



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:01 PM UTC

PDB ID : 1KLU / pdb_00001klu
Title : Crystal structure of HLA-DR1/TPI(23-37) complexed with staphylococcal enterotoxin C3 variant 3B2 (SEC3-3B2)
Authors : Sundberg, E.J.; Sawicki, M.W.; Andersen, P.S.; Sidney, J.; Sette, A.; Mariuzza, R.A.
Deposited on : 2001-12-12
Resolution : 1.93 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

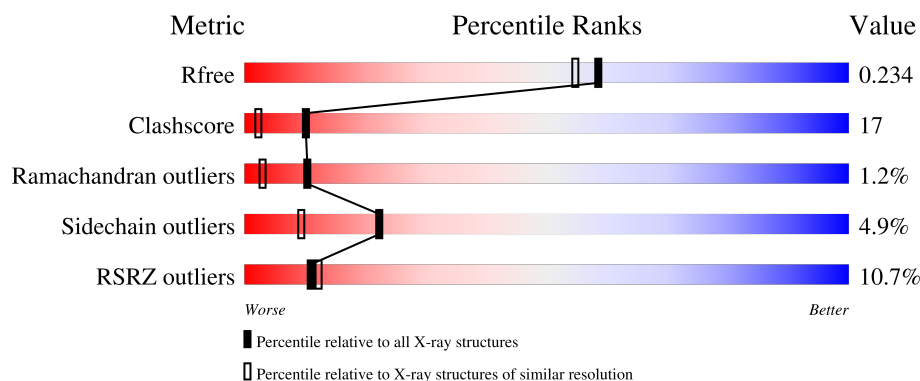
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	179	7% 74% 24% .
2	B	190	15% 65% 27% 7% .
3	C	15	20% 67% 33%
4	D	239	10% 69% 23% 5% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5415 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	179	Total	C	N	O	S	0	0	0
			1470	952	239	274	5			

- Molecule 2 is a protein called HLA CLASS II HISTOCOMPATIBILITY ANTIGEN, DR-1 BETA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	190	Total	C	N	O	S	0	0	0
			1557	979	279	293	6			

- Molecule 3 is a protein called Triosephosphate isomerase peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	15	Total	C	N	O	0	0	0
			103	64	17	22			

- Molecule 4 is a protein called ENTEROTOXIN TYPE C-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	230	Total	C	N	O	S	0	0	0
			1880	1194	306	370	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	43	SER	LYS	engineered mutation	UNP P0A0L5
D	45	PHE	LEU	engineered mutation	UNP P0A0L5
D	46	LYS	ALA	engineered mutation	UNP P0A0L5
D	47	TRP	HIS	engineered mutation	UNP P0A0L5

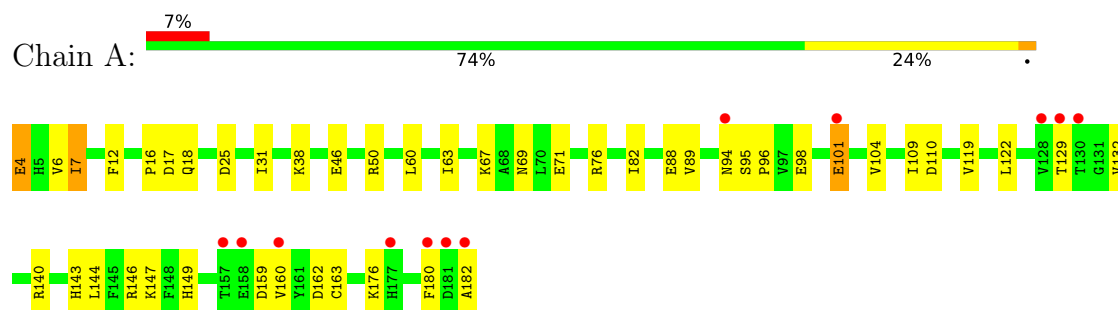
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	123	Total 123	O 123	0	0
5	B	106	Total 106	O 106	0	0
5	C	25	Total 25	O 25	0	0
5	D	151	Total 151	O 151	0	0

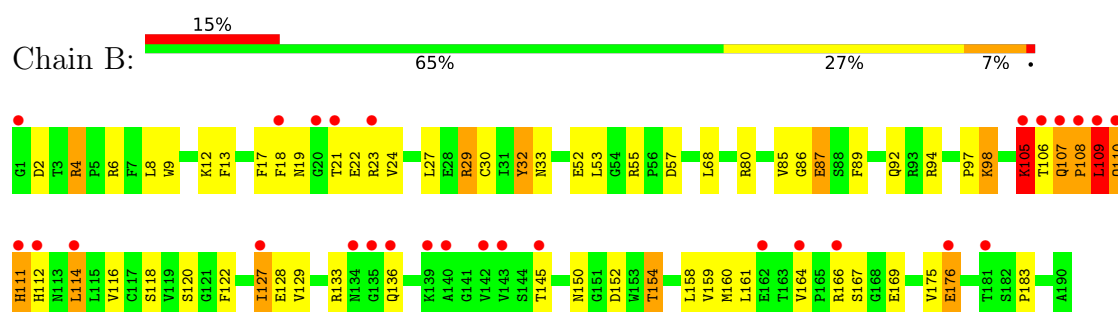
3 Residue-property plots [i](#)

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

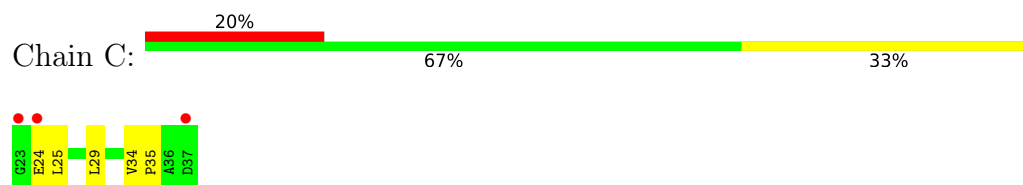
- Molecule 1: HLA CLASS II HISTOCOMPATIBILITY ANTIGEN, DR ALPHA CHAIN



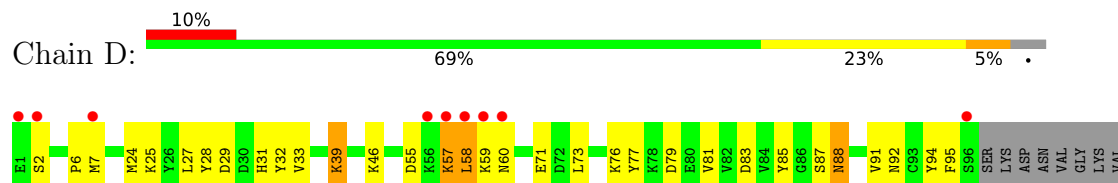
- Molecule 2: HLA CLASS II HISTOCOMPATIBILITY ANTIGEN, DR-1 BETA CHAIN

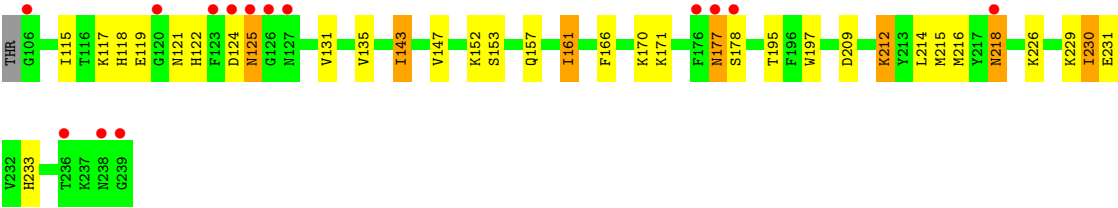


- Molecule 3: Triosephosphate isomerase peptide



- Molecule 4: ENTEROTOXIN TYPE C-3





4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	171.30Å 171.30Å 121.02Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	14.95 – 1.93 14.95 – 1.93	Depositor EDS
% Data completeness (in resolution range)	84.3 (14.95-1.93) 95.5 (14.95-1.93)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 1.94Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.208 , 0.220 0.223 , 0.234	Depositor DCC
R_{free} test set	5019 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	19.8	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.025 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5415	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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	0.39	0/1515	0.91	6/2065 (0.3%)
2	B	0.43	0/1597	1.02	8/2168 (0.4%)
3	C	0.59	0/103	1.15	1/138 (0.7%)
4	D	0.35	0/1922	0.88	5/2587 (0.2%)
All	All	0.39	0/5137	0.94	20/6958 (0.3%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	109	LEU	N-CA-C	16.96	136.98	110.70
2	B	164	VAL	CA-C-N	6.54	126.57	119.90
2	B	164	VAL	C-N-CA	6.54	126.57	119.90
2	B	105	LYS	N-CA-C	-6.34	105.83	112.93
4	D	7	MET	N-CA-C	-5.96	102.59	110.40
1	A	31	ILE	N-CA-C	-5.88	105.98	111.45
1	A	110	ASP	N-CA-C	5.71	118.09	109.41
1	A	163	CYS	N-CA-C	-5.53	98.74	108.20
4	D	147	VAL	N-CA-C	-5.53	100.62	109.20
3	C	25	LEU	N-CA-C	5.35	117.44	109.25
4	D	115	ILE	N-CA-C	5.29	115.58	108.17
1	A	159	ASP	N-CA-C	5.27	117.52	110.35
2	B	120	SER	N-CA-C	5.22	116.93	109.14
1	A	88	GLU	N-CA-C	-5.15	100.34	108.73
4	D	24	MET	N-CA-C	-5.13	105.31	112.45
4	D	81	VAL	N-CA-C	-5.11	100.97	108.54
1	A	144	LEU	N-CA-C	-5.06	102.02	109.86
2	B	109	LEU	N-CA-CB	-5.04	102.52	110.49
2	B	107	GLN	CA-C-N	-5.01	113.58	119.84
2	B	107	GLN	C-N-CA	-5.01	113.58	119.84

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	1470	0	1406	42	0
2	B	1557	0	1488	84	0
3	C	103	0	105	5	0
4	D	1880	0	1807	59	0
5	A	123	0	0	1	0
5	B	106	0	0	0	0
5	C	25	0	0	0	0
5	D	151	0	0	1	0
All	All	5415	0	4806	169	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 17.

All (169) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:150:ASN:HD22	2:B:154:THR:HG22	1.12	1.14
2:B:127:ILE:HD11	2:B:175:VAL:HG13	1.39	1.04
2:B:116:VAL:HG13	2:B:160:MET:HE2	1.42	1.02
2:B:127:ILE:HD13	2:B:128:GLU:N	1.84	0.91
2:B:52:GLU:HA	2:B:55:ARG:HD2	1.54	0.90
2:B:109:LEU:HD12	2:B:112:HIS:CD2	2.07	0.89
4:D:117:LYS:HE2	4:D:119:GLU:HB3	1.53	0.87
4:D:121:ASN:HD21	4:D:153:SER:H	1.20	0.86
2:B:150:ASN:HD22	2:B:154:THR:CG2	1.91	0.83
1:A:76:ARG:HD3	5:A:305:HOH:O	1.77	0.83
2:B:158:LEU:HB3	2:B:160:MET:HE1	1.60	0.82
2:B:114:LEU:HD21	2:B:160:MET:HB3	1.62	0.81
1:A:7:ILE:HD13	2:B:17:PHE:CE2	2.18	0.79
2:B:109:LEU:HD12	2:B:112:HIS:CG	2.18	0.78
2:B:150:ASN:ND2	2:B:154:THR:HG22	1.97	0.77
4:D:215:MET:O	4:D:218:ASN:HB2	1.85	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:39:LYS:HZ1	4:D:79:ASP:C	1.93	0.76
2:B:110:GLN:O	2:B:112:HIS:N	2.18	0.75
4:D:157:GLN:O	4:D:161:ILE:HD13	1.87	0.74
1:A:16:PRO:HD2	2:B:6:ARG:HD3	1.70	0.73
2:B:159:VAL:C	2:B:160:MET:HE3	2.14	0.73
2:B:105:LYS:HE3	2:B:105:LYS:H	1.52	0.73
4:D:39:LYS:NZ	4:D:79:ASP:HA	2.04	0.73
4:D:166:PHE:CZ	4:D:170:LYS:HD2	2.23	0.72
2:B:109:LEU:O	2:B:112:HIS:CD2	2.42	0.72
4:D:229:LYS:C	4:D:230:ILE:HD12	2.14	0.72
1:A:122:LEU:HB2	1:A:162:ASP:HB2	1.73	0.70
1:A:94:ASN:ND2	1:A:104:VAL:HB	2.07	0.70
2:B:8:LEU:O	2:B:32:TYR:O	2.09	0.69
2:B:21:THR:O	2:B:80:ARG:NH1	2.25	0.69
4:D:88:ASN:H	4:D:88:ASN:HD22	1.40	0.68
1:A:16:PRO:HD2	2:B:6:ARG:HH11	1.58	0.67
1:A:147:LYS:HE3	1:A:149:HIS:HE1	1.59	0.67
4:D:131:VAL:HG12	4:D:230:ILE:HD13	1.78	0.66
2:B:106:THR:HG22	2:B:108:PRO:HD3	1.78	0.65
1:A:7:ILE:HD13	2:B:17:PHE:HE2	1.60	0.65
2:B:94:ARG:HH11	2:B:94:ARG:HG3	1.61	0.64
4:D:39:LYS:NZ	4:D:79:ASP:C	2.54	0.64
4:D:28:TYR:HE2	4:D:161:ILE:HD12	1.62	0.63
2:B:98:LYS:HB2	2:B:98:LYS:NZ	2.13	0.63
2:B:176:GLU:OE2	2:B:183:PRO:HB3	1.99	0.62
1:A:6:VAL:C	1:A:7:ILE:HD12	2.24	0.62
4:D:39:LYS:HZ1	4:D:79:ASP:CA	2.12	0.62
2:B:109:LEU:O	2:B:110:GLN:O	2.17	0.62
2:B:116:VAL:HG13	2:B:160:MET:CE	2.23	0.62
2:B:110:GLN:O	2:B:111:HIS:C	2.41	0.62
4:D:230:ILE:HD12	4:D:230:ILE:N	2.15	0.61
2:B:85:VAL:HG13	3:C:24:GLU:HB2	1.82	0.60
2:B:114:LEU:CD2	2:B:160:MET:HB3	2.29	0.60
4:D:143:ILE:HD12	4:D:171:LYS:HE3	1.82	0.60
4:D:2:SER:CB	4:D:195:THR:H	2.16	0.59
2:B:133:ARG:O	2:B:136:GLN:HG2	2.03	0.58
1:A:94:ASN:HD22	1:A:104:VAL:HB	1.67	0.58
2:B:152:ASP:OD1	2:B:154:THR:HB	2.04	0.58
4:D:121:ASN:ND2	4:D:153:SER:H	1.97	0.58
2:B:105:LYS:H	2:B:105:LYS:CE	2.16	0.57
2:B:160:MET:HE3	2:B:160:MET:N	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:97:PRO:HB3	2:B:122:PHE:HB3	1.86	0.57
2:B:109:LEU:CD1	2:B:112:HIS:CE1	2.88	0.57
2:B:109:LEU:HD11	2:B:112:HIS:CE1	2.39	0.56
2:B:127:ILE:HD13	2:B:128:GLU:H	1.69	0.56
4:D:88:ASN:H	4:D:88:ASN:ND2	2.02	0.56
4:D:2:SER:HB2	4:D:195:THR:H	1.68	0.56
2:B:94:ARG:HG3	2:B:94:ARG:NH1	2.20	0.56
2:B:114:LEU:HD23	2:B:161:LEU:C	2.30	0.56
2:B:116:VAL:CG1	2:B:160:MET:HE2	2.28	0.56
1:A:89:VAL:HG12	1:A:176:LYS:HG3	1.86	0.56
4:D:131:VAL:CG1	4:D:230:ILE:HD13	2.36	0.56
1:A:7:ILE:HD12	1:A:7:ILE:N	2.20	0.56
2:B:145:THR:HG22	2:B:158:LEU:H	1.71	0.55
4:D:57:LYS:HB3	4:D:58:LEU:HD22	1.89	0.55
4:D:135:VAL:HB	4:D:143:ILE:CD1	2.37	0.55
2:B:2:ASP:OD1	2:B:4:ARG:HD3	2.06	0.55
4:D:122:HIS:O	4:D:152:LYS:HE3	2.07	0.54
2:B:109:LEU:C	2:B:110:GLN:O	2.51	0.53
4:D:58:LEU:HD22	4:D:58:LEU:N	2.23	0.53
1:A:60:LEU:O	1:A:63:ILE:HG22	2.08	0.53
2:B:24:VAL:HG23	2:B:80:ARG:NH1	2.24	0.53
1:A:38:LYS:O	4:D:212:LYS:HE2	2.09	0.52
2:B:109:LEU:O	2:B:112:HIS:HD2	1.91	0.52
2:B:108:PRO:HG2	2:B:109:LEU:H	1.75	0.52
2:B:109:LEU:CD1	2:B:112:HIS:CG	2.91	0.52
2:B:111:HIS:O	2:B:112:HIS:CD2	2.63	0.52
1:A:143:HIS:HD2	2:B:12:LYS:NZ	2.08	0.51
2:B:57:ASP:OD1	3:C:35:PRO:HD2	2.10	0.51
4:D:39:LYS:HZ2	4:D:79:ASP:HA	1.73	0.51
2:B:114:LEU:HD23	2:B:161:LEU:O	2.11	0.51
2:B:166:ARG:O	2:B:169:GLU:HG3	2.11	0.51
4:D:39:LYS:HZ1	4:D:79:ASP:HA	1.70	0.51
4:D:57:LYS:CB	4:D:58:LEU:HD22	2.41	0.50
4:D:87:SER:H	4:D:157:GLN:NE2	2.10	0.50
4:D:46:LYS:HB2	4:D:71:GLU:HG3	1.94	0.50
2:B:86:GLY:HA2	2:B:89:PHE:CE1	2.47	0.49
1:A:67:LYS:O	1:A:71:GLU:HG2	2.13	0.49
4:D:161:ILE:CD1	4:D:161:ILE:N	2.76	0.49
4:D:83:ASP:OD1	4:D:118:HIS:HD2	1.95	0.49
1:A:16:PRO:HD2	2:B:6:ARG:CD	2.39	0.49
2:B:106:THR:HG22	2:B:108:PRO:CD	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:127:ILE:HD12	2:B:129:VAL:HG23	1.93	0.48
4:D:76:LYS:HD2	4:D:77:TYR:CE1	2.48	0.48
4:D:226:LYS:HG2	5:D:262:HOH:O	2.11	0.48
2:B:127:ILE:HD12	2:B:129:VAL:CG2	2.43	0.48
1:A:82:ILE:HG13	2:B:33:ASN:HB3	1.96	0.48
2:B:107:GLN:O	2:B:108:PRO:C	2.53	0.47
2:B:98:LYS:HB2	2:B:98:LYS:HZ2	1.79	0.47
4:D:231:GLU:HB3	4:D:233:HIS:CE1	2.49	0.47
1:A:160:VAL:O	1:A:160:VAL:HG23	2.13	0.47
1:A:16:PRO:HD2	2:B:6:ARG:NH1	2.27	0.47
1:A:140:ARG:HG3	1:A:146:ARG:HG3	1.98	0.46
1:A:16:PRO:HG2	2:B:6:ARG:NH1	2.31	0.46
1:A:89:VAL:CG1	1:A:176:LYS:HG3	2.45	0.46
2:B:176:GLU:CD	2:B:183:PRO:HB3	2.41	0.46
4:D:31:HIS:O	4:D:32:TYR:HB3	2.16	0.45
1:A:12:PHE:CD1	1:A:12:PHE:C	2.94	0.45
1:A:7:ILE:N	1:A:7:ILE:CD1	2.79	0.45
4:D:28:TYR:CE2	4:D:161:ILE:HD12	2.48	0.45
1:A:46:GLU:OE2	1:A:50:ARG:NE	2.44	0.45
1:A:69:ASN:OD1	3:C:34:VAL:HG22	2.17	0.45
1:A:129:THR:O	1:A:132:VAL:HG22	2.17	0.45
2:B:27:LEU:HG	2:B:29:ARG:HD2	1.99	0.45
4:D:209:ASP:CG	4:D:212:LYS:HB2	2.43	0.44
4:D:55:ASP:O	4:D:59:LYS:HE3	2.17	0.44
4:D:39:LYS:NZ	4:D:79:ASP:CA	2.70	0.44
1:A:17:ASP:O	1:A:18:GLN:HB2	2.17	0.44
1:A:16:PRO:CD	2:B:6:ARG:HH11	2.29	0.44
4:D:33:VAL:O	4:D:85:TYR:HA	2.17	0.44
4:D:121:ASN:HD21	4:D:153:SER:N	2.00	0.43
4:D:55:ASP:HB3	4:D:60:ASN:H	1.83	0.43
4:D:230:ILE:N	4:D:230:ILE:CD1	2.80	0.43
4:D:25:LYS:HE2	4:D:29:ASP:OD1	2.19	0.43
1:A:95:SER:HB2	1:A:96:PRO:HD2	2.00	0.43
4:D:87:SER:H	4:D:157:GLN:HE21	1.65	0.43
4:D:125:ASN:HD22	4:D:125:ASN:HA	1.53	0.43
2:B:13:PHE:CD2	3:C:29:LEU:HD23	2.54	0.43
1:A:4:GLU:OE1	2:B:19:ASN:HA	2.20	0.42
1:A:109:ILE:CD1	1:A:119:VAL:HG21	2.48	0.42
2:B:2:ASP:OD1	2:B:4:ARG:CD	2.67	0.42
4:D:216:MET:C	4:D:218:ASN:H	2.27	0.42
1:A:180:PHE:CE1	1:A:182:ALA:HB3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:19:ASN:HB3	2:B:22:GLU:HB3	2.02	0.42
1:A:46:GLU:CD	1:A:46:GLU:C	2.87	0.42
4:D:58:LEU:HD22	4:D:58:LEU:H	1.84	0.42
4:D:166:PHE:CE1	4:D:170:LYS:HD2	2.55	0.42
1:A:7:ILE:HD13	2:B:17:PHE:CD2	2.54	0.42
4:D:27:LEU:HD22	4:D:214:LEU:HD11	2.02	0.42
2:B:24:VAL:CG2	2:B:80:ARG:NH1	2.83	0.42
1:A:76:ARG:NH2	2:B:57:ASP:OD2	2.43	0.42
1:A:101:GLU:OE1	1:A:101:GLU:O	2.36	0.42
2:B:127:ILE:HD13	2:B:127:ILE:C	2.43	0.42
4:D:212:LYS:HD3	4:D:212:LYS:HA	1.64	0.42
4:D:161:ILE:HD13	4:D:161:ILE:H	1.85	0.42
4:D:231:GLU:OE1	4:D:233:HIS:HE1	2.03	0.42
2:B:118:SER:HA	2:B:158:LEU:HD23	2.02	0.42
4:D:91:VAL:O	4:D:92:ASN:HB2	2.20	0.42
4:D:94:TYR:O	4:D:95:PHE:HB3	2.20	0.42
2:B:9:TRP:CH2	2:B:30:CYS:HB3	2.54	0.41
1:A:76:ARG:HH22	2:B:57:ASP:CG	2.26	0.41
2:B:109:LEU:CD1	2:B:112:HIS:CD2	2.92	0.41
4:D:177:ASN:O	4:D:178:SER:HB2	2.19	0.41
4:D:6:PRO:HB3	4:D:197:TRP:CZ2	2.56	0.41
2:B:87:GLU:HG3	2:B:92:GLN:NE2	2.35	0.41
2:B:166:ARG:O	2:B:167:SER:C	2.62	0.41
4:D:135:VAL:HB	4:D:143:ILE:HD13	2.02	0.41
1:A:143:HIS:HD2	2:B:12:LYS:HZ1	1.68	0.41
1:A:17:ASP:OD1	2:B:6:ARG:HD2	2.21	0.41
1:A:60:LEU:N	1:A:60:LEU:HD22	2.36	0.40
2:B:18:PHE:CG	2:B:23:ARG:NH2	2.89	0.40
2:B:109:LEU:HD11	2:B:112:HIS:ND1	2.36	0.40
3:C:29:LEU:HD12	3:C:29:LEU:HA	1.89	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	177/179 (99%)	175 (99%)	2 (1%)	0	100	100
2	B	188/190 (99%)	174 (93%)	10 (5%)	4 (2%)	5	1
3	C	13/15 (87%)	13 (100%)	0	0	100	100
4	D	226/239 (95%)	216 (96%)	7 (3%)	3 (1%)	9	3
All	All	604/623 (97%)	578 (96%)	19 (3%)	7 (1%)	10	3

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	110	GLN
2	B	111	HIS
4	D	57	LYS
4	D	125	ASN
2	B	32	TYR
4	D	124	ASP
2	B	108	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	163/163 (100%)	158 (97%)	5 (3%)	35	21
2	B	171/171 (100%)	159 (93%)	12 (7%)	14	3
3	C	10/10 (100%)	10 (100%)	0	100	100
4	D	211/220 (96%)	201 (95%)	10 (5%)	23	10
All	All	555/564 (98%)	528 (95%)	27 (5%)	22	9

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU

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Mol	Chain	Res	Type
1	A	7	ILE
1	A	25	ASP
1	A	98	GLU
1	A	101	GLU
2	B	4	ARG
2	B	29	ARG
2	B	53	LEU
2	B	68	LEU
2	B	87	GLU
2	B	98	LYS
2	B	105	LYS
2	B	109	LEU
2	B	114	LEU
2	B	127	ILE
2	B	154	THR
2	B	176	GLU
4	D	39	LYS
4	D	58	LEU
4	D	73	LEU
4	D	88	ASN
4	D	143	ILE
4	D	161	ILE
4	D	177	ASN
4	D	212	LYS
4	D	218	ASN
4	D	230	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	HIS
1	A	149	HIS
2	B	70	GLN
2	B	92	GLN
2	B	112	HIS
2	B	113	ASN
2	B	150	ASN
2	B	156	GLN
4	D	52	ASN
4	D	60	ASN
4	D	88	ASN
4	D	92	ASN

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Mol	Chain	Res	Type
4	D	118	HIS
4	D	121	ASN
4	D	125	ASN
4	D	129	GLN
4	D	130	ASN
4	D	148	GLN
4	D	157	GLN
4	D	194	ASN
4	D	218	ASN
4	D	233	HIS

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	179/179 (100%)	0.22	12 (6%) 24 27	11, 20, 48, 65	0
2	B	190/190 (100%)	0.62	28 (14%) 6 6	11, 25, 57, 76	0
3	C	15/15 (100%)	0.62	3 (20%) 3 3	14, 19, 43, 46	0
4	D	230/239 (96%)	0.37	23 (10%) 12 14	13, 23, 50, 67	0
All	All	614/623 (98%)	0.41	66 (10%) 11 12	11, 23, 53, 76	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	106	THR	5.7
1	A	182	ALA	5.6
4	D	239	GLY	5.6
4	D	57	LYS	5.4
2	B	109	LEU	5.3
2	B	108	PRO	5.2
4	D	177	ASN	5.1
1	A	101	GLU	5.1
4	D	96	SER	5.0
4	D	126	GLY	5.0
2	B	111	HIS	4.7
2	B	112	HIS	4.6
4	D	56	LYS	4.5
2	B	21	THR	4.2
1	A	181	ASP	4.2
3	C	23	GLY	4.2
4	D	1	GLU	4.1
2	B	1	GLY	4.1
2	B	110	GLN	3.9
2	B	145	THR	3.8
4	D	178	SER	3.8

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Mol	Chain	Res	Type	RSRZ
2	B	105	LYS	3.8
4	D	125	ASN	3.6
2	B	136	GLN	3.6
4	D	58	LEU	3.5
1	A	160	VAL	3.5
1	A	157	THR	3.4
1	A	128	VAL	3.3
4	D	123	PHE	3.3
4	D	106	GLY	3.2
4	D	176	PHE	3.1
1	A	180	PHE	3.0
3	C	37	ASP	3.0
4	D	127	ASN	3.0
1	A	158	GLU	3.0
4	D	238	ASN	3.0
4	D	120	GLY	2.9
2	B	107	GLN	2.9
2	B	134	ASN	2.8
4	D	124	ASP	2.8
2	B	140	ALA	2.8
4	D	59	LYS	2.7
4	D	218	ASN	2.7
1	A	177	HIS	2.7
2	B	20	GLY	2.7
1	A	94	ASN	2.6
1	A	130	THR	2.6
4	D	60	ASN	2.5
2	B	142	VAL	2.5
2	B	164	VAL	2.5
2	B	143	VAL	2.4
2	B	18	PHE	2.4
1	A	129	THR	2.4
4	D	236	THR	2.4
4	D	7	MET	2.4
2	B	139	LYS	2.3
2	B	162	GLU	2.3
2	B	166	ARG	2.2
2	B	114	LEU	2.2
2	B	127	ILE	2.2
3	C	24	GLU	2.2
4	D	2	SER	2.1
2	B	135	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	181	THR	2.1
2	B	176	GLU	2.1
2	B	23	ARG	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.