



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:21 PM UTC

PDB ID : 1KLG / pdb_00001klg
Title : Crystal structure of HLA-DR1/TPI(23-37, Thr28->Ile mutant) 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 : 2001-12-11
Resolution : 2.40 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

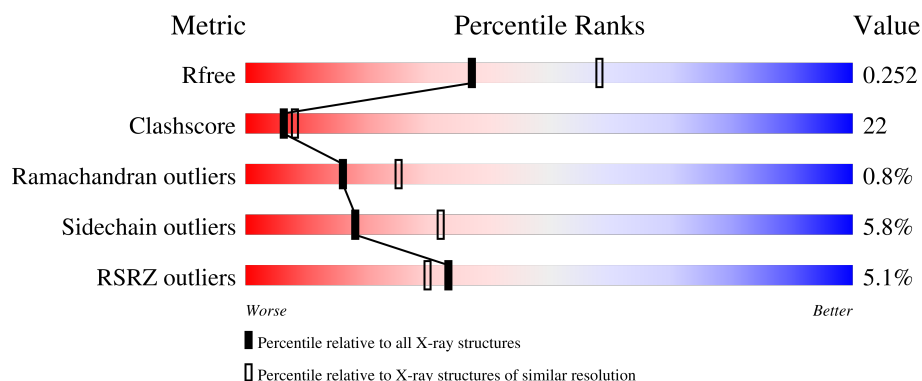
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	177	 2% 61% 36% .
2	B	190	 5% 54% 38% 6% ..
3	C	15	 13% 53% 27% 20%
4	D	239	 7% 64% 29% . .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5260 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	177	Total	C	N	O	S	0	0	0
			1457	945	237	270	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	187	Total	C	N	O	S	0	0	0
			1533	963	275	289	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
			104	66	17	21			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	231	Total	C	N	O	S	0	0	0
			1891	1201	308	372	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

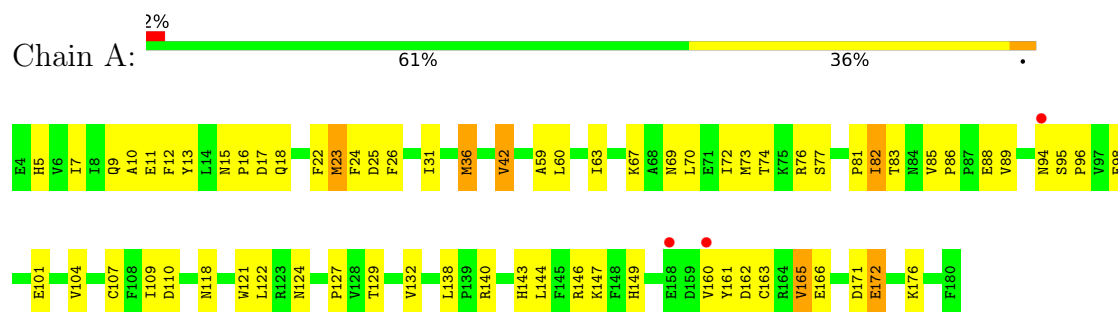
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	91	Total 91	O 91	0	0
5	B	67	Total 67	O 67	0	0
5	C	15	Total 15	O 15	0	0
5	D	102	Total 102	O 102	0	0

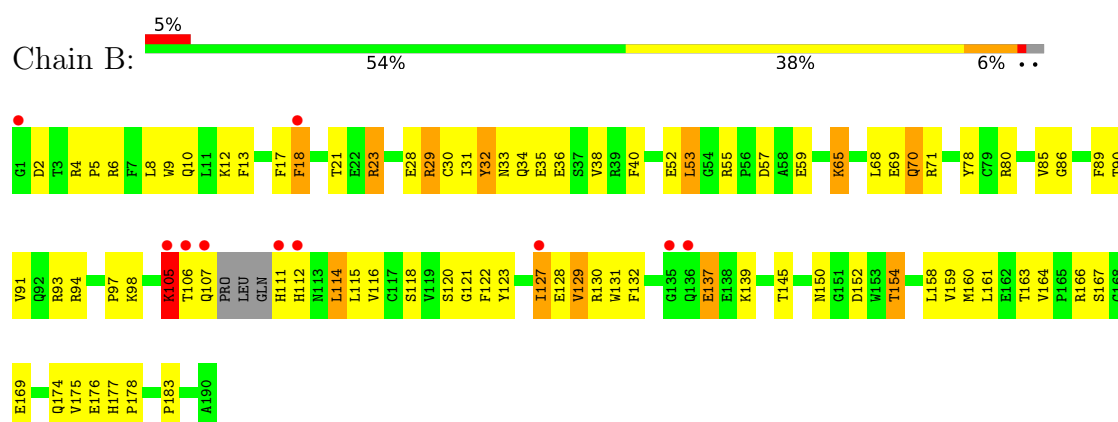
3 Residue-property plots [i](#)

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

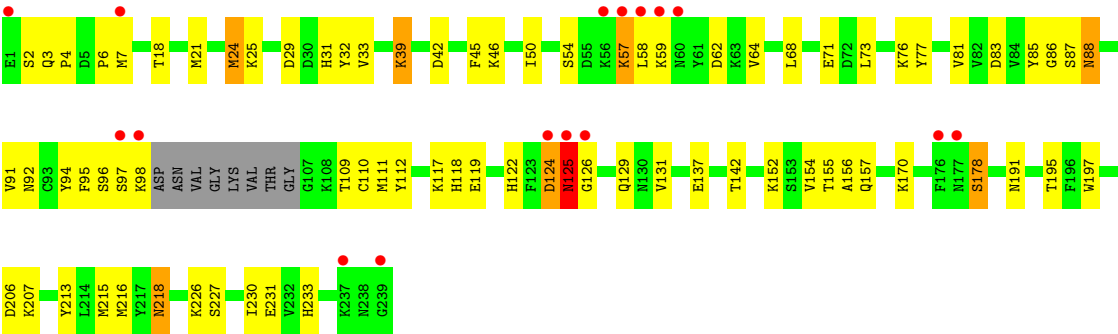


- Molecule 3: Triosephosphate isomerase peptide



- 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	172.99Å 172.99Å 121.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	14.96 – 2.40 14.96 – 2.40	Depositor EDS
% Data completeness (in resolution range)	80.5 (14.96-2.40) 92.9 (14.96-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.39Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.206 , 0.246 0.216 , 0.252	Depositor DCC
R_{free} test set	2658 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	27.7	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.033 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5260	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% 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	0.49	0/1502	1.07	12/2047 (0.6%)
2	B	0.42	0/1571	0.93	4/2130 (0.2%)
3	C	1.98	2/104 (1.9%)	1.79	4/139 (2.9%)
4	D	0.39	0/1933	0.92	5/2601 (0.2%)
All	All	0.51	2/5110 (0.0%)	0.99	25/6917 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	27	GLY	C-N	13.90	1.55	1.33
3	C	28	ILE	C-N	-13.21	1.16	1.33

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	28	ILE	O-C-N	-13.09	106.66	122.36
3	C	28	ILE	CA-C-N	7.95	133.96	123.00
3	C	28	ILE	C-N-CA	7.95	133.96	123.00
1	A	172	GLU	N-CA-C	7.08	114.59	108.22
2	B	105	LYS	N-CA-C	-6.53	101.60	110.68
1	A	144	LEU	N-CA-C	-6.51	100.33	110.17
2	B	120	SER	N-CA-C	6.31	118.54	109.14
4	D	7	MET	N-CA-C	-6.27	102.24	110.39
1	A	77	SER	N-CA-C	-6.19	105.31	112.92
4	D	81	VAL	N-CA-C	-6.17	99.69	108.58
1	A	85	VAL	CA-C-N	6.17	124.16	119.66
1	A	85	VAL	C-N-CA	6.17	124.16	119.66
2	B	164	VAL	CA-C-N	6.01	126.32	119.83
2	B	164	VAL	C-N-CA	6.01	126.32	119.83
1	A	88	GLU	N-CA-C	-5.76	99.33	108.73
4	D	156	ALA	N-CA-C	-5.70	105.07	111.28
1	A	163	CYS	N-CA-C	-5.68	98.48	108.20
1	A	165	VAL	N-CA-C	5.43	115.71	108.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	25	LEU	N-CA-C	5.41	117.53	109.25
4	D	126	GLY	N-CA-C	-5.27	106.79	114.90
1	A	15	ASN	N-CA-C	-5.21	102.94	110.40
1	A	138	LEU	CA-C-N	5.18	125.16	120.03
1	A	138	LEU	C-N-CA	5.18	125.16	120.03
1	A	86	PRO	N-CA-C	5.15	115.42	110.47
4	D	24	MET	N-CA-C	-5.11	105.35	112.45

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	1457	0	1397	71	0
2	B	1533	0	1461	105	0
3	C	104	0	108	13	0
4	D	1891	0	1822	58	0
5	A	91	0	0	3	0
5	B	67	0	0	1	0
5	C	15	0	0	1	0
5	D	102	0	0	1	0
All	All	5260	0	4788	217	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 22.

All (217) 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:127:ILE:HD11	2:B:175:VAL:HG13	1.19	1.15
1:A:82:ILE:HD13	2:B:33:ASN:HB3	1.17	1.12
2:B:150:ASN:HD22	2:B:154:THR:HG22	1.19	1.05
2:B:127:ILE:HD13	2:B:128:GLU:N	1.79	0.96
2:B:116:VAL:HG13	2:B:160:MET:HE2	1.49	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:158:LEU:HB3	2:B:160:MET:HE1	1.51	0.93
2:B:129:VAL:HG21	2:B:159:VAL:HG21	1.49	0.92
1:A:10:ALA:HB3	1:A:23:MET:HE2	1.51	0.92
1:A:82:ILE:HD12	1:A:83:THR:H	1.35	0.91
2:B:127:ILE:HD11	2:B:175:VAL:CG1	2.03	0.89
4:D:46:LYS:HB3	4:D:71:GLU:HG3	1.54	0.88
2:B:114:LEU:HD21	2:B:160:MET:HB3	1.58	0.84
1:A:22:PHE:CZ	3:C:28:ILE:HD11	2.16	0.80
2:B:150:ASN:ND2	2:B:154:THR:HG22	1.95	0.80
2:B:105:LYS:HE3	2:B:105:LYS:H	1.46	0.79
4:D:88:ASN:HD22	4:D:88:ASN:H	1.29	0.78
1:A:82:ILE:CD1	2:B:33:ASN:HB3	2.08	0.78
2:B:21:THR:O	2:B:80:ARG:NH1	2.18	0.76
4:D:215:MET:O	4:D:218:ASN:HB2	1.86	0.75
1:A:94:ASN:HD22	1:A:104:VAL:HB	1.52	0.74
2:B:31:ILE:CD1	2:B:36:GLU:HA	2.17	0.74
2:B:127:ILE:HD13	2:B:128:GLU:H	1.50	0.74
2:B:129:VAL:CG2	2:B:159:VAL:HG21	2.18	0.73
2:B:85:VAL:HG13	3:C:24:GLU:HB2	1.72	0.72
4:D:87:SER:H	4:D:157:GLN:HE21	1.35	0.72
1:A:7:ILE:HD13	2:B:17:PHE:CE2	2.24	0.72
2:B:31:ILE:HD13	2:B:36:GLU:HA	1.72	0.71
1:A:73:MET:HE1	2:B:53:LEU:HD22	1.74	0.70
4:D:64:VAL:HG22	4:D:109:THR:HG22	1.72	0.69
4:D:117:LYS:HE2	4:D:119:GLU:HB3	1.73	0.69
2:B:152:ASP:OD1	2:B:154:THR:HB	1.92	0.69
2:B:2:ASP:CG	2:B:6:ARG:HH22	2.01	0.69
1:A:109:ILE:HD12	1:A:165:VAL:HG21	1.74	0.69
2:B:8:LEU:O	2:B:32:TYR:O	2.11	0.68
1:A:129:THR:O	1:A:132:VAL:HG22	1.95	0.67
1:A:82:ILE:HD12	1:A:83:THR:N	2.08	0.67
2:B:94:ARG:HG3	2:B:94:ARG:HH11	1.60	0.66
1:A:16:PRO:HD2	2:B:6:ARG:HD3	1.77	0.66
4:D:87:SER:H	4:D:157:GLN:NE2	1.92	0.66
1:A:147:LYS:HE3	1:A:149:HIS:HE1	1.62	0.65
2:B:86:GLY:HA2	2:B:89:PHE:CE1	2.32	0.65
4:D:64:VAL:HG22	4:D:109:THR:CG2	2.27	0.64
2:B:159:VAL:C	2:B:160:MET:HE3	2.21	0.64
1:A:82:ILE:HD13	2:B:33:ASN:CB	2.11	0.64
2:B:106:THR:O	2:B:107:GLN:O	2.16	0.64
1:A:7:ILE:HD13	2:B:17:PHE:HE2	1.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:105:LYS:H	2:B:105:LYS:CE	2.10	0.63
4:D:57:LYS:HB3	4:D:58:LEU:HD22	1.81	0.63
2:B:13:PHE:CD2	3:C:29:LEU:HD23	2.35	0.62
2:B:150:ASN:HD22	2:B:154:THR:CG2	2.04	0.62
1:A:23:MET:HE1	1:A:25:ASP:HB2	1.81	0.62
1:A:36:MET:HG2	4:D:45:PHE:CZ	2.35	0.61
2:B:160:MET:HE3	2:B:160:MET:N	2.15	0.61
4:D:76:LYS:HD2	4:D:77:TYR:CE1	2.36	0.61
1:A:89:VAL:HG13	1:A:109:ILE:HD13	1.81	0.61
2:B:23:ARG:HH21	2:B:23:ARG:HG2	1.63	0.61
4:D:122:HIS:O	4:D:152:LYS:HE3	2.00	0.61
2:B:2:ASP:OD1	2:B:4:ARG:HD3	2.00	0.61
2:B:70:GLN:HG3	2:B:71:ARG:N	2.16	0.61
2:B:93:ARG:O	2:B:94:ARG:HG3	2.01	0.60
1:A:22:PHE:CE2	3:C:28:ILE:HD11	2.35	0.60
1:A:72:ILE:HG13	5:A:270:HOH:O	2.02	0.60
2:B:145:THR:HG22	2:B:158:LEU:H	1.65	0.60
1:A:76:ARG:HH22	2:B:57:ASP:CG	2.09	0.60
1:A:95:SER:HB2	1:A:96:PRO:HD2	1.84	0.60
2:B:30:CYS:O	2:B:31:ILE:HD13	2.02	0.60
2:B:28:GLU:HB3	2:B:40:PHE:HB3	1.85	0.59
4:D:58:LEU:HD22	4:D:58:LEU:N	2.18	0.58
4:D:231:GLU:HB3	4:D:233:HIS:CE1	2.39	0.58
1:A:73:MET:HG3	2:B:9:TRP:CZ3	2.39	0.57
5:A:270:HOH:O	3:C:34:VAL:HG22	2.04	0.57
1:A:143:HIS:HD2	2:B:12:LYS:NZ	2.02	0.57
2:B:97:PRO:HB3	2:B:122:PHE:HB3	1.86	0.57
2:B:116:VAL:HG13	2:B:160:MET:CE	2.30	0.57
4:D:191:ASN:HB3	4:D:227:SER:HB2	1.86	0.57
2:B:111:HIS:CG	2:B:112:HIS:H	2.24	0.56
4:D:231:GLU:HB3	4:D:233:HIS:HE1	1.70	0.56
2:B:65:LYS:H	2:B:65:LYS:HD2	1.70	0.56
1:A:13:TYR:CZ	1:A:67:LYS:HD2	2.41	0.56
2:B:31:ILE:HD12	2:B:36:GLU:HA	1.86	0.56
1:A:94:ASN:ND2	1:A:104:VAL:HB	2.20	0.55
1:A:109:ILE:CD1	1:A:165:VAL:HG21	2.37	0.55
2:B:161:LEU:HG	2:B:163:THR:HG23	1.89	0.55
4:D:88:ASN:H	4:D:88:ASN:ND2	2.02	0.55
1:A:107:CYS:HB2	1:A:121:TRP:CH2	2.42	0.55
1:A:26:PHE:HB2	1:A:31:ILE:HD11	1.89	0.54
2:B:31:ILE:HD12	2:B:36:GLU:CA	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:VAL:HG13	1:A:109:ILE:CD1	2.38	0.53
3:C:34:VAL:HB	3:C:35:PRO:HD2	1.91	0.53
1:A:36:MET:HG2	4:D:45:PHE:CE1	2.43	0.53
1:A:73:MET:CE	2:B:53:LEU:HD13	2.39	0.53
4:D:6:PRO:HD3	4:D:197:TRP:CE2	2.44	0.53
2:B:29:ARG:NH2	2:B:36:GLU:OE1	2.40	0.52
2:B:94:ARG:HG3	2:B:94:ARG:NH1	2.23	0.52
4:D:216:MET:C	4:D:218:ASN:H	2.17	0.52
2:B:161:LEU:HG	2:B:163:THR:CG2	2.40	0.52
1:A:107:CYS:SG	1:A:109:ILE:HD11	2.50	0.51
2:B:107:GLN:HE21	2:B:114:LEU:H	1.59	0.51
2:B:114:LEU:CD2	2:B:160:MET:HB3	2.34	0.51
2:B:166:ARG:N	2:B:169:GLU:OE1	2.44	0.51
2:B:127:ILE:HD12	2:B:129:VAL:CG1	2.41	0.51
4:D:42:ASP:OD1	4:D:97:SER:HA	2.11	0.50
2:B:52:GLU:OE1	2:B:55:ARG:HD2	2.11	0.50
2:B:93:ARG:HG2	2:B:123:TYR:CD1	2.47	0.50
1:A:10:ALA:CB	1:A:23:MET:HE2	2.34	0.50
4:D:6:PRO:HB3	4:D:197:TRP:CZ2	2.47	0.50
1:A:23:MET:O	1:A:23:MET:HE3	2.12	0.50
2:B:116:VAL:HG22	2:B:160:MET:HG3	1.94	0.50
5:B:217:HOH:O	3:C:27:GLY:HA3	2.11	0.50
1:A:17:ASP:C	1:A:18:GLN:HG2	2.38	0.49
2:B:166:ARG:O	2:B:169:GLU:HG3	2.13	0.49
2:B:57:ASP:OD1	3:C:34:VAL:HB	2.13	0.49
4:D:110:CYS:HB3	5:D:242:HOH:O	2.11	0.49
1:A:98:GLU:HB2	1:A:101:GLU:HG3	1.94	0.48
2:B:130:ARG:HB2	2:B:132:PHE:HE1	1.77	0.48
2:B:145:THR:CG2	2:B:158:LEU:H	2.25	0.48
2:B:18:PHE:CD1	2:B:18:PHE:N	2.80	0.48
2:B:31:ILE:CD1	2:B:36:GLU:CA	2.89	0.48
2:B:127:ILE:HG12	2:B:177:HIS:HB2	1.96	0.48
4:D:33:VAL:O	4:D:85:TYR:HA	2.14	0.48
1:A:81:PRO:HB3	2:B:5:PRO:HB2	1.94	0.48
2:B:129:VAL:HG21	2:B:159:VAL:CG2	2.32	0.48
4:D:124:ASP:O	4:D:125:ASN:CB	2.62	0.48
4:D:131:VAL:CG1	4:D:230:ILE:HD13	2.43	0.48
4:D:83:ASP:OD1	4:D:118:HIS:HD2	1.97	0.48
2:B:118:SER:HA	2:B:158:LEU:HD23	1.96	0.48
4:D:97:SER:O	4:D:98:LYS:CB	2.62	0.47
1:A:23:MET:HE3	1:A:23:MET:C	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:LEU:HD23	1:A:127:PRO:HA	1.96	0.47
4:D:6:PRO:HB3	4:D:197:TRP:CH2	2.49	0.47
4:D:39:LYS:NZ	4:D:39:LYS:HB2	2.29	0.47
4:D:86:GLY:HA3	4:D:111:MET:HE3	1.95	0.47
4:D:54:SER:HA	4:D:62:ASP:HA	1.97	0.47
1:A:22:PHE:CG	1:A:59:ALA:HB1	2.50	0.47
2:B:129:VAL:HA	2:B:174:GLN:O	2.15	0.47
4:D:42:ASP:HB3	4:D:50:ILE:HB	1.96	0.47
4:D:97:SER:O	4:D:98:LYS:HB2	2.13	0.47
1:A:42:VAL:HG22	1:A:42:VAL:O	2.13	0.47
4:D:2:SER:HB2	4:D:195:THR:H	1.80	0.47
4:D:96:SER:O	4:D:97:SER:HB3	2.15	0.47
1:A:17:ASP:O	1:A:18:GLN:HG2	2.15	0.46
2:B:166:ARG:O	2:B:167:SER:C	2.59	0.46
1:A:118:ASN:HB3	1:A:166:GLU:HB2	1.96	0.46
1:A:36:MET:HG3	1:A:60:LEU:HD11	1.97	0.46
1:A:36:MET:HG3	1:A:60:LEU:CD1	2.46	0.46
2:B:121:GLY:HA2	2:B:154:THR:HG23	1.97	0.46
4:D:21:MET:CB	4:D:178:SER:H	2.29	0.46
1:A:74:THR:HA	2:B:32:TYR:OH	2.16	0.46
2:B:78:TYR:CD2	3:C:29:LEU:HD22	2.50	0.46
4:D:94:TYR:O	4:D:95:PHE:HB3	2.16	0.46
1:A:12:PHE:HB3	2:B:10:GLN:HG2	1.98	0.46
1:A:124:ASN:OD1	1:A:160:VAL:HG22	2.15	0.46
2:B:65:LYS:HD2	2:B:65:LYS:N	2.31	0.46
2:B:111:HIS:HD1	2:B:112:HIS:H	1.61	0.46
1:A:171:ASP:O	1:A:172:GLU:HG3	2.15	0.46
1:A:70:LEU:HD13	2:B:9:TRP:HB2	1.98	0.45
2:B:57:ASP:OD1	3:C:35:PRO:HD2	2.16	0.45
4:D:129:GLN:OE1	4:D:226:LYS:HA	2.16	0.45
1:A:76:ARG:NH2	2:B:57:ASP:OD2	2.50	0.45
1:A:12:PHE:CD1	1:A:12:PHE:C	2.94	0.45
4:D:18:THR:HB	4:D:206:ASP:HA	1.99	0.45
4:D:59:LYS:HA	4:D:59:LYS:HE2	1.98	0.45
2:B:176:GLU:HG2	2:B:183:PRO:HB3	1.99	0.45
4:D:21:MET:HB2	4:D:178:SER:H	1.81	0.45
4:D:119:GLU:OE2	4:D:119:GLU:HA	2.17	0.45
1:A:22:PHE:CZ	3:C:28:ILE:CD1	2.95	0.45
1:A:122:LEU:HB2	1:A:162:ASP:HB2	1.98	0.45
2:B:130:ARG:NH1	2:B:174:GLN:HE21	2.15	0.45
4:D:68:LEU:HB3	4:D:73:LEU:HD23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:137:GLU:HB2	4:D:142:THR:HG21	1.99	0.44
1:A:17:ASP:CG	2:B:6:ARG:HH11	2.26	0.44
4:D:31:HIS:O	4:D:32:TYR:HB3	2.17	0.44
4:D:88:ASN:HA	4:D:110:CYS:O	2.17	0.44
2:B:137:GLU:HG2	2:B:139:LYS:HE3	2.00	0.44
1:A:110:ASP:OD1	1:A:140:ARG:HD2	2.17	0.44
2:B:129:VAL:O	2:B:129:VAL:HG23	2.16	0.44
4:D:25:LYS:HE2	4:D:29:ASP:OD1	2.18	0.43
4:D:91:VAL:O	4:D:92:ASN:HB2	2.18	0.43
1:A:162:ASP:HA	1:A:176:LYS:O	2.17	0.43
2:B:115:LEU:O	2:B:160:MET:HA	2.18	0.43
1:A:73:MET:HE1	2:B:53:LEU:HD13	1.99	0.43
2:B:23:ARG:HG2	2:B:23:ARG:NH2	2.32	0.43
1:A:36:MET:CE	1:A:63:ILE:HG13	2.48	0.43
2:B:31:ILE:CD1	2:B:36:GLU:CB	2.96	0.43
4:D:46:LYS:HD3	4:D:71:GLU:CD	2.44	0.43
4:D:112:TYR:CE2	4:D:215:MET:HG3	2.54	0.43
4:D:170:LYS:HA	4:D:170:LYS:HD3	1.92	0.43
1:A:36:MET:HE3	1:A:63:ILE:HG13	2.00	0.42
4:D:86:GLY:CA	4:D:157:GLN:NE2	2.82	0.42
2:B:129:VAL:CG2	2:B:129:VAL:O	2.66	0.42
2:B:65:LYS:HA	2:B:65:LYS:HE2	2.01	0.42
4:D:3:GLN:HA	4:D:4:PRO:HD3	1.84	0.42
1:A:69:ASN:CG	3:C:34:VAL:HG13	2.45	0.42
1:A:143:HIS:HD2	2:B:12:LYS:HZ1	1.68	0.42
1:A:9:GLN:HG3	1:A:24:PHE:CE1	2.55	0.42
1:A:22:PHE:CD1	1:A:22:PHE:C	2.98	0.42
4:D:24:MET:HE1	4:D:213:TYR:CZ	2.55	0.42
4:D:39:LYS:NZ	4:D:39:LYS:CB	2.83	0.41
2:B:31:ILE:HD12	2:B:36:GLU:CB	2.50	0.41
2:B:31:ILE:HD12	2:B:36:GLU:HB2	2.02	0.41
2:B:130:ARG:HB2	2:B:132:PHE:CE1	2.54	0.41
2:B:177:HIS:CG	2:B:178:PRO:HD2	2.55	0.41
1:A:42:VAL:O	1:A:42:VAL:CG2	2.67	0.41
2:B:131:TRP:HB3	2:B:161:LEU:HD22	2.02	0.41
3:C:29:LEU:HD12	3:C:29:LEU:HA	1.80	0.41
2:B:166:ARG:HB2	2:B:169:GLU:CD	2.45	0.41
1:A:23:MET:HE1	1:A:25:ASP:CB	2.48	0.41
2:B:2:ASP:OD1	2:B:6:ARG:NH2	2.53	0.41
2:B:90:THR:OG1	2:B:91:VAL:N	2.54	0.41
1:A:76:ARG:HB3	5:C:275:HOH:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2:ASP:CG	2:B:6:ARG:NH2	2.73	0.41
2:B:111:HIS:CG	2:B:112:HIS:N	2.89	0.41
4:D:154:VAL:HG22	4:D:155:THR:N	2.36	0.41
1:A:72:ILE:CG1	5:A:270:HOH:O	2.66	0.40
1:A:60:LEU:N	1:A:60:LEU:HD22	2.36	0.40
1:A:122:LEU:O	1:A:161:TYR:HA	2.21	0.40
4:D:131:VAL:HG11	4:D:230:ILE:HD13	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	175/177 (99%)	167 (95%)	8 (5%)	0	100	100
2	B	183/190 (96%)	173 (94%)	9 (5%)	1 (0%)	24	37
3	C	13/15 (87%)	13 (100%)	0	0	100	100
4	D	227/239 (95%)	212 (93%)	11 (5%)	4 (2%)	6	9
All	All	598/621 (96%)	565 (94%)	28 (5%)	5 (1%)	16	25

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	57	LYS
4	D	125	ASN
4	D	124	ASP
2	B	32	TYR
4	D	178	SER

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	162/162 (100%)	155 (96%)	7 (4%)	26	44
2	B	168/171 (98%)	149 (89%)	19 (11%)	5	8
3	C	10/10 (100%)	9 (90%)	1 (10%)	7	11
4	D	213/220 (97%)	208 (98%)	5 (2%)	44	66
All	All	553/563 (98%)	521 (94%)	32 (6%)	18	32

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	HIS
1	A	11	GLU
1	A	23	MET
1	A	36	MET
1	A	42	VAL
1	A	82	ILE
1	A	146	ARG
2	B	18	PHE
2	B	23	ARG
2	B	29	ARG
2	B	34	GLN
2	B	35	GLU
2	B	38	VAL
2	B	53	LEU
2	B	59	GLU
2	B	65	LYS
2	B	68	LEU
2	B	69	GLU
2	B	70	GLN
2	B	98	LYS
2	B	105	LYS
2	B	114	LEU
2	B	127	ILE
2	B	129	VAL

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Mol	Chain	Res	Type
2	B	137	GLU
2	B	154	THR
3	C	24	GLU
4	D	39	LYS
4	D	88	ASN
4	D	125	ASN
4	D	207	LYS
4	D	218	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	143	HIS
1	A	149	HIS
2	B	10	GLN
2	B	34	GLN
2	B	70	GLN
2	B	92	GLN
2	B	107	GLN
2	B	113	ASN
2	B	134	ASN
2	B	150	ASN
2	B	156	GLN
2	B	174	GLN
4	D	60	ASN
4	D	88	ASN
4	D	92	ASN
4	D	118	HIS
4	D	125	ASN
4	D	129	GLN
4	D	141	ASN
4	D	157	GLN
4	D	218	ASN
4	D	233	HIS

5.3.3 RNA ⓘ

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	28:ILE	C	29:LEU	N	1.16

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	177/177 (100%)	-0.16	3 (1%) 69 65	12, 25, 51, 63	0
2	B	187/190 (98%)	0.26	10 (5%) 32 28	13, 33, 66, 86	0
3	C	15/15 (100%)	0.48	2 (13%) 7 5	19, 26, 61, 64	0
4	D	231/239 (96%)	0.03	16 (6%) 23 19	16, 31, 64, 75	0
All	All	610/621 (98%)	0.06	31 (5%) 33 30	12, 31, 62, 86	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	107	GLN	5.1
2	B	106	THR	5.0
2	B	111	HIS	4.9
3	C	23	GLY	4.5
4	D	239	GLY	3.6
4	D	56	LYS	3.5
4	D	1	GLU	3.5
4	D	97	SER	3.3
4	D	177	ASN	3.0
4	D	125	ASN	3.0
3	C	24	GLU	2.9
2	B	136	GLN	2.9
1	A	94	ASN	2.6
4	D	237	LYS	2.6
2	B	1	GLY	2.5
4	D	59	LYS	2.5
4	D	124	ASP	2.4
2	B	112	HIS	2.4
2	B	135	GLY	2.4
4	D	126	GLY	2.3
4	D	57	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
4	D	98	LYS	2.3
4	D	58	LEU	2.2
1	A	158	GLU	2.1
2	B	105	LYS	2.1
4	D	7	MET	2.1
2	B	18	PHE	2.1
4	D	60	ASN	2.0
1	A	160	VAL	2.0
4	D	176	PHE	2.0
2	B	127	ILE	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.