



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:57 PM UTC

PDB ID : 5WKF / pdb_00005wkf
Title : D30 TCR in complex with HLA-A*11:01-GTS1
Authors : Gras, S.; Rossjohn, J.
Deposited on : 2017-07-25
Resolution : 2.95 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

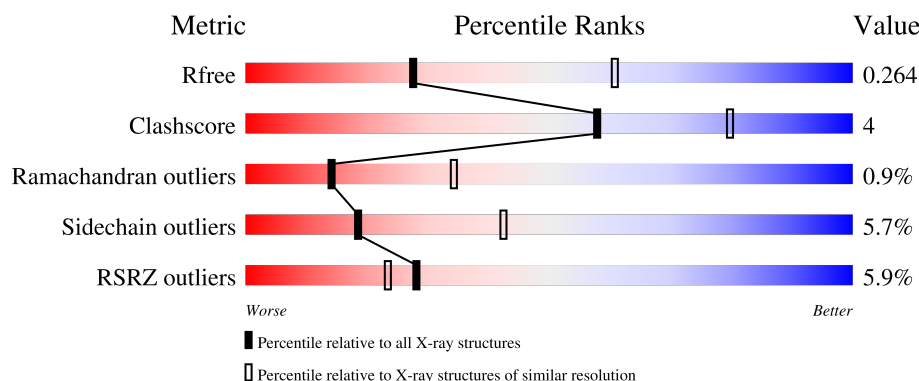
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	274	85% 14% .
1	F	274	91% 9%
2	B	100	89% 10% .
2	G	100	3% 83% 17%
3	C	10	80% 20%

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Mol	Chain	Length	Quality of chain
3	H	10	90% 10%
4	D	198	17% 78% 19% • •
4	I	198	17% 78% 18% • •
5	E	244	6% 81% 18% •
5	J	244	3% 80% 16% •

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13580 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called HLA class I histocompatibility antigen, A-11 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2235	1389	407	430	9			
1	F	274	Total	C	N	O	S	0	0	0
			2235	1389	407	430	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
			829	528	140	158	3			
2	G	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
G	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called GTS1 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			69	40	14	15			
3	H	10	Total	C	N	O	0	0	0
			69	40	14	15			

- Molecule 4 is a protein called T-cell receptor alpha variable 30,T-cell receptor, sp3.4 alpha chain Chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	198	Total 1536	C 962	N 259	O 305	S 10	0	0	0
4	I	198	Total 1536	C 962	N 259	O 305	S 10	0	0	0

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	20	ILE	VAL	conflict	UNP A0A087WSZ9
D	44	ILE	VAL	conflict	UNP A0A087WSZ9
D	91	LEU	-	linker	UNP A0A087WSZ9
D	92	GLY	-	linker	UNP A0A087WSZ9
D	93	ASP	-	linker	UNP A0A087WSZ9
D	94	ALA	-	linker	UNP A0A087WSZ9
D	95	GLY	-	linker	UNP A0A087WSZ9
D	96	ASN	-	linker	UNP A0A087WSZ9
D	97	MET	-	linker	UNP A0A087WSZ9
D	98	LEU	-	linker	UNP A0A087WSZ9
D	99	THR	-	linker	UNP A0A087WSZ9
D	100	PHE	-	linker	UNP A0A087WSZ9
D	101	GLY	-	linker	UNP A0A087WSZ9
D	102	GLY	-	linker	UNP A0A087WSZ9
D	103	GLY	-	linker	UNP A0A087WSZ9
D	104	THR	-	linker	UNP A0A087WSZ9
D	105	ARG	-	linker	UNP A0A087WSZ9
D	106	LEU	-	linker	UNP A0A087WSZ9
D	107	MET	-	linker	UNP A0A087WSZ9
D	108	VAL	-	linker	UNP A0A087WSZ9
D	109	LYS	-	linker	UNP A0A087WSZ9
D	110	PRO	-	linker	UNP A0A087WSZ9
D	111	HIS	-	linker	UNP A0A087WSZ9
I	20	ILE	VAL	conflict	UNP A0A087WSZ9
I	51	ILE	VAL	conflict	UNP A0A087WSZ9
I	102	LEU	-	linker	UNP A0A087WSZ9
I	103	GLY	-	linker	UNP A0A087WSZ9
I	105	ASP	-	linker	UNP A0A087WSZ9
I	106	ALA	-	linker	UNP A0A087WSZ9
I	107	GLY	-	linker	UNP A0A087WSZ9
I	108	ASN	-	linker	UNP A0A087WSZ9
I	109	MET	-	linker	UNP A0A087WSZ9
I	110	LEU	-	linker	UNP A0A087WSZ9
I	111	THR	-	linker	UNP A0A087WSZ9
I	112	PHE	-	linker	UNP A0A087WSZ9

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Chain	Residue	Modelled	Actual	Comment	Reference
I	113	GLY	-	linker	UNP A0A087WSZ9
I	114	GLY	-	linker	UNP A0A087WSZ9
I	115	GLY	-	linker	UNP A0A087WSZ9
I	116	THR	-	linker	UNP A0A087WSZ9
I	117	ARG	-	linker	UNP A0A087WSZ9
I	118	LEU	-	linker	UNP A0A087WSZ9
I	119	MET	-	linker	UNP A0A087WSZ9
I	120	VAL	-	linker	UNP A0A087WSZ9
I	121	LYS	-	linker	UNP A0A087WSZ9
I	122	PRO	-	linker	UNP A0A087WSZ9
I	123	HIS	-	linker	UNP A0A087WSZ9

- Molecule 5 is a protein called D30 TCR beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	244	Total	C	N	O	S	0	0	0
			1935	1229	336	365	5			
5	J	244	Total	C	N	O	S	0	0	0
			1935	1229	336	365	5			

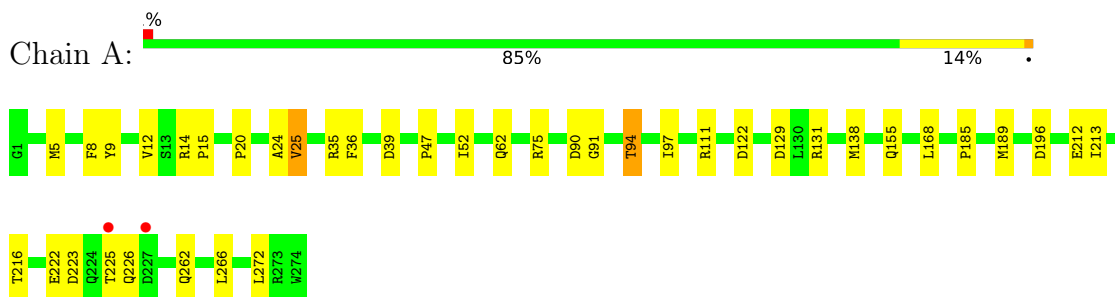
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	60	Total	O	0	0
			60	60		
6	B	28	Total	O	0	0
			28	28		
6	C	4	Total	O	0	0
			4	4		
6	D	35	Total	O	0	0
			35	35		
6	E	36	Total	O	0	0
			36	36		
6	F	69	Total	O	0	0
			69	69		
6	G	34	Total	O	0	0
			34	34		
6	H	2	Total	O	0	0
			2	2		
6	I	31	Total	O	0	0
			31	31		
6	J	65	Total	O	0	0
			65	65		

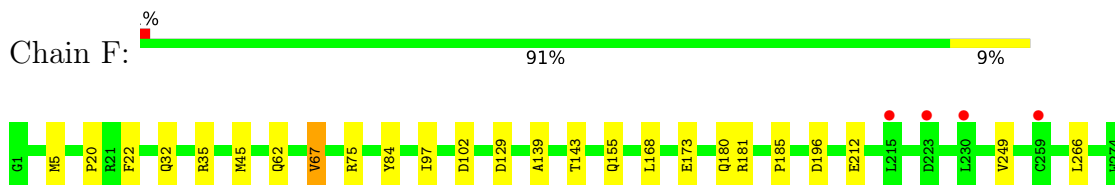
3 Residue-property plots

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.

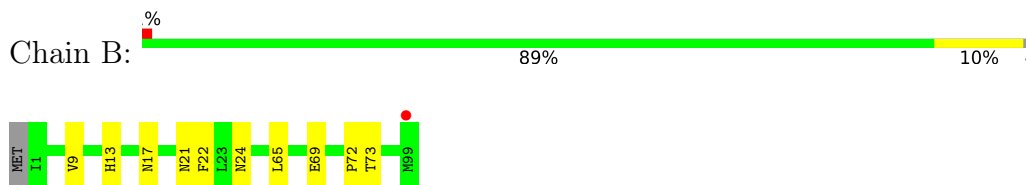
- Molecule 1: HLA class I histocompatibility antigen, A-11 alpha chain



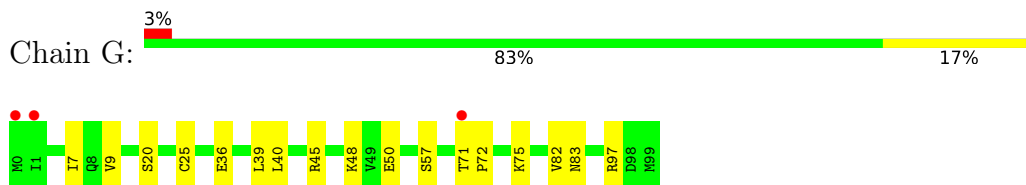
- Molecule 1: HLA class I histocompatibility antigen, A-11 alpha chain



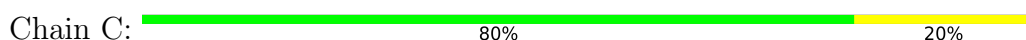
- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin

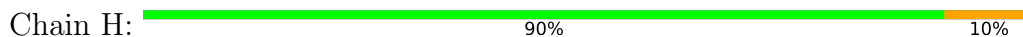


- Molecule 3: GTS1 peptide

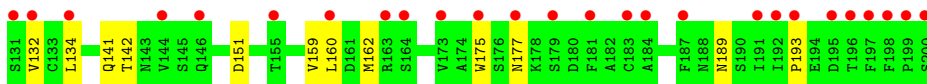
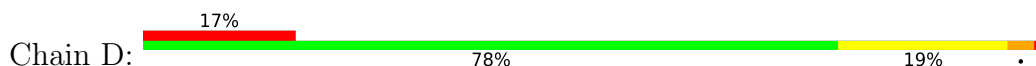




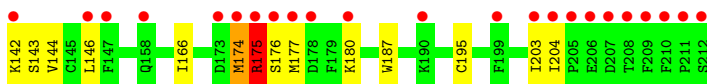
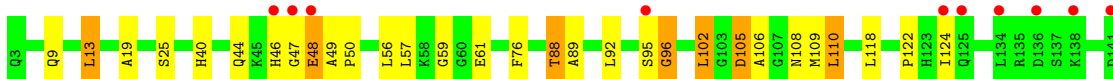
- Molecule 3: GTS1 peptide



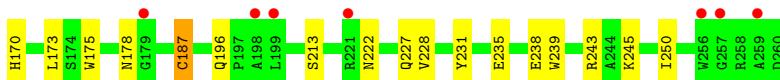
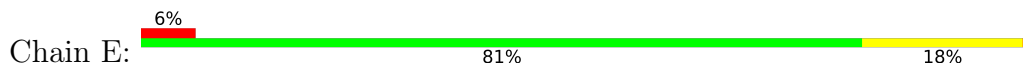
- Molecule 4: T-cell receptor alpha variable 30,T-cell receptor, sp3.4 alpha chain Chimera



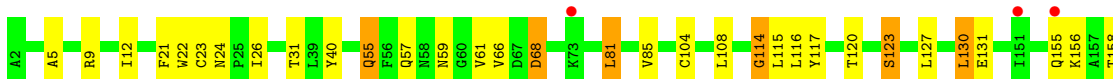
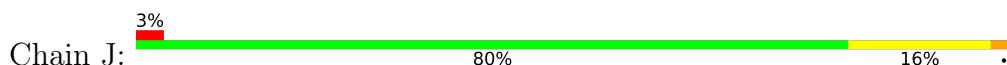
- Molecule 4: T-cell receptor alpha variable 30,T-cell receptor, sp3.4 alpha chain Chimera

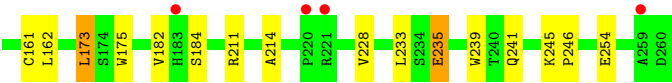


- Molecule 5: D30 TCR beta chain



- Molecule 5: D30 TCR beta chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.06Å 146.38Å 167.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.17 – 2.95 47.17 – 2.95	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.17-2.95) 100.0 (47.17-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.96Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.190 , 0.253 0.201 , 0.264	Depositor DCC
R_{free} test set	2276 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	41.0	Xtriage
Anisotropy	0.587	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13580	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.67	0/2296	1.23	5/3117 (0.2%)
1	F	0.66	0/2296	1.24	3/3117 (0.1%)
2	B	0.67	0/852	1.11	0/1152
2	G	0.67	0/860	1.15	0/1162
3	C	0.52	0/69	1.13	0/91
3	H	0.57	0/69	1.15	0/91
4	D	0.83	1/1569 (0.1%)	1.23	6/2117 (0.3%)
4	I	0.76	0/1569	1.24	3/2117 (0.1%)
5	E	0.69	0/1986	1.14	4/2702 (0.1%)
5	J	0.70	0/1986	1.15	6/2702 (0.2%)
All	All	0.70	1/13552 (0.0%)	1.20	27/18368 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	98	LEU	CA-C	-5.12	1.46	1.52

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	53	GLY	CA-C-N	6.70	133.76	121.70
4	D	53	GLY	C-N-CA	6.70	133.76	121.70
1	A	225	THR	N-CA-C	-6.38	105.51	113.55
1	A	111	ARG	CA-C-N	6.31	126.94	122.33
1	A	111	ARG	C-N-CA	6.31	126.94	122.33
4	I	175	ARG	CA-C-N	5.92	132.85	121.54
4	I	175	ARG	C-N-CA	5.92	132.85	121.54
4	I	105	ASP	CA-CB-CG	5.73	118.33	112.60
1	A	129	ASP	CA-CB-CG	5.65	118.25	112.60
1	F	196	ASP	N-CA-C	-5.60	107.69	114.75
1	F	129	ASP	CA-CB-CG	5.49	118.09	112.60
4	D	89	CYS	N-CA-C	-5.36	100.51	109.24
5	J	114	GLY	CA-C-N	5.26	131.60	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	J	114	GLY	C-N-CA	5.26	131.60	121.54
5	E	80	ARG	CA-C-N	5.26	127.59	120.38
5	E	80	ARG	C-N-CA	5.26	127.59	120.38
4	D	91	LEU	N-CA-C	5.23	117.63	109.52
5	J	108	LEU	CA-C-N	5.21	126.87	120.51
5	J	108	LEU	C-N-CA	5.21	126.87	120.51
1	F	102	ASP	CA-CB-CG	5.17	117.77	112.60
5	J	66	VAL	N-CA-C	-5.14	107.14	111.56
5	E	96	LEU	CA-C-N	5.08	127.59	120.28
5	E	96	LEU	C-N-CA	5.08	127.59	120.28
5	J	68	ASP	CA-CB-CG	5.05	117.66	112.60
1	A	122	ASP	CA-CB-CG	5.00	117.61	112.60
4	D	177	ASN	CA-C-N	5.00	127.96	120.90
4	D	177	ASN	C-N-CA	5.00	127.96	120.90

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	2235	0	2087	17	0
1	F	2235	0	2087	10	0
2	B	829	0	794	4	0
2	G	837	0	803	7	0
3	C	69	0	71	2	0
3	H	69	0	71	3	0
4	D	1536	0	1479	27	0
4	I	1536	0	1477	21	0
5	E	1935	0	1874	21	0
5	J	1935	0	1874	20	0
6	A	60	0	0	0	0
6	B	28	0	0	0	0
6	C	4	0	0	0	0
6	D	35	0	0	0	0
6	E	36	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	F	69	0	0	1	0
6	G	34	0	0	0	0
6	H	2	0	0	0	0
6	I	31	0	0	0	0
6	J	65	0	0	0	0
All	All	13580	0	12617	114	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 4.

All (114) 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)
4:D:53:GLY:HA2	4:D:54:GLU:HB2	1.26	1.11
1:A:62:GLN:HE22	4:D:94:ALA:H	1.06	0.97
4:D:53:GLY:HA2	4:D:54:GLU:CB	2.09	0.82
4:D:53:GLY:CA	4:D:54:GLU:HB2	2.10	0.77
1:F:155:GLN:NE2	5:J:114:GLY:O	2.21	0.74
4:D:8:PRO:HD2	4:D:21:ILE:HG22	1.70	0.74
1:A:62:GLN:NE2	4:D:94:ALA:H	1.87	0.68
4:D:132:VAL:HG23	4:D:175:TRP:HB3	1.75	0.67
4:I:47:GLY:HA2	4:I:48:GLU:HB3	1.76	0.67
5:J:173:LEU:HG	5:J:228:VAL:HG22	1.76	0.67
5:J:12:ILE:HG23	5:J:130:LEU:HD23	1.77	0.65
5:E:114:GLY:HA2	5:E:117:TYR:CD2	2.31	0.65
1:F:97:ILE:HD11	3:H:10:ARG:HH22	1.60	0.65
5:J:158:THR:HB	5:J:211:ARG:HD3	1.81	0.62
1:A:15:PRO:HG2	1:A:91:GLY:H	1.65	0.61
1:A:20:PRO:HG2	1:A:75:ARG:HG2	1.83	0.61
1:F:5:MET:HB2	1:F:168:LEU:HD13	1.83	0.60
4:D:54:GLU:HA	4:D:65:PHE:HB3	1.82	0.59
5:J:239:TRP:HB2	5:J:245:LYS:HD3	1.86	0.58
4:I:57:LEU:HD21	5:J:116:LEU:HD23	1.85	0.58
5:E:161:CYS:HB2	5:E:175:TRP:CZ2	2.39	0.57
5:J:233:LEU:HD12	5:J:246:PRO:HD2	1.86	0.56
2:B:13:HIS:HB2	2:B:21:ASN:HD21	1.71	0.56
4:D:160:LEU:HG	5:E:187:CYS:HB3	1.87	0.55
3:C:6:PRO:HA	5:E:113:GLN:O	2.06	0.55
4:D:93:ASP:HB3	4:D:97:MET:H	1.72	0.55
5:J:61:VAL:HG11	5:J:68:ASP:HB3	1.89	0.55
4:I:105:ASP:HB2	4:I:109:MET:H	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:239:TRP:HB2	5:E:245:LYS:HD3	1.87	0.55
1:F:143:THR:HG23	3:H:10:ARG:HB2	1.89	0.54
5:J:26:ILE:HD11	5:J:120:THR:HG21	1.89	0.54
4:D:50:LEU:HD12	4:D:50:LEU:N	2.23	0.54
1:F:62:GLN:HE22	4:I:106:ALA:H	1.55	0.54
1:F:185:PRO:HD2	1:F:266:LEU:HD13	1.89	0.54
4:D:110:PRO:HG3	4:D:159:VAL:HG11	1.90	0.53
4:I:61:GLU:HA	4:I:76:PHE:HB3	1.92	0.52
5:E:40:TYR:HB2	5:E:105:ALA:HB3	1.91	0.52
4:I:49:ALA:HB2	5:J:123:SER:HA	1.90	0.52
1:A:223:ASP:HB2	4:I:175:ARG:NH1	2.24	0.52
4:I:88:THR:HG23	4:I:89:ALA:H	1.75	0.52
1:A:24:ALA:HB3	1:A:36:PHE:HB3	1.92	0.52
1:F:22:PHE:HE2	1:F:67:VAL:HG22	1.75	0.52
1:A:8:PHE:HB2	1:A:25:VAL:HG23	1.92	0.51
4:I:174:MET:HA	4:I:177:MET:HE2	1.91	0.51
5:E:40:TYR:CE1	5:E:55:GLN:HG3	2.45	0.51
1:A:12:VAL:HG22	1:A:94:THR:HG23	1.92	0.51
5:E:170:HIS:HB3	5:E:231:TYR:HB2	1.92	0.51
1:A:185:PRO:HD2	1:A:266:LEU:HD13	1.93	0.50
4:D:21:ILE:HG21	4:D:104:THR:HG21	1.93	0.50
2:G:25:CYS:HB2	2:G:39:LEU:HD21	1.93	0.50
1:A:189:MET:HE3	1:A:272:LEU:HB3	1.94	0.50
4:I:25:SER:HB2	4:I:102:LEU:HD21	1.94	0.49
4:I:92:LEU:HD11	4:I:122:PRO:HB3	1.94	0.49
4:D:33:HIS:HB3	4:D:35:TYR:CE1	2.48	0.48
4:D:126:LYS:HG3	4:D:127:SER:H	1.79	0.48
5:E:75:ARG:HG2	5:E:93:PRO:HD2	1.95	0.48
1:A:97:ILE:HD11	3:C:10:ARG:HH22	1.79	0.48
1:F:20:PRO:HD2	1:F:75:ARG:HG2	1.96	0.48
4:I:175:ARG:HA	5:J:184:SER:HB3	1.95	0.48
4:D:33:HIS:CG	4:D:98:LEU:HD11	2.48	0.47
5:J:5:ALA:HB3	5:J:24:ASN:HB3	1.97	0.47
4:D:33:HIS:ND1	4:D:98:LEU:HD11	2.29	0.47
4:I:57:LEU:HD12	4:I:59:GLY:H	1.80	0.47
5:J:22:TRP:CH2	5:J:24:ASN:HB2	2.50	0.47
1:A:9:TYR:HB2	1:A:97:ILE:HB	1.96	0.46
5:E:156:LYS:HG3	5:E:213:SER:HA	1.97	0.46
4:D:93:ASP:HB3	4:D:97:MET:N	2.31	0.46
5:J:40:TYR:CE1	5:J:55:GLN:HG3	2.50	0.46
4:I:95:SER:HA	4:I:96:GLY:HA2	1.74	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:J:155:GLN:O	5:J:214:ALA:HB2	2.16	0.45
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.98	0.45
2:B:17:ASN:HA	2:B:72:PRO:O	2.16	0.45
5:E:71:LEU:HD22	5:E:76:PHE:HB3	1.97	0.45
4:D:81:LEU:HD21	4:D:110:PRO:HB3	1.98	0.45
5:E:178:ASN:HD21	5:E:222:ASN:HD22	1.64	0.45
2:G:40:LEU:HD23	2:G:45:ARG:HA	1.99	0.45
4:I:40:HIS:CG	4:I:110:LEU:HD11	2.52	0.45
5:E:21:PHE:HZ	5:E:127:LEU:HD22	1.81	0.45
5:E:136:VAL:O	5:E:243:ARG:NH2	2.46	0.45
4:D:162:MET:HE1	5:E:156:LYS:HG2	1.99	0.45
5:E:22:TRP:CH2	5:E:24:ASN:HB2	2.52	0.44
5:J:156:LYS:HB2	5:J:211:ARG:HH21	1.83	0.44
5:E:15:LYS:O	5:E:16:ARG:HB2	2.17	0.44
4:I:144:VAL:HG12	4:I:187:TRP:HB3	2.00	0.44
4:D:142:THR:HG21	4:D:193:PRO:HG3	1.99	0.44
4:I:56:LEU:HB3	4:I:76:PHE:HB2	2.00	0.43
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.99	0.43
5:J:161:CYS:HB2	5:J:175:TRP:CZ2	2.53	0.43
2:G:20:SER:HA	2:G:71:THR:HG22	2.00	0.43
1:A:62:GLN:HE22	4:D:94:ALA:N	1.91	0.43
2:B:22:PHE:CE1	2:B:69:GLU:HG2	2.53	0.43
4:D:19:ALA:HB3	4:D:76:LEU:HB2	2.00	0.42
4:I:108:ASN:HB2	5:J:115:LEU:H	1.83	0.42
5:E:173:LEU:HG	5:E:228:VAL:HG22	2.01	0.42
2:G:7:ILE:HG12	2:G:82:VAL:HG21	2.01	0.42
4:D:50:LEU:HD12	4:D:50:LEU:H	1.83	0.42
4:I:13:LEU:HD11	4:I:19:ALA:HB2	2.01	0.42
5:J:235:GLU:CD	5:J:235:GLU:H	2.28	0.42
4:I:142:LYS:HG3	4:I:143:SER:H	1.85	0.42
4:D:4:PRO:HB3	4:D:91:LEU:HG	2.02	0.42
1:F:84:TYR:HB3	1:F:139:ALA:HB1	2.02	0.42
4:I:44:GLN:HB3	4:I:50:PRO:HA	2.02	0.42
1:A:47:PRO:HB3	1:A:52:ILE:HG23	2.02	0.41
5:E:227:GLN:HG3	5:E:250:ILE:HG23	2.01	0.41
4:D:52:GLY:HA3	6:F:354:HOH:O	2.21	0.41
1:A:14:ARG:HH22	1:A:39:ASP:CG	2.28	0.41
2:G:36:GLU:HG3	2:G:83:ASN:HB3	2.02	0.41
5:J:21:PHE:HZ	5:J:127:LEU:HD22	1.84	0.41
1:F:143:THR:CG2	3:H:10:ARG:HB2	2.49	0.41
5:E:45:ILE:O	5:E:48:GLN:HG2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:110:PRO:HG3	4:D:159:VAL:CG1	2.51	0.40
2:G:75:LYS:H	2:G:75:LYS:HG3	1.68	0.40
2:G:72:PRO:HB2	2:G:97:ARG:HH12	1.86	0.40
1:A:155:GLN:NE2	5:E:116:LEU:H	2.20	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	272/274 (99%)	265 (97%)	7 (3%)	0	100	100
1	F	272/274 (99%)	260 (96%)	12 (4%)	0	100	100
2	B	97/100 (97%)	93 (96%)	4 (4%)	0	100	100
2	G	98/100 (98%)	93 (95%)	5 (5%)	0	100	100
3	C	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
3	H	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
4	D	196/198 (99%)	174 (89%)	17 (9%)	5 (3%)	4	12
4	I	196/198 (99%)	179 (91%)	11 (6%)	6 (3%)	3	8
5	E	242/244 (99%)	228 (94%)	12 (5%)	2 (1%)	16	37
5	J	242/244 (99%)	226 (93%)	14 (6%)	2 (1%)	16	37
All	All	1631/1652 (99%)	1532 (94%)	84 (5%)	15 (1%)	14	34

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	54	GLU
4	I	176	SER
5	J	81	LEU

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Mol	Chain	Res	Type
4	D	92	GLY
4	I	48	GLU
4	D	39	HIS
4	D	93	ASP
4	D	189	ASN
4	I	46	HIS
5	E	16	ARG
4	I	88	THR
4	I	96	GLY
5	J	85	VAL
4	I	203	ILE
5	E	85	VAL

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	231/231 (100%)	218 (94%)	13 (6%)	19	43
1	F	231/231 (100%)	222 (96%)	9 (4%)	28	54
2	B	94/95 (99%)	92 (98%)	2 (2%)	47	69
2	G	95/95 (100%)	91 (96%)	4 (4%)	26	53
3	C	8/8 (100%)	8 (100%)	0	100	100
3	H	8/8 (100%)	7 (88%)	1 (12%)	4	13
4	D	173/173 (100%)	162 (94%)	11 (6%)	16	38
4	I	173/173 (100%)	160 (92%)	13 (8%)	12	31
5	E	209/209 (100%)	199 (95%)	10 (5%)	23	48
5	J	209/209 (100%)	191 (91%)	18 (9%)	10	25
All	All	1431/1432 (100%)	1350 (94%)	81 (6%)	18	42

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	VAL

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Mol	Chain	Res	Type
1	A	35	ARG
1	A	90	ASP
1	A	94	THR
1	A	131	ARG
1	A	138	MET
1	A	196	ASP
1	A	212	GLU
1	A	213	ILE
1	A	216	THR
1	A	222	GLU
1	A	226	GLN
1	A	262	GLN
2	B	9	VAL
2	B	73	THR
4	D	11	VAL
4	D	12	ILE
4	D	13	LEU
4	D	38	LYS
4	D	44	ILE
4	D	50	LEU
4	D	84	SER
4	D	91	LEU
4	D	134	LEU
4	D	141	GLN
4	D	151	ASP
5	E	16	ARG
5	E	31	THR
5	E	46	LEU
5	E	51	LYS
5	E	115	LEU
5	E	156	LYS
5	E	187	CYS
5	E	196	GLN
5	E	235	GLU
5	E	238	GLU
1	F	32	GLN
1	F	35	ARG
1	F	45	MET
1	F	67	VAL
1	F	173	GLU
1	F	180	GLN
1	F	181	ARG

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Mol	Chain	Res	Type
1	F	212	GLU
1	F	249	VAL
2	G	9	VAL
2	G	48	LYS
2	G	50	GLU
2	G	57	SER
3	H	10	ARG
4	I	9	GLN
4	I	13	LEU
4	I	102	LEU
4	I	110	LEU
4	I	118	LEU
4	I	124	ILE
4	I	146	LEU
4	I	166	ILE
4	I	174	MET
4	I	175	ARG
4	I	180	LYS
4	I	195	CYS
4	I	204	ILE
5	J	9	ARG
5	J	23	CYS
5	J	31	THR
5	J	55	GLN
5	J	57	GLN
5	J	59	ASN
5	J	81	LEU
5	J	104	CYS
5	J	117	TYR
5	J	123	SER
5	J	130	LEU
5	J	131	GLU
5	J	162	LEU
5	J	173	LEU
5	J	182	VAL
5	J	235	GLU
5	J	241	GLN
5	J	254	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	62	GLN
1	A	72	GLN
1	A	155	GLN
2	B	21	ASN
4	D	3	GLN
4	D	33	HIS
4	D	37	GLN
4	D	39	HIS
4	D	70	GLN
4	D	71	GLN
4	D	121	GLN
4	D	141	GLN
4	D	185	ASN
5	E	44	GLN
5	E	92	GLN
5	E	113	GLN
5	E	218	GLN
5	E	219	ASN
5	E	222	ASN
5	E	229	GLN
1	F	62	GLN
1	F	66	ASN
1	F	87	GLN
1	F	115	GLN
1	F	155	GLN
1	F	180	GLN
3	H	9	ASN
4	I	40	HIS
4	I	81	GLN
4	I	82	GLN
4	I	153	GLN
4	I	201	ASN
5	J	43	GLN
5	J	57	GLN
5	J	58	ASN
5	J	113	GLN
5	J	134	ASN
5	J	219	ASN
5	J	223	HIS
5	J	229	GLN
5	J	236	ASN
5	J	241	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no 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	274/274 (100%)	0.02	2 (0%) 84 80	18, 37, 70, 96	0
1	F	274/274 (100%)	0.04	4 (1%) 72 65	16, 36, 85, 100	0
2	B	99/100 (99%)	0.24	1 (1%) 79 74	23, 48, 72, 81	0
2	G	100/100 (100%)	0.14	3 (3%) 52 45	22, 39, 63, 71	0
3	C	10/10 (100%)	-0.14	0 100 100	23, 24, 30, 47	0
3	H	10/10 (100%)	-0.65	0 100 100	17, 19, 26, 37	0
4	D	198/198 (100%)	0.90	33 (16%) 4 4	25, 59, 130, 163	1 (0%)
4	I	198/198 (100%)	0.92	33 (16%) 4 4	18, 54, 137, 166	1 (0%)
5	E	244/244 (100%)	0.29	14 (5%) 29 24	18, 40, 86, 102	6 (2%)
5	J	244/244 (100%)	0.13	7 (2%) 53 46	15, 38, 76, 106	6 (2%)
All	All	1651/1652 (99%)	0.31	97 (5%) 28 23	15, 41, 96, 166	14 (0%)

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	I	205	PRO	8.7
4	I	207	ASP	5.4
4	I	175	ARG	5.1
4	I	177	MET	5.0
4	D	132	VAL	5.0
4	I	208	THR	4.6
4	D	197	PHE	4.3
4	I	203	ILE	4.2
4	D	193	PRO	4.0
4	D	192	ILE	3.7
4	D	134	LEU	3.7
4	I	211	PRO	3.7
4	I	136	ASP	3.6

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Mol	Chain	Res	Type	RSRZ
4	I	204	ILE	3.6
4	I	178	ASP	3.6
4	D	164	SER	3.5
4	I	209	PHE	3.5
4	I	206	GLU	3.5
4	D	199	PRO	3.5
5	E	155	GLN	3.4
4	I	212	SER	3.4
4	D	175	TRP	3.3
4	I	210	PHE	3.2
4	I	146	LEU	3.2
4	D	198	PHE	3.2
2	B	99	MET	3.2
5	J	151	ILE	3.1
5	J	259	ALA	3.1
4	D	173	VAL	3.0
2	G	0	MET	3.0
4	D	200	SER	3.0
4	D	191	ILE	3.0
4	D	146	GLN	2.9
5	E	151	ILE	2.8
4	I	46	HIS	2.7
5	E	162	LEU	2.7
4	D	40	GLY	2.7
5	J	155	GLN	2.7
4	D	163	ARG	2.7
4	I	47	GLY	2.7
4	D	183	CYS	2.6
4	I	199	PHE	2.6
5	E	256	TRP	2.5
4	D	195	ASP	2.5
4	I	176	SER	2.5
4	D	160	LEU	2.5
4	I	138	LYS	2.5
5	J	183	HIS	2.5
4	I	124	ILE	2.4
1	F	230	LEU	2.4
5	E	257	GLY	2.4
4	D	187	PHE	2.4
1	F	215	LEU	2.4
5	E	259	ALA	2.4
4	I	95	SER	2.4

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Mol	Chain	Res	Type	RSRZ
4	D	196	THR	2.4
5	E	130	LEU	2.4
4	I	180	LYS	2.3
4	D	184	ALA	2.3
5	E	148	GLU	2.3
4	D	177	ASN	2.3
4	D	181	PHE	2.3
4	I	174	MET	2.3
4	D	112	ILE	2.3
4	D	127	SER	2.3
4	D	131	SER	2.3
5	E	221	ARG	2.3
2	G	1	ILE	2.2
5	J	220	PRO	2.2
4	D	118	ALA	2.2
4	I	190	LYS	2.2
5	E	199	LEU	2.2
4	I	125	GLN	2.2
4	I	48	GLU	2.2
4	D	144	VAL	2.2
4	D	155	THR	2.2
4	D	130	LYS	2.2
5	J	73	LYS	2.1
5	E	179	GLY	2.1
4	I	134	LEU	2.1
5	E	156	LYS	2.1
1	A	227	ASP	2.1
1	F	223	ASP	2.1
4	I	147	PHE	2.1
4	I	158	GLN	2.1
4	I	142	LYS	2.1
2	G	71	THR	2.1
4	I	141	ASP	2.1
4	D	179	SER	2.1
4	D	92	GLY	2.1
1	A	225	THR	2.1
5	J	221	ARG	2.1
5	E	2	ALA	2.0
4	D	128	SER	2.0
1	F	259	CYS	2.0
5	E	198	ALA	2.0
4	I	173	ASP	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.