRO	3.6
2	F	48	LYS	3.6
1	E	252	GLY	3.6
1	A	253	LYS	3.6
1	C	191	HIS	3.5
1	E	247	VAL	3.5
1	C	178	THR	3.5
1	C	193	PRO	3.4
1	E	42	ASN	3.4
1	A	272	LEU	3.4
1	A	255	GLN	3.4
1	E	181	ARG	3.4
1	A	249	VAL	3.3
1	E	218	GLN	3.2
1	G	270	LEU	3.2
1	E	228	MET	3.2
1	G	190	THR	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	201	LEU	3.2
1	E	272	LEU	3.2
1	G	191	HIS	3.1
1	A	181	ARG	3.1
1	G	181	ARG	3.1
1	A	197	GLY	3.1
1	C	221	GLY	3.1
1	G	200	THR	3.1
1	C	181	ARG	3.1
1	G	246	SER	3.1
1	C	180	LEU	3.1
1	E	222	GLU	3.0
1	G	254	GLU	3.0
2	F	1	ILE	3.0
1	E	201	LEU	3.0
2	F	47	PRO	3.0
1	A	225	THR	3.0
1	A	256	ASN	2.9
1	A	258	THR	2.9
1	G	17	LEU	2.9
1	E	230	LEU	2.9
1	G	260	ARG	2.9
1	C	17	LEU	2.9
1	G	226	GLN	2.8
1	C	198	GLU	2.8
1	A	194	ARG	2.8
2	D	48	LYS	2.7
2	B	47	PRO	2.7
1	E	249	VAL	2.7
1	C	254	GLU	2.6
1	C	228	MET	2.6
1	E	225	THR	2.6
1	A	219	LEU	2.6
1	C	226	GLN	2.6
1	C	200	THR	2.6
1	A	224	LEU	2.6
1	C	222	GLU	2.6
1	G	16	GLY	2.5
1	C	54	GLN	2.5
1	G	198	GLU	2.5
1	A	257	TYR	2.4
1	G	217	TRP	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	273	ARG	2.4
1	C	258	THR	2.4
1	E	251	LEU	2.3
1	E	194	ARG	2.3
1	C	15	PRO	2.3
1	A	248	VAL	2.3
1	E	248	VAL	2.3
1	E	191	HIS	2.3
2	F	44	LYS	2.3
1	A	251	LEU	2.3
1	C	272	LEU	2.3
1	E	176	ASN	2.2
1	A	222	GLU	2.2
1	E	198	GLU	2.2
1	A	201	LEU	2.2
1	G	53	GLU	2.2
1	G	189	VAL	2.2
1	G	15	PRO	2.2
1	E	258	THR	2.2
1	G	192	HIS	2.2
1	A	218	GLN	2.1
1	A	247	VAL	2.1
1	E	262	TYR	2.1
1	A	221	GLY	2.1
1	C	176	ASN	2.1
1	E	254	GLU	2.1
2	F	46	ILE	2.1
2	H	48	LYS	2.1
1	E	217	TRP	2.1
2	H	1	ILE	2.0
1	C	175	GLY	2.0
1	A	18	GLU	2.0
2	B	1	ILE	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 carbohydrates in this entry.

## 6.4 Ligands

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.