



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:58 AM UTC

PDB ID : 2Q6W / pdb_00002q6w
Title : The structure of HLA-DRA, DRB3*0101 (DR52a) with bound platelet integrin peptide associated with fetal and neonatal alloimmune thrombocytopenia
Authors : Parry, C.S.
Deposited on : 2007-06-05
Resolution : 2.25 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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

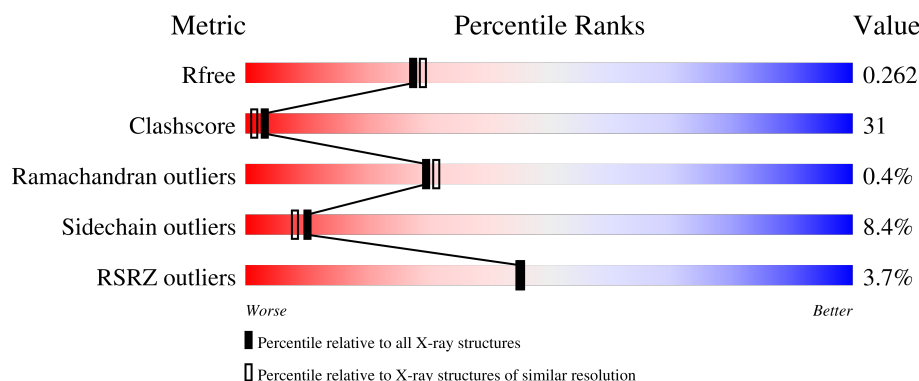
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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

Mol	Chain	Length	Quality of chain
1	A	182	8% 45% 41% 13% .
1	D	182	3% 40% 44% 14% ..
2	B	190	2% 44% 39% 15% ..
2	E	190	8% 44% 46% 7% ..
3	C	12	8% 58% 33% 8%

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Mol	Chain	Length	Quality of chain
3	F	12	8% 25% 75%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6530 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 II histocompatibility antigen, DR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	180	Total	C	N	O	S	0	0	1
			1442	939	235	263	5			
1	D	180	Total	C	N	O	S	3	0	0
			1462	949	236	272	5			

- Molecule 2 is a protein called HLA class II histocompatibility antigen, DRB3-1 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	188	Total	C	N	O	S	0	0	0
			1485	943	267	270	5			
2	E	188	Total	C	N	O	S	0	0	0
			1477	934	271	267	5			

- Molecule 3 is a protein called Integrin beta-3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	11	Total	C	N	O	0	0	0
			85	54	15	16			
3	F	12	Total	C	N	O	2	0	0
			92	57	16	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	26	ARG	CYS	engineered mutation	UNP P05106
F	26	ARG	CYS	engineered mutation	UNP P05106

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	131	Total	O	0	0
			131	131		

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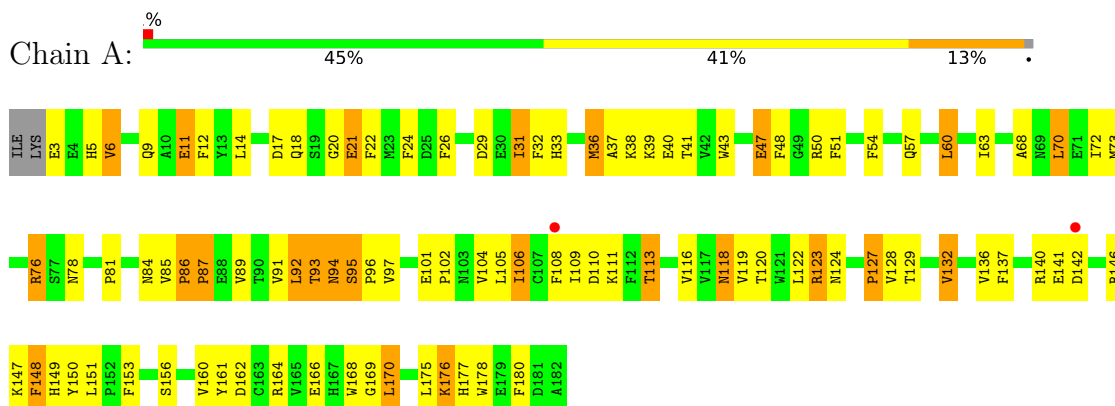
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	124	Total 124	O 124	0	0
4	C	4	Total 4	O 4	0	0
4	D	122	Total 122	O 122	0	0
4	E	99	Total 99	O 99	0	0
4	F	7	Total 7	O 7	0	0

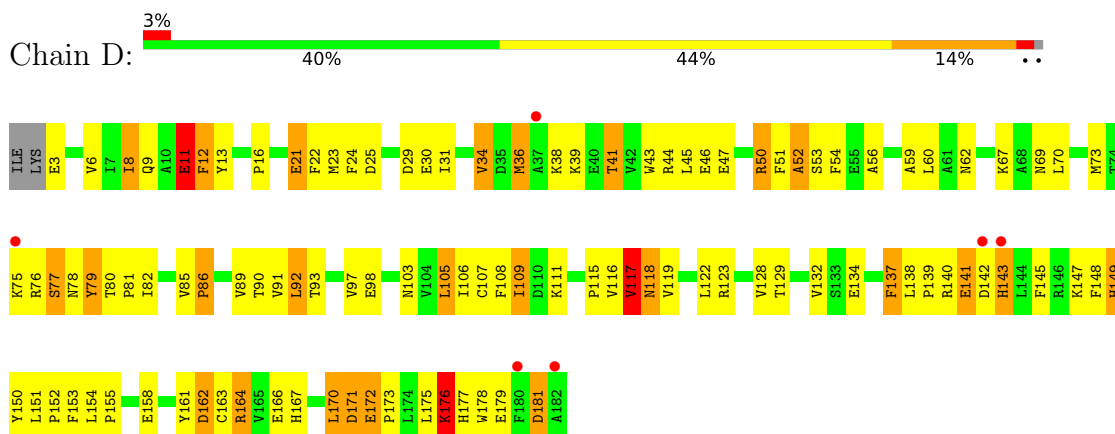
3 Residue-property plots

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

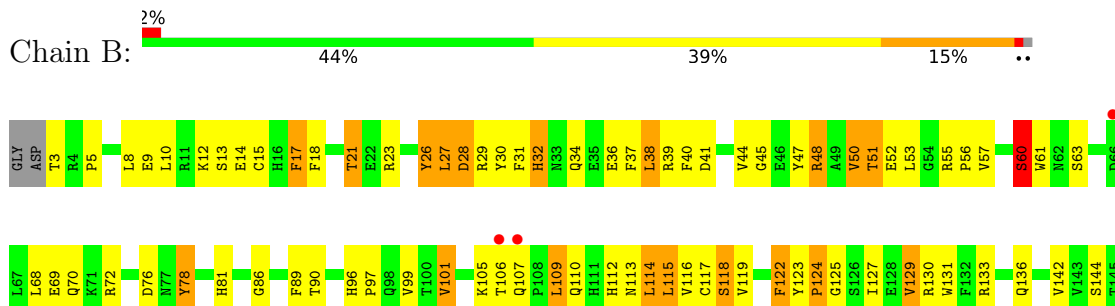
- Molecule 1: HLA class II histocompatibility antigen, DR alpha chain



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

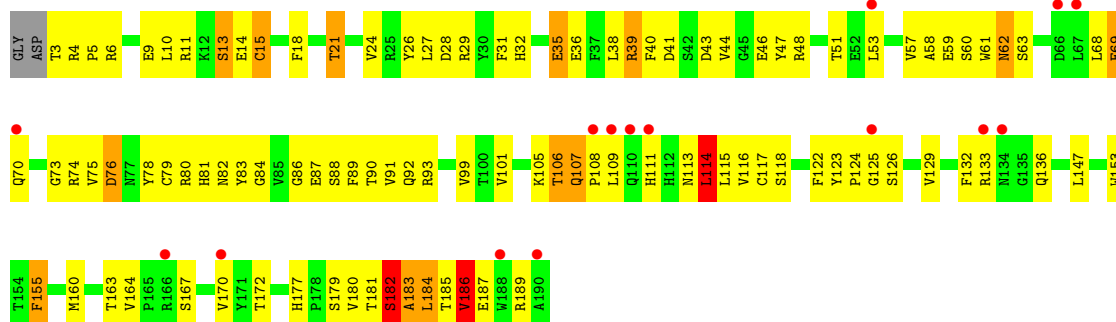


- Molecule 2: HLA class II histocompatibility antigen, DRB3-1 beta chain

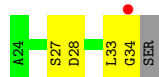




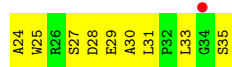
- Molecule 2: HLA class II histocompatibility antigen, DRB3-1 beta chain



- Molecule 3: Integrin beta-3



- Molecule 3: Integrin beta-3



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	92.17Å 92.17Å 248.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.05 – 2.25 20.05 – 2.25	Depositor EDS
% Data completeness (in resolution range)	93.3 (20.05-2.25) 93.1 (20.05-2.25)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.79 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.210 , 0.265 0.206 , 0.262	Depositor DCC
R_{free} test set	3776 reflections (6.83%)	wwPDB-VP
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 31.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6530	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	1.74	16/1485 (1.1%)	1.50	12/2027 (0.6%)
1	D	1.75	23/1506 (1.5%)	1.52	21/2052 (1.0%)
2	B	1.95	35/1523 (2.3%)	1.55	14/2068 (0.7%)
2	E	1.57	8/1517 (0.5%)	1.41	12/2059 (0.6%)
3	C	1.41	0/87	1.40	0/118
3	F	1.44	1/94 (1.1%)	1.44	1/126 (0.8%)
All	All	1.75	83/6212 (1.3%)	1.49	60/8450 (0.7%)

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

All (83) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	70	GLN	C-O	9.48	1.34	1.24
1	D	149	HIS	CA-C	-9.30	1.41	1.52
1	D	82	ILE	CA-CB	8.98	1.63	1.54
3	F	35	SER	CA-CB	8.54	1.70	1.53
2	B	177	HIS	C-O	-8.11	1.15	1.24
2	B	76	ASP	CA-C	-7.83	1.45	1.53
1	A	153	PHE	N-CA	-7.75	1.37	1.46
2	B	122	PHE	N-CA	-7.48	1.37	1.45
2	B	118	SER	N-CA	-7.41	1.37	1.46
2	B	142	VAL	CA-CB	7.31	1.62	1.53
1	D	52	ALA	CA-C	-7.22	1.43	1.52
1	D	8	ILE	CA-C	-6.87	1.44	1.52
1	D	109	ILE	C-O	-6.67	1.17	1.24
1	A	106	ILE	N-CA	-6.66	1.38	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	171	TYR	CA-C	-6.65	1.44	1.52
1	D	164	ARG	N-CA	-6.65	1.38	1.46
1	D	150	TYR	CA-C	-6.59	1.44	1.52
2	B	130	ARG	C-O	-6.52	1.16	1.23
2	B	101	VAL	CA-CB	-6.49	1.46	1.54
2	E	15	CYS	C-O	-6.46	1.16	1.24
2	B	119	VAL	C-O	-6.38	1.17	1.24
2	B	116	VAL	C-O	-6.30	1.16	1.24
1	A	118	ASN	CA-C	-6.26	1.45	1.52
2	B	26	TYR	CA-C	6.24	1.60	1.52
1	D	107	CYS	N-CA	-6.22	1.38	1.46
2	B	78	TYR	C-O	-6.20	1.16	1.24
1	D	107	CYS	C-O	-6.19	1.16	1.24
1	A	89	VAL	CA-CB	-6.19	1.46	1.54
1	A	150	TYR	CA-C	-6.13	1.45	1.52
1	A	76	ARG	CA-C	-6.05	1.44	1.52
1	D	137	PHE	CA-C	-6.04	1.44	1.52
1	D	53	SER	N-CA	-6.00	1.38	1.45
1	D	178	TRP	N-CA	-5.91	1.38	1.46
2	B	96	HIS	CA-CB	5.90	1.63	1.53
2	B	81	HIS	N-CA	-5.89	1.39	1.46
2	E	76	ASP	CA-C	-5.88	1.46	1.52
1	A	148	PHE	CA-C	-5.87	1.45	1.52
2	B	129	VAL	CA-CB	-5.82	1.47	1.54
1	A	92	LEU	C-O	-5.79	1.16	1.23
2	B	148	ILE	N-CA	-5.76	1.39	1.46
2	B	175	VAL	N-CA	-5.72	1.39	1.46
1	D	12	PHE	CA-C	-5.71	1.45	1.52
1	A	123	ARG	NE-CZ	5.70	1.39	1.33
2	B	29	ARG	CA-C	-5.68	1.45	1.52
2	B	28	ASP	CA-C	5.67	1.59	1.52
1	D	22	PHE	CA-C	-5.67	1.45	1.52
1	D	163	CYS	N-CA	-5.62	1.39	1.46
2	E	99	VAL	CA-CB	-5.62	1.47	1.54
1	A	86	PRO	CA-C	5.58	1.55	1.52
2	E	117	CYS	N-CA	-5.52	1.39	1.46
1	A	43	TRP	C-O	-5.49	1.17	1.23
1	D	162	ASP	CA-C	-5.49	1.46	1.52
2	B	60	SER	CA-C	-5.47	1.46	1.52
2	E	118	SER	N-CA	-5.44	1.39	1.46
2	B	99	VAL	C-O	-5.44	1.18	1.24
1	A	6	VAL	CA-C	-5.43	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	162	ASP	C-O	-5.42	1.17	1.23
2	E	155	PHE	N-CA	-5.41	1.38	1.45
1	A	31	ILE	N-CA	-5.39	1.38	1.46
2	B	90	THR	CA-CB	5.37	1.60	1.53
1	D	21	GLU	CA-C	-5.37	1.45	1.52
1	A	156	SER	CA-C	-5.32	1.45	1.52
2	E	187	GLU	CA-C	-5.30	1.46	1.52
1	D	106	ILE	CA-CB	5.30	1.60	1.54
1	D	34	VAL	N-CA	-5.27	1.39	1.46
1	A	21	GLU	CA-C	-5.22	1.46	1.52
1	D	162	ASP	N-CA	-5.21	1.39	1.45
2	B	172	THR	C-O	-5.20	1.18	1.24
2	B	174	GLN	CA-C	-5.19	1.46	1.52
2	B	115	LEU	N-CA	-5.19	1.39	1.46
2	B	187	GLU	N-CA	-5.12	1.39	1.45
2	B	86	GLY	C-O	-5.11	1.18	1.23
2	B	125	GLY	CA-C	-5.10	1.45	1.52
2	B	44	VAL	CA-CB	5.08	1.60	1.55
2	E	129	VAL	N-CA	-5.08	1.40	1.46
1	A	127	PRO	N-CA	-5.05	1.40	1.47
2	B	32	HIS	CA-C	-5.05	1.46	1.52
2	B	160	MET	N-CA	-5.04	1.40	1.45
1	D	8	ILE	N-CA	-5.01	1.40	1.46
1	D	118	ASN	CA-C	-5.01	1.46	1.52
2	B	32	HIS	C-O	-5.01	1.18	1.24
2	B	17	PHE	CA-C	-5.00	1.46	1.52
2	B	153	TRP	C-O	-5.00	1.16	1.23

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	69	GLU	CA-C-O	9.66	131.01	120.20
1	D	36	MET	CA-CB-CG	-9.24	95.62	114.10
1	A	146	ARG	N-CA-C	-8.88	95.05	108.99
1	A	31	ILE	N-CA-C	-7.97	104.04	111.45
3	F	35	SER	N-CA-CB	7.63	123.48	110.50
1	A	68	ALA	N-CA-C	-7.57	103.11	111.36
2	B	70	GLN	N-CA-C	-7.54	102.75	110.97
2	B	96	HIS	CA-C-N	7.34	127.10	119.76
2	B	96	HIS	C-N-CA	7.34	127.10	119.76
1	D	86	PRO	CA-C-N	6.97	126.73	119.76
1	D	86	PRO	C-N-CA	6.97	126.73	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	126	SER	N-CA-CB	-6.96	98.47	109.19
2	E	21	THR	N-CA-C	-6.85	102.82	113.02
2	E	126	SER	N-CA-C	6.75	120.63	110.64
2	B	107	GLN	N-CA-C	-6.66	100.08	110.14
1	A	95	SER	N-CA-C	-6.55	98.90	109.58
2	B	51	THR	N-CA-C	-6.50	99.23	109.50
2	B	173	CYS	N-CA-C	-6.50	97.92	108.52
1	A	94	ASN	N-CA-C	6.05	119.92	112.54
1	D	50	ARG	CB-CG-CD	-6.02	97.45	111.30
2	E	21	THR	CB-CA-C	5.95	120.44	110.44
1	D	122	LEU	CA-C-N	-5.92	114.65	123.00
1	D	122	LEU	C-N-CA	-5.92	114.65	123.00
1	D	118	ASN	N-CA-C	-5.92	98.76	108.41
2	E	114	LEU	N-CA-C	-5.81	98.94	108.41
1	A	70	LEU	N-CA-C	-5.81	104.86	111.07
2	B	45	GLY	N-CA-C	-5.80	107.38	114.92
1	D	172	GLU	CA-C-N	5.80	126.14	119.93
1	D	172	GLU	C-N-CA	5.80	126.14	119.93
2	E	186	VAL	CB-CA-C	5.79	119.41	110.50
2	B	44	VAL	N-CA-C	-5.75	106.21	111.67
2	E	179	SER	N-CA-C	-5.75	105.93	113.17
2	E	183	ALA	N-CA-C	-5.74	99.06	108.76
1	A	11	GLU	CA-C-N	-5.64	113.33	122.36
1	A	11	GLU	C-N-CA	-5.64	113.33	122.36
1	D	90	THR	N-CA-C	-5.63	99.84	109.24
1	A	93	THR	N-CA-C	-5.62	101.70	110.42
2	B	21	THR	CB-CA-C	5.61	119.40	109.75
2	E	93	ARG	NE-CZ-NH1	-5.61	115.89	121.50
1	D	176	LYS	CB-CG-CD	-5.58	98.47	111.30
1	A	168	TRP	N-CA-C	5.56	118.10	111.71
1	D	22	PHE	N-CA-C	-5.55	98.87	108.69
1	D	105	LEU	N-CA-C	-5.47	100.33	109.24
2	E	84	GLY	N-CA-C	-5.40	105.85	112.77
1	D	25	ASP	N-CA-C	-5.34	99.98	109.06
1	D	11	GLU	CA-C-N	-5.34	113.56	122.21
1	D	11	GLU	C-N-CA	-5.34	113.56	122.21
1	D	149	HIS	N-CA-C	-5.32	101.10	109.50
2	B	159	VAL	N-CA-C	-5.27	98.90	107.28
1	D	31	ILE	N-CA-C	-5.26	105.71	110.82
1	D	117	VAL	CB-CA-C	-5.26	104.49	111.80
1	D	85	VAL	CA-C-N	-5.25	114.97	120.38
1	D	85	VAL	C-N-CA	-5.25	114.97	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	124	PRO	CA-C-N	5.23	126.68	120.14
2	B	124	PRO	C-N-CA	5.23	126.68	120.14
2	B	63	SER	N-CA-C	-5.15	106.25	112.54
2	E	185	THR	N-CA-C	-5.15	100.06	108.76
1	A	176	LYS	CD-CE-NZ	-5.10	95.58	111.90
1	A	47	GLU	N-CA-C	5.07	118.23	111.75
2	B	52	GLU	CB-CA-C	5.04	120.45	110.17

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	87	PRO	Peptide

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	1442	0	1372	93	0
1	D	1462	0	1378	109	0
2	B	1485	0	1386	70	0
2	E	1477	0	1367	116	0
3	C	85	0	79	9	0
3	F	92	0	84	19	0
4	A	131	0	0	19	0
4	B	124	0	0	17	0
4	C	4	0	0	1	0
4	D	122	0	0	20	0
4	E	99	0	0	20	0
4	F	7	0	0	0	0
All	All	6530	0	5666	364	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 31.

All (364) 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:D:92:LEU:CD2	1:D:108:PHE:CE1	1.92	1.50
1:D:92:LEU:CD2	1:D:108:PHE:HE1	1.19	1.49
1:D:92:LEU:HD21	1:D:108:PHE:CE1	1.51	1.36
2:E:3:THR:HG21	4:E:279:HOH:O	1.05	1.24
2:B:53:LEU:CB	4:B:304:HOH:O	1.87	1.20
1:D:73:MET:HE3	2:E:53:LEU:HD13	1.21	1.14
1:D:92:LEU:HD21	1:D:108:PHE:CZ	1.84	1.11
1:D:92:LEU:HD22	1:D:108:PHE:CE1	1.86	1.11
1:D:36:MET:C	4:D:253:HOH:O	1.95	1.07
1:A:37:ALA:O	1:A:38:LYS:HB3	1.53	1.06
4:D:303:HOH:O	3:F:25:TRP:CE3	2.10	1.04
1:D:73:MET:CE	2:E:53:LEU:HD13	1.89	1.03
1:D:36:MET:O	4:D:253:HOH:O	1.74	1.02
4:E:280:HOH:O	3:F:25:TRP:CE2	2.17	0.97
4:E:280:HOH:O	3:F:25:TRP:CD2	2.15	0.97
1:D:92:LEU:HD23	1:D:108:PHE:HE1	1.26	0.96
1:D:73:MET:HE1	2:E:53:LEU:HB3	1.44	0.96
1:D:81:PRO:HB3	2:E:5:PRO:HB2	1.48	0.95
2:E:53:LEU:HB2	4:E:281:HOH:O	1.67	0.92
1:D:86:PRO:HB2	1:D:170:LEU:HD13	1.52	0.92
2:B:68:LEU:O	2:B:72:ARG:HG3	1.68	0.92
2:B:48:ARG:HG2	2:B:48:ARG:HH11	1.35	0.91
1:D:46:GLU:CA	1:D:47:GLU:N	2.34	0.90
4:D:303:HOH:O	3:F:25:TRP:HE3	1.46	0.89
2:E:81:HIS:HE1	3:F:24:ALA:O	1.56	0.87
1:A:37:ALA:HB1	4:A:303:HOH:O	1.73	0.87
1:A:166:GLU:CG	4:A:229:HOH:O	2.22	0.87
1:D:47:GLU:N	4:D:294:HOH:O	2.07	0.86
2:E:29:ARG:HD2	2:E:36:GLU:OE2	1.74	0.86
1:D:73:MET:O	1:D:77:SER:HB2	1.75	0.86
1:D:118:ASN:ND2	4:D:295:HOH:O	1.95	0.86
4:D:241:HOH:O	2:E:32:HIS:HE1	1.59	0.85
4:D:303:HOH:O	3:F:25:TRP:HB2	1.76	0.85
2:E:113:ASN:HD22	2:E:114:LEU:H	1.21	0.84
1:A:149:HIS:HD2	4:A:262:HOH:O	1.60	0.83
4:A:301:HOH:O	2:B:32:HIS:HE1	1.62	0.83
1:A:37:ALA:O	1:A:38:LYS:CB	2.27	0.82
2:E:26:TYR:OH	3:F:28:ASP:OD2	1.96	0.81
4:D:287:HOH:O	2:E:155:PHE:CE1	2.34	0.78
1:A:149:HIS:CD2	4:A:262:HOH:O	2.35	0.77
2:E:164:VAL:HB	4:E:236:HOH:O	1.85	0.76
1:D:139:PRO:HB3	4:D:273:HOH:O	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:MET:HE3	1:A:60:LEU:HD23	1.69	0.74
2:B:3:THR:N	4:B:295:HOH:O	2.20	0.74
1:D:98:GLU:O	1:D:155:PRO:HG2	1.88	0.74
2:E:167:SER:CA	4:E:259:HOH:O	2.37	0.73
2:E:101:VAL:HG23	2:E:186:VAL:HG22	1.72	0.71
1:A:92:LEU:HB2	1:A:108:PHE:HE1	1.54	0.71
2:E:73:GLY:HA3	4:E:282:HOH:O	1.89	0.70
2:B:36:GLU:O	2:B:50:VAL:HG22	1.91	0.70
1:D:46:GLU:CA	4:D:294:HOH:O	2.39	0.70
1:D:147:LYS:HE3	1:D:149:HIS:HE1	1.55	0.70
2:E:101:VAL:CG2	2:E:186:VAL:HG22	2.22	0.69
4:B:308:HOH:O	1:D:176:LYS:HG2	1.93	0.69
2:E:109:LEU:H	2:E:109:LEU:HD12	1.58	0.69
1:D:86:PRO:CB	1:D:170:LEU:HD13	2.23	0.69
1:A:3:GLU:OE2	1:A:6:VAL:CG2	2.41	0.69
2:E:47:TYR:H	2:E:62:ASN:HD21	1.42	0.68
2:E:43:ASP:OD2	4:E:267:HOH:O	2.11	0.68
2:E:9:GLU:OE1	2:E:32:HIS:HD2	1.77	0.68
2:B:10:LEU:HD21	2:B:12:LYS:HE3	1.76	0.68
2:E:107:GLN:HE21	2:E:107:GLN:N	1.90	0.68
1:A:141:GLU:CB	4:A:269:HOH:O	2.41	0.68
1:D:147:LYS:HE3	1:D:149:HIS:CE1	2.29	0.68
2:E:3:THR:HA	4:E:264:HOH:O	1.94	0.67
2:B:48:ARG:HG2	2:B:48:ARG:NH1	2.08	0.67
1:A:92:LEU:HB2	1:A:108:PHE:CE1	2.30	0.66
2:E:109:LEU:O	4:E:261:HOH:O	2.13	0.66
2:E:3:THR:CG2	2:E:6:ARG:HH12	2.09	0.65
2:B:113:ASN:HD22	2:B:114:LEU:H	1.44	0.65
1:D:76:ARG:HD3	4:E:277:HOH:O	1.96	0.65
1:D:81:PRO:HB3	2:E:5:PRO:CB	2.26	0.65
2:B:160:MET:HE2	4:B:258:HOH:O	1.96	0.65
2:E:57:VAL:HG11	3:F:33:LEU:HD22	1.79	0.64
1:A:86:PRO:HB2	1:A:170:LEU:HD13	1.80	0.64
1:A:70:LEU:HB2	2:B:9:GLU:HB2	1.79	0.64
1:D:50:ARG:HH11	1:D:50:ARG:CG	2.09	0.63
1:A:96:PRO:HG3	2:B:118:SER:OG	1.99	0.62
2:E:3:THR:CG2	4:E:279:HOH:O	1.87	0.62
1:D:129:THR:O	1:D:132:VAL:HG22	2.00	0.62
2:E:113:ASN:ND2	2:E:114:LEU:H	1.94	0.62
1:A:21:GLU:OE1	4:A:283:HOH:O	2.16	0.62
2:B:109:LEU:O	2:B:110:GLN:HB2	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:18:PHE:HB3	4:E:248:HOH:O	2.01	0.61
2:E:111:HIS:HB2	4:E:289:HOH:O	2.01	0.61
1:D:123:ARG:HG3	1:D:161:TYR:CE2	2.36	0.60
1:D:43:TRP:HH2	4:D:303:HOH:O	1.83	0.60
2:B:174:GLN:HG3	2:B:183:ALA:HB1	1.84	0.59
1:D:23:MET:HE2	1:D:30:GLU:HG3	1.84	0.59
1:A:109:ILE:HD11	1:A:119:VAL:HG21	1.84	0.59
2:E:74:ARG:NH2	4:E:256:HOH:O	2.34	0.59
1:D:91:VAL:HG23	1:D:176:LYS:HB3	1.84	0.59
1:A:51:PHE:CB	4:A:310:HOH:O	2.50	0.59
1:A:109:ILE:CD1	1:A:119:VAL:HG21	2.33	0.58
2:E:184:LEU:HD22	2:E:184:LEU:H	1.69	0.58
2:B:57:VAL:HG11	3:C:33:LEU:HD23	1.85	0.58
2:E:101:VAL:HA	2:E:116:VAL:O	2.04	0.58
2:B:68:LEU:O	2:B:72:ARG:CG	2.46	0.58
1:A:51:PHE:CB	4:A:313:HOH:O	2.52	0.58
2:B:133:ARG:NH2	4:B:267:HOH:O	2.36	0.57
1:D:54:PHE:CZ	3:F:27:SER:HB3	2.40	0.57
2:E:28:ASP:O	2:E:39:ARG:HA	2.04	0.57
1:D:141:GLU:OE1	2:E:29:ARG:NH2	2.36	0.57
1:A:37:ALA:CA	4:A:303:HOH:O	2.53	0.57
1:A:37:ALA:CB	4:A:308:HOH:O	2.53	0.57
1:D:176:LYS:HD3	4:D:296:HOH:O	2.05	0.56
2:E:123:TYR:O	2:E:177:HIS:HE1	1.89	0.56
2:B:117:CYS:HB2	2:B:131:TRP:CZ2	2.41	0.56
2:E:3:THR:HG22	2:E:6:ARG:HH12	1.69	0.56
1:A:37:ALA:HB3	4:A:308:HOH:O	2.05	0.56
1:D:176:LYS:CE	4:D:296:HOH:O	2.54	0.56
1:A:111:LYS:HG2	1:A:140:ARG:CZ	2.35	0.56
1:A:57:GLN:HA	1:A:60:LEU:HD12	1.87	0.55
2:B:167:SER:HB3	4:B:270:HOH:O	2.06	0.55
2:B:188:TRP:CH2	2:B:190:ALA:HA	2.41	0.55
1:D:73:MET:HE1	2:E:53:LEU:CB	2.27	0.55
2:E:133:ARG:O	2:E:170:VAL:O	2.24	0.55
1:D:128:VAL:HG11	1:D:151:LEU:HD11	1.88	0.55
1:D:13:TYR:CE2	1:D:67:LYS:HG3	2.42	0.55
2:B:101:VAL:HG11	2:B:188:TRP:HB2	1.88	0.54
1:D:118:ASN:HB3	4:D:298:HOH:O	2.07	0.54
1:D:115:PRO:HD3	1:D:145:PHE:CE2	2.43	0.54
1:A:93:THR:HG21	1:A:97:VAL:CG2	2.38	0.54
2:E:69:GLU:HA	4:E:225:HOH:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:47:TYR:N	2:E:62:ASN:HD21	2.04	0.54
2:E:81:HIS:CE1	3:F:24:ALA:O	2.48	0.54
2:E:31:PHE:HA	2:E:35:GLU:O	2.08	0.54
1:A:123:ARG:O	1:A:124:ASN:C	2.46	0.54
2:B:113:ASN:ND2	2:B:114:LEU:H	2.06	0.54
2:E:108:PRO:HD2	4:E:289:HOH:O	2.07	0.54
2:B:9:GLU:OE1	2:B:32:HIS:HD2	1.91	0.53
1:D:93:THR:HG22	1:D:105:LEU:HD23	1.90	0.53
1:D:3:GLU:HB3	2:E:18:PHE:CE2	2.43	0.53
1:D:16:PRO:HD2	2:E:6:ARG:NE	2.24	0.53
1:A:128:VAL:HG12	1:A:151:LEU:HD11	1.90	0.53
2:B:50:VAL:O	2:E:51:THR:HG22	2.09	0.53
2:E:122:PHE:HE2	2:E:125:GLY:O	1.91	0.53
1:D:77:SER:HB3	4:D:241:HOH:O	2.08	0.52
1:D:79:TYR:N	1:D:79:TYR:CD1	2.77	0.52
1:A:160:VAL:HB	1:A:177:HIS:CE1	2.45	0.52
2:B:185:THR:HG23	4:B:239:HOH:O	2.09	0.52
1:A:38:LYS:NZ	1:A:40:GLU:CD	2.67	0.52
2:B:27:LEU:HD12	2:B:41:ASP:HA	1.89	0.52
1:D:111:LYS:HG2	1:D:140:ARG:CZ	2.39	0.52
2:B:30:TYR:HB2	2:B:38:LEU:HB3	1.92	0.52
2:B:38:LEU:HD11	2:B:61:TRP:CZ3	2.44	0.52
1:D:3:GLU:HA	2:E:18:PHE:CD2	2.45	0.52
2:E:74:ARG:O	2:E:75:VAL:C	2.51	0.52
2:E:123:TYR:CD1	2:E:124:PRO:HA	2.45	0.52
1:D:50:ARG:HH11	1:D:50:ARG:HG2	1.73	0.52
2:E:75:VAL:CG1	2:E:80:ARG:NH1	2.73	0.52
1:D:117:VAL:HG22	1:D:137:PHE:HE2	1.75	0.51
2:E:109:LEU:HD12	2:E:109:LEU:N	2.24	0.51
1:A:180:PHE:HE1	4:A:306:HOH:O	1.93	0.51
2:E:9:GLU:OE1	2:E:32:HIS:CD2	2.61	0.51
1:A:118:ASN:HB2	1:A:166:GLU:HB2	1.93	0.51
1:A:147:LYS:HE2	1:A:149:HIS:HE1	1.75	0.51
1:A:37:ALA:C	4:A:303:HOH:O	2.52	0.51
2:B:38:LEU:CD1	2:B:61:TRP:CZ3	2.94	0.50
2:E:41:ASP:HB3	2:E:44:VAL:HB	1.93	0.50
1:D:50:ARG:HD2	4:D:233:HOH:O	2.11	0.50
2:E:113:ASN:HD22	2:E:114:LEU:N	2.00	0.50
2:E:189:ARG:HG3	4:E:255:HOH:O	2.11	0.50
1:A:122:LEU:HD23	1:A:127:PRO:HA	1.93	0.50
2:E:90:THR:OG1	2:E:91:VAL:N	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:ARG:NH2	4:A:311:HOH:O	2.45	0.50
2:B:34:GLN:NE2	4:B:227:HOH:O	2.40	0.50
1:D:103:ASN:O	1:D:152:PRO:HA	2.12	0.50
3:C:34:GLY:HA3	4:C:315:HOH:O	2.12	0.50
1:A:38:LYS:HZ1	1:A:40:GLU:CD	2.19	0.50
2:B:28:ASP:O	2:B:39:ARG:HA	2.11	0.50
2:B:57:VAL:HG11	3:C:33:LEU:CD2	2.42	0.50
2:E:39:ARG:HD2	4:E:260:HOH:O	2.11	0.49
2:E:105:LYS:H	2:E:113:ASN:HD21	1.60	0.49
2:B:162:GLU:OE2	1:D:177:HIS:ND1	2.45	0.49
1:A:76:ARG:HH22	3:C:34:GLY:C	2.20	0.49
1:A:94:ASN:CB	4:B:247:HOH:O	2.60	0.49
2:E:132:PHE:HA	2:E:136:GLN:O	2.13	0.49
1:A:91:VAL:HG11	1:A:178:TRP:HB2	1.93	0.49
2:B:166:ARG:HD3	4:B:293:HOH:O	2.13	0.49
4:B:308:HOH:O	1:D:176:LYS:CG	2.55	0.49
2:B:18:PHE:HB3	2:B:23:ARG:HH11	1.77	0.49
2:B:55:ARG:O	2:B:56:PRO:C	2.55	0.49
1:D:12:PHE:HA	2:E:9:GLU:O	2.13	0.49
2:E:14:GLU:HB2	2:E:27:LEU:HB2	1.94	0.49
2:E:109:LEU:H	2:E:109:LEU:CD1	2.24	0.49
1:D:62:ASN:HD21	3:F:29:GLU:HA	1.77	0.48
2:E:132:PHE:HB2	2:E:172:THR:HB	1.95	0.48
1:D:8:ILE:O	1:D:24:PHE:HA	2.13	0.48
1:D:11:GLU:HA	1:D:21:GLU:O	2.13	0.48
2:E:76:ASP:HA	2:E:80:ARG:HB2	1.95	0.48
1:D:143:HIS:N	1:D:143:HIS:CD2	2.80	0.48
1:D:73:MET:CE	2:E:53:LEU:CD1	2.78	0.48
1:A:24:PHE:HB3	1:A:31:ILE:HD12	1.95	0.48
1:A:78:ASN:ND2	4:A:304:HOH:O	2.39	0.48
1:D:132:VAL:HG12	1:D:151:LEU:HD13	1.95	0.48
2:E:107:GLN:HE21	2:E:107:GLN:CA	2.27	0.48
1:A:36:MET:HE2	1:A:63:ILE:CG1	2.44	0.48
1:A:40:GLU:HG2	1:A:41:THR:O	2.14	0.48
2:E:147:LEU:HG	2:E:155:PHE:CD2	2.49	0.48
1:D:142:ASP:CG	4:D:224:HOH:O	2.56	0.47
2:E:83:TYR:O	2:E:87:GLU:HB2	2.14	0.47
2:B:148:ILE:O	2:B:155:PHE:HA	2.14	0.47
2:E:115:LEU:O	2:E:160:MET:HA	2.14	0.47
2:B:172:THR:HG21	4:B:225:HOH:O	2.14	0.47
1:D:62:ASN:ND2	3:F:30:ALA:H	2.12	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:3:THR:HG22	2:E:4:ARG:N	2.29	0.47
2:E:62:ASN:HD22	2:E:68:LEU:HD21	1.79	0.47
1:A:93:THR:HA	1:A:104:VAL:O	2.15	0.47
1:A:24:PHE:CD1	1:A:32:PHE:CZ	3.02	0.47
1:A:48:PHE:CD1	2:B:89:PHE:CD1	3.02	0.47
1:A:86:PRO:HB3	1:A:169:GLY:C	2.39	0.47
1:D:50:ARG:CG	1:D:50:ARG:NH1	2.74	0.47
2:E:78:TYR:CZ	3:F:28:ASP:HB2	2.50	0.47
2:E:26:TYR:O	2:E:27:LEU:HD12	2.16	0.46
1:A:106:ILE:HB	1:A:108:PHE:CE1	2.51	0.46
2:E:88:SER:HA	2:E:92:GLN:NE2	2.30	0.46
1:D:11:GLU:HG3	2:E:11:ARG:HB3	1.98	0.46
1:D:105:LEU:HD23	1:D:105:LEU:HA	1.73	0.46
1:D:108:PHE:CD1	1:D:108:PHE:N	2.84	0.46
1:D:138:LEU:HD12	1:D:148:PHE:HE1	1.80	0.46
1:A:11:GLU:HA	1:A:21:GLU:O	2.16	0.46
2:B:56:PRO:O	2:B:60:SER:OG	2.32	0.46
2:B:57:VAL:HG13	3:C:34:GLY:H	1.81	0.46
2:E:68:LEU:C	2:E:70:GLN:N	2.73	0.46
2:E:106:THR:C	2:E:107:GLN:HE21	2.23	0.46
2:E:10:LEU:HB3	2:E:31:PHE:HB2	1.96	0.46
2:E:29:ARG:CD	2:E:36:GLU:OE2	2.57	0.46
1:A:6:VAL:HA	2:B:15:CYS:O	2.16	0.46
1:D:21:GLU:OE2	1:D:137:PHE:O	2.34	0.46
1:A:37:ALA:CB	4:A:303:HOH:O	2.44	0.46
2:B:48:ARG:NH1	2:B:48:ARG:CG	2.71	0.46
2:E:107:GLN:CA	2:E:107:GLN:NE2	2.79	0.46
2:E:31:PHE:CE2	2:E:36:GLU:HB2	2.51	0.45
1:D:9:GLN:NE2	1:D:11:GLU:OE2	2.42	0.45
1:D:41:THR:HG22	1:D:56:ALA:HB2	1.98	0.45
2:E:61:TRP:CH2	3:F:33:LEU:HD23	2.52	0.45
2:B:109:LEU:O	2:B:110:GLN:CB	2.65	0.45
1:A:5:HIS:HD2	4:B:191:HOH:O	2.00	0.45
1:A:140:ARG:O	2:B:12:LYS:NZ	2.50	0.45
1:D:92:LEU:HD22	1:D:108:PHE:CD1	2.45	0.45
1:D:38:LYS:N	4:D:253:HOH:O	2.50	0.45
1:D:103:ASN:HB3	1:D:153:PHE:CE1	2.51	0.45
2:E:68:LEU:C	2:E:70:GLN:H	2.24	0.45
2:B:31:PHE:CE2	2:B:36:GLU:HB2	2.52	0.45
2:B:112:HIS:HB2	4:B:311:HOH:O	2.16	0.45
2:E:24:VAL:HG11	2:E:79:CYS:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:44:VAL:HG12	2:E:46:GLU:H	1.82	0.45
1:A:14:LEU:CD2	1:A:116:VAL:CG2	2.94	0.45
1:A:26:PHE:HB2	1:A:31:ILE:HD11	1.99	0.45
2:E:13:SER:HA	2:E:27:LEU:O	2.17	0.45
1:A:162:ASP:HB3	1:A:175:LEU:HD22	1.99	0.45
1:A:12:PHE:CE2	1:A:21:GLU:HB3	2.52	0.44
1:A:33:HIS:CG	1:A:136:VAL:HG11	2.52	0.44
1:D:41:THR:HG21	1:D:54:PHE:HB3	1.99	0.44
1:D:105:LEU:HG	1:D:153:PHE:CE1	2.52	0.44
1:A:12:PHE:HB2	2:B:8:LEU:HD11	1.99	0.44
1:A:38:LYS:O	1:A:39:LYS:C	2.58	0.44
1:A:81:PRO:HB3	2:B:5:PRO:CB	2.47	0.44
1:A:47:GLU:HA	1:A:50:ARG:HD2	1.99	0.44
1:D:123:ARG:HB2	1:D:128:VAL:HG21	2.00	0.44
1:D:134:GLU:HA	1:D:148:PHE:O	2.18	0.44
2:E:46:GLU:OE2	2:E:48:ARG:NE	2.33	0.44
2:E:86:GLY:HA2	2:E:89:PHE:CE2	2.52	0.44
2:E:123:TYR:CG	2:E:124:PRO:HA	2.52	0.44
1:A:164:ARG:HB2	1:A:175:LEU:CD2	2.47	0.44
2:E:115:LEU:HD22	2:E:163:THR:HG21	2.00	0.44
2:E:116:VAL:HG22	2:E:160:MET:HG2	2.00	0.44
1:D:70:LEU:HB2	2:E:9:GLU:HB2	1.98	0.44
1:D:167:HIS:HB3	1:D:170:LEU:HD22	1.99	0.44
1:D:171:ASP:OD1	1:D:171:ASP:N	2.39	0.44
1:D:172:GLU:HB2	1:D:173:PRO:HD2	2.00	0.44
1:A:110:ASP:CG	1:A:111:LYS:HG3	2.42	0.44
2:B:36:GLU:O	2:B:51:THR:HG23	2.18	0.44
1:D:30:GLU:O	1:D:44:ARG:CB	2.66	0.44
1:D:69:ASN:CG	3:F:33:LEU:HG	2.43	0.44
2:B:136:GLN:CG	4:B:267:HOH:O	2.66	0.44
1:D:54:PHE:HZ	3:F:27:SER:HB3	1.80	0.44
1:D:118:ASN:HB2	1:D:166:GLU:HB2	2.00	0.44
1:A:129:THR:O	1:A:132:VAL:HB	2.18	0.43
1:D:91:VAL:H	1:D:176:LYS:NZ	2.16	0.43
2:B:101:VAL:CG1	2:B:188:TRP:HB2	2.48	0.43
1:D:54:PHE:CE1	3:F:27:SER:HB3	2.54	0.43
1:A:9:GLN:O	2:B:12:LYS:HA	2.18	0.43
1:A:72:ILE:HG22	1:A:76:ARG:HD2	2.01	0.43
2:B:78:TYR:CZ	3:C:28:ASP:HB2	2.53	0.43
2:E:39:ARG:O	2:E:39:ARG:HG3	2.18	0.43
1:A:175:LEU:O	4:A:263:HOH:O	2.22	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:117:VAL:HG22	1:D:137:PHE:CE2	2.52	0.43
1:D:176:LYS:NZ	4:D:296:HOH:O	2.51	0.43
2:B:27:LEU:HA	2:B:40:PHE:O	2.18	0.43
2:B:26:TYR:OH	3:C:28:ASP:OD2	2.23	0.43
1:D:12:PHE:CD1	1:D:12:PHE:C	2.97	0.43
2:E:147:LEU:HD12	2:E:147:LEU:HA	1.86	0.43
2:E:182:SER:HB2	2:E:183:ALA:H	1.60	0.43
1:D:6:VAL:HA	2:E:15:CYS:O	2.19	0.43
1:D:34:VAL:HG21	1:D:59:ALA:HB2	2.01	0.43
1:D:75:LYS:HG2	1:D:79:TYR:OH	2.18	0.43
1:A:101:GLU:HA	1:A:102:PRO:HD3	1.92	0.42
2:B:127:ILE:HD11	2:B:175:VAL:HG13	2.01	0.42
2:E:48:ARG:HG3	2:E:48:ARG:HH11	1.83	0.42
1:A:21:GLU:OE2	1:A:137:PHE:O	2.36	0.42
1:A:12:PHE:O	1:A:20:GLY:HA2	2.20	0.42
1:A:76:ARG:HH11	1:A:76:ARG:HD3	1.65	0.42
1:A:92:LEU:HD23	1:A:92:LEU:HA	1.84	0.42
2:B:123:TYR:CD1	2:B:124:PRO:HA	2.55	0.42
1:D:76:ARG:HB3	2:E:53:LEU:CD2	2.49	0.42
1:D:162:ASP:HA	1:D:176:LYS:O	2.19	0.42
1:A:3:GLU:OE2	1:A:6:VAL:HG21	2.15	0.42
1:A:176:LYS:HD2	1:A:176:LYS:HA	1.76	0.42
2:B:40:PHE:HB2	2:B:47:TYR:CE2	2.54	0.42
1:D:77:SER:O	1:D:78:ASN:HB2	2.18	0.42
1:A:3:GLU:OE2	1:A:6:VAL:HG23	2.18	0.42
1:A:22:PHE:CD1	1:A:22:PHE:C	2.97	0.42
2:E:57:VAL:HG11	3:F:33:LEU:CD2	2.46	0.42
1:A:29:ASP:HB3	2:B:153:TRP:CE2	2.54	0.42
1:A:85:VAL:HB	1:A:113:THR:HG22	2.02	0.42
1:D:116:VAL:O	1:D:167:HIS:HD2	2.03	0.42
1:A:105:LEU:HB2	1:A:151:LEU:HB3	2.02	0.42
2:E:24:VAL:HG21	2:E:80:ARG:HG3	2.02	0.42
1:A:122:LEU:O	1:A:161:TYR:HA	2.20	0.42
1:D:29:ASP:HB3	2:E:153:TRP:CE2	2.55	0.42
1:D:67:LYS:O	1:D:70:LEU:HB3	2.19	0.42
1:D:109:ILE:CD1	1:D:119:VAL:HG21	2.50	0.42
2:E:61:TRP:CZ2	3:F:31:LEU:HB2	2.55	0.41
2:E:115:LEU:HD12	2:E:115:LEU:HA	1.82	0.41
1:A:17:ASP:O	1:A:18:GLN:HB2	2.20	0.41
1:A:73:MET:SD	3:C:33:LEU:HD13	2.60	0.41
2:B:37:PHE:CD1	2:B:38:LEU:HB2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:177:HIS:CD2	2:B:178:PRO:HD2	2.55	0.41
2:E:83:TYR:CZ	2:E:91:VAL:HG21	2.55	0.41
1:A:38:LYS:N	4:A:303:HOH:O	2.54	0.41
2:B:15:CYS:HB3	2:B:17:PHE:CZ	2.56	0.41
1:A:128:VAL:CG1	1:A:151:LEU:HD11	2.50	0.41
2:B:97:PRO:HA	2:B:122:PHE:HB3	2.03	0.41
2:B:106:THR:HA	4:B:286:HOH:O	2.19	0.41
2:B:144:SER:C	2:B:146:GLY:H	2.28	0.41
2:B:170:VAL:HG22	2:B:189:ARG:HG2	2.01	0.41
2:E:86:GLY:HA2	2:E:89:PHE:CZ	2.55	0.41
1:A:94:ASN:HB2	4:B:247:HOH:O	2.21	0.41
1:D:164:ARG:HB2	1:D:175:LEU:CD2	2.50	0.41
1:A:108:PHE:CD2	1:A:148:PHE:HE2	2.39	0.41
1:A:140:ARG:HH21	1:A:142:ASP:CG	2.26	0.41
2:B:129:VAL:HB	2:B:159:VAL:HG21	2.03	0.41
1:D:45:LEU:HD11	2:E:153:TRP:HB2	2.02	0.41
1:D:51:PHE:O	1:D:52:ALA:HB2	2.21	0.41
2:E:24:VAL:HB	2:E:75:VAL:HG13	2.03	0.41
2:E:27:LEU:HA	2:E:40:PHE:O	2.21	0.41
1:A:87:PRO:HB3	1:A:109:ILE:HG22	2.03	0.41
1:A:177:HIS:CD2	2:E:114:LEU:HD21	2.56	0.41
1:D:154:LEU:O	1:D:155:PRO:C	2.64	0.41
1:A:84:ASN:HB3	1:A:169:GLY:CA	2.51	0.40
1:A:54:PHE:CE1	3:C:27:SER:HB2	2.57	0.40
2:E:24:VAL:HG11	2:E:79:CYS:CB	2.51	0.40
2:B:14:GLU:HB2	2:B:27:LEU:HB2	2.04	0.40
2:B:115:LEU:O	2:B:160:MET:HA	2.21	0.40
1:D:147:LYS:CE	1:D:149:HIS:HE1	2.29	0.40
2:E:3:THR:HG22	2:E:6:ARG:NH1	2.35	0.40
2:E:82:ASN:HA	4:E:280:HOH:O	2.21	0.40
1:A:91:VAL:CG1	1:A:178:TRP:HB2	2.51	0.40
1:D:89:VAL:HA	1:D:108:PHE:O	2.21	0.40
1:D:164:ARG:HD2	1:D:175:LEU:HD21	2.03	0.40
1:D:179:GLU:OE2	1:D:181:ASP:OD1	2.40	0.40
2:E:59:GLU:O	2:E:60:SER:C	2.64	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	177/182 (97%)	170 (96%)	7 (4%)	0	100	100
1	D	176/182 (97%)	169 (96%)	7 (4%)	0	100	100
2	B	186/190 (98%)	175 (94%)	11 (6%)	0	100	100
2	E	186/190 (98%)	169 (91%)	14 (8%)	3 (2%)	7	4
3	C	9/12 (75%)	9 (100%)	0	0	100	100
3	F	10/12 (83%)	10 (100%)	0	0	100	100
All	All	744/768 (97%)	702 (94%)	39 (5%)	3 (0%)	30	31

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	182	SER
2	E	58	ALA
2	E	62	ASN

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	154/166 (93%)	147 (96%)	7 (4%)	24	29
1	D	158/166 (95%)	141 (89%)	17 (11%)	6	4
2	B	152/170 (89%)	137 (90%)	15 (10%)	7	5
2	E	151/170 (89%)	137 (91%)	14 (9%)	8	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	8/9 (89%)	8 (100%)	0	100	100
3	F	9/9 (100%)	9 (100%)	0	100	100
All	All	632/690 (92%)	579 (92%)	53 (8%)	10	8

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	MET
1	A	60	LEU
1	A	95	SER
1	A	113	THR
1	A	120	THR
1	A	132	VAL
1	A	170	LEU
2	B	13	SER
2	B	21	THR
2	B	27	LEU
2	B	38	LEU
2	B	48	ARG
2	B	50	VAL
2	B	60	SER
2	B	69	GLU
2	B	105	LYS
2	B	109	LEU
2	B	114	LEU
2	B	172	THR
2	B	184	LEU
2	B	185	THR
2	B	187	GLU
1	D	11	GLU
1	D	39	LYS
1	D	41	THR
1	D	60	LEU
1	D	77	SER
1	D	79	TYR
1	D	80	THR
1	D	92	LEU
1	D	97	VAL
1	D	117	VAL
1	D	141	GLU
1	D	143	HIS

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Mol	Chain	Res	Type
1	D	158	GLU
1	D	170	LEU
1	D	171	ASP
1	D	176	LYS
1	D	181	ASP
2	E	13	SER
2	E	21	THR
2	E	35	GLU
2	E	38	LEU
2	E	39	ARG
2	E	63	SER
2	E	106	THR
2	E	107	GLN
2	E	114	LEU
2	E	180	VAL
2	E	181	THR
2	E	182	SER
2	E	184	LEU
2	E	186	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	62	ASN
1	A	149	HIS
2	B	32	HIS
2	B	64	GLN
2	B	70	GLN
2	B	113	ASN
2	B	150	ASN
1	D	15	ASN
1	D	62	ASN
1	D	149	HIS
1	D	167	HIS
2	E	32	HIS
2	E	62	ASN
2	E	81	HIS
2	E	92	GLN
2	E	107	GLN
2	E	113	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	180/182 (98%)	-0.08	2 (1%) 78 79	25, 39, 54, 63	0
1	D	180/182 (98%)	0.12	6 (3%) 49 49	25, 41, 62, 71	1 (0%)
2	B	188/190 (98%)	0.01	3 (1%) 70 71	23, 37, 59, 69	0
2	E	188/190 (98%)	0.43	15 (7%) 18 16	28, 47, 75, 81	0
3	C	11/12 (91%)	0.81	1 (9%) 15 14	39, 40, 60, 69	0
3	F	12/12 (100%)	0.83	1 (8%) 17 16	45, 54, 72, 81	1 (8%)
All	All	759/768 (98%)	0.15	28 (3%) 45 45	23, 41, 65, 81	2 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	125	GLY	4.3
2	E	133	ARG	4.2
1	D	182	ALA	4.0
3	C	34	GLY	3.8
2	E	109	LEU	3.7
2	E	190	ALA	3.6
2	E	134	ASN	3.4
1	A	142	ASP	2.9
1	D	142	ASP	2.8
2	E	166	ARG	2.7
2	E	111	HIS	2.7
1	A	108	PHE	2.7
3	F	34	GLY	2.7
1	D	37	ALA	2.7
2	E	188	TRP	2.7
2	E	70	GLN	2.6
2	E	110	GLN	2.5
2	B	66	ASP	2.4
2	E	53	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
2	E	66	ASP	2.3
1	D	143	HIS	2.3
2	B	107	GLN	2.3
2	B	106	THR	2.2
1	D	180	PHE	2.1
2	E	108	PRO	2.1
1	D	75	LYS	2.0
2	E	67	LEU	2.0
2	E	170	VAL	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.