



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:04 AM GMT

PDB ID : 3CVH
Title : How TCR-like antibody recognizes MHC-bound peptide
Authors : Mareeva, T.; Martinez-Hackert, E.; Sykulev, Y.
Deposited on : 2008-04-18
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

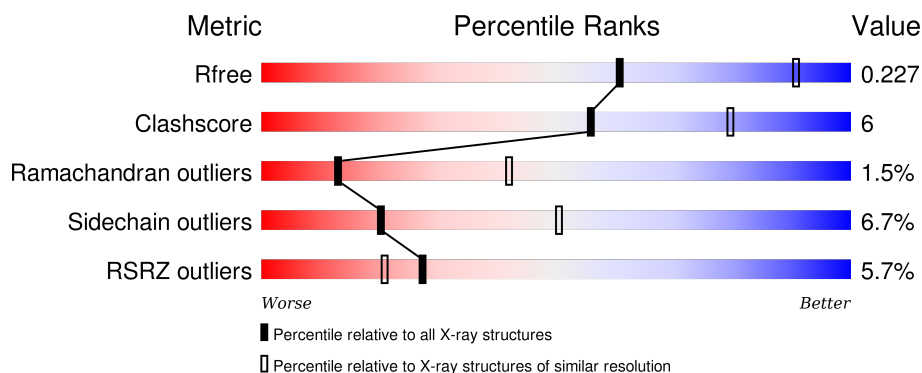
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>0%</div> <div> <div></div> <div>80%</div> <div>18%</div> <div>•</div> </div> </div>
1	M	274	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>13%</div> <div>•</div> </div> </div>
2	B	99	<div> <div></div> <div> <div></div> <div>84%</div> <div>15%</div> <div>•</div> </div> </div>
2	N	99	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>26%</div> <div>•</div> </div> </div>
3	C	8	<div> <div></div> <div> <div></div> <div>75%</div> <div>25%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	O	8	<div><div></div><div>75%</div><div>25%</div></div>
4	H	219	<div><div></div><div>77%</div><div>16%</div><div></div><div></div></div>
4	Q	219	<div><div>13%</div><div></div><div>81%</div><div>13%</div><div></div><div></div></div>
5	L	209	<div><div></div><div>83%</div><div>16%</div><div></div></div>
5	R	209	<div><div>24%</div><div></div><div>73%</div><div>22%</div><div>5%</div></div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12886 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 H-2 class I histocompatibility antigen, K-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2232	1408	393	422	9			
1	M	274	Total	C	N	O	S	0	0	0
			2232	1408	393	422	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
			821	524	138	152	7			
2	N	99	Total	C	N	O	S	0	0	0
			821	524	138	152	7			

- Molecule 3 is a protein called Ovalbumin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	0	0	0
			68	45	10	13			
3	O	8	Total	C	N	O	0	0	0
			68	45	10	13			

- Molecule 4 is a protein called 25-D1.16 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	213	Total	C	N	O	S	0	0	0
			1618	1038	260	313	7			
4	Q	210	Total	C	N	O	S	0	0	0
			1600	1028	257	309	6			

- Molecule 5 is a protein called 25-D1.16 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	L	209	Total 1625	C 1014	N 272	O 334	S 5	0	0	0
5	R	209	Total 1625	C 1014	N 272	O 334	S 5	0	0	0

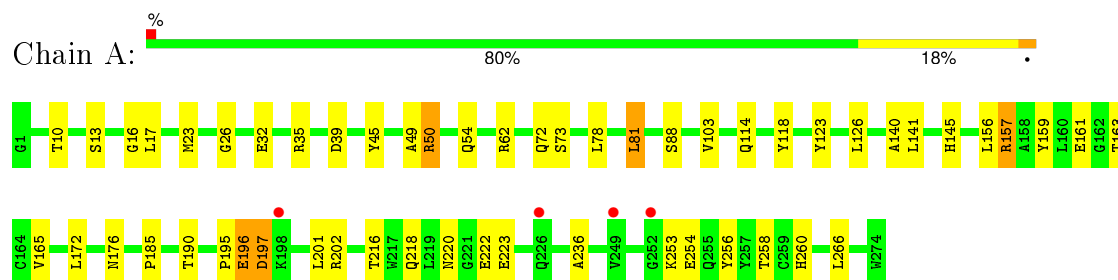
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	42	Total 42	O 42	0	0
6	B	14	Total 14	O 14	0	0
6	C	2	Total 2	O 2	0	0
6	H	39	Total 39	O 39	0	0
6	L	36	Total 36	O 36	0	0
6	M	19	Total 19	O 19	0	0
6	N	8	Total 8	O 8	0	0
6	Q	8	Total 8	O 8	0	0
6	R	8	Total 8	O 8	0	0

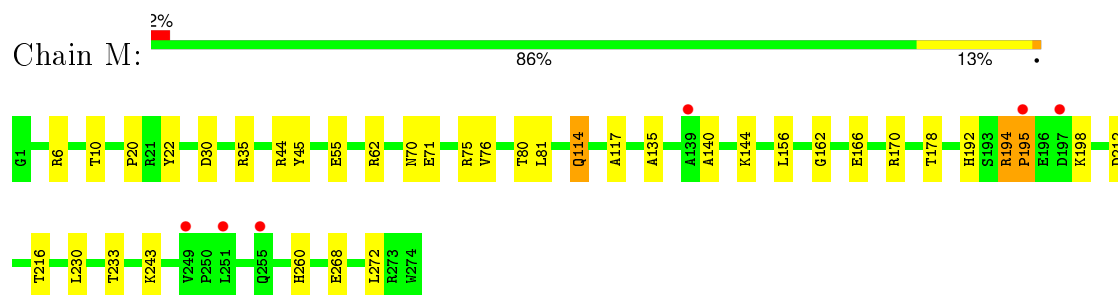
3 Residue-property plots [i](#)

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

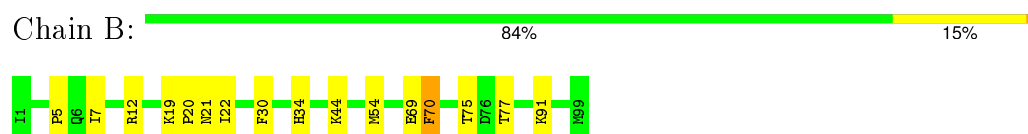
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain



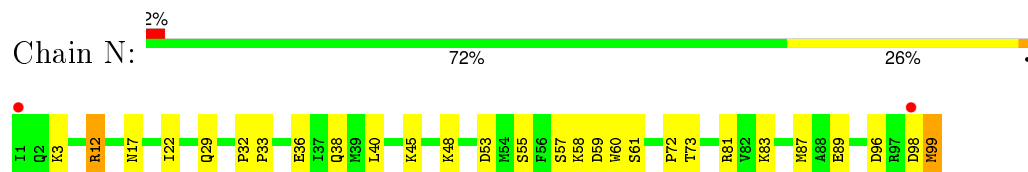
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin

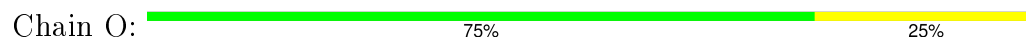


- Molecule 3: Ovalbumin

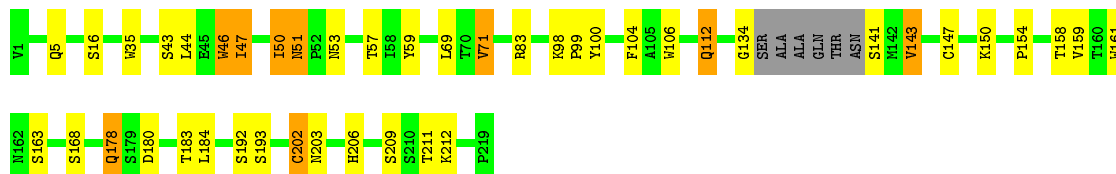
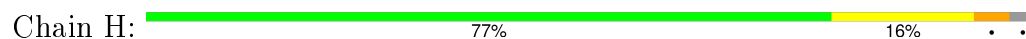




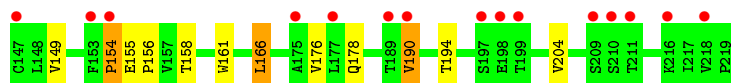
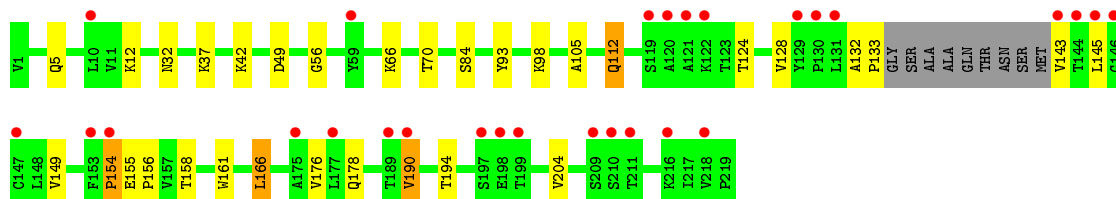
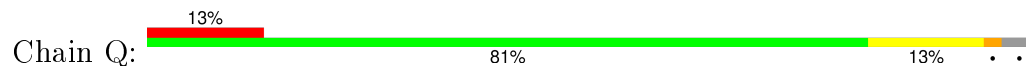
• Molecule 3: Ovalbumin



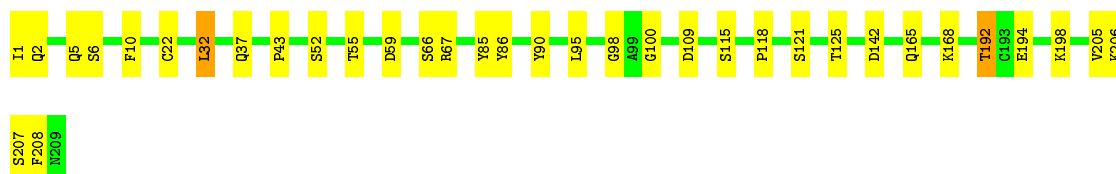
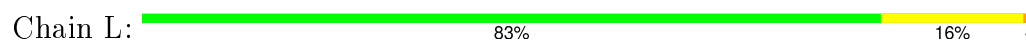
• Molecule 4: 25-D1.16 heavy chain



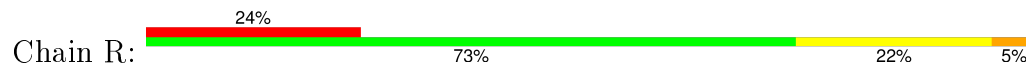
• Molecule 4: 25-D1.16 heavy chain



• Molecule 5: 25-D1.16 light chain



• Molecule 5: 25-D1.16 light chain



G193	
E194	
A195	
T196	
H197	
K198	
T199	
S200	
T201	
S202	
P203	
I204	
V205	
N209	

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.57Å 111.34Å 219.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.98 – 2.90 49.63 – 2.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.98-2.90) 98.0 (49.63-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.27 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.219 , 0.296 0.220 , 0.227	Depositor DCC
R_{free} test set	2189 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	52.0	Xtriage
Anisotropy	0.333	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 60.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 43663 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	12886	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.44	0/2293	0.58	0/3113
1	M	0.39	0/2293	0.52	0/3113
2	B	0.43	0/847	0.52	0/1148
2	N	0.39	0/847	0.53	0/1148
3	C	0.47	0/68	0.74	0/88
3	O	0.44	0/68	0.59	0/88
4	H	0.46	0/1665	0.59	0/2277
4	Q	0.39	0/1647	0.53	0/2254
5	L	0.45	0/1660	0.57	0/2256
5	R	0.38	0/1660	0.52	0/2256
All	All	0.42	0/13048	0.55	0/17741

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2232	0	2123	34	0
1	M	2232	0	2123	19	0
2	B	821	0	796	7	0
2	N	821	0	796	16	0
3	C	68	0	74	1	0
3	O	68	0	74	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	1618	0	1587	39	0
4	Q	1600	0	1570	16	0
5	L	1625	0	1569	12	0
5	R	1625	0	1569	24	0
6	A	42	0	0	5	0
6	B	14	0	0	1	0
6	C	2	0	0	0	0
6	H	39	0	0	7	0
6	L	36	0	0	0	0
6	M	19	0	0	1	0
6	N	8	0	0	0	0
6	Q	8	0	0	0	0
6	R	8	0	0	1	0
All	All	12886	0	12281	156	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 6.

All (156) 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:H:112:GLN:HE21	4:H:112:GLN:H	1.05	0.95
1:A:126:LEU:HD22	1:A:156:LEU:HD23	1.57	0.85
4:H:141:SER:HB3	6:H:222:HOH:O	1.75	0.84
4:H:5:GLN:H	4:H:112:GLN:HE22	1.26	0.81
4:H:134:GLY:HA2	6:H:233:HOH:O	1.79	0.81
1:A:72:GLN:HE22	4:H:59:TYR:H	1.28	0.78
4:H:163:SER:H	4:H:203:ASN:HD21	1.30	0.78
1:M:20:PRO:HG2	1:M:75:ARG:HG2	1.67	0.77
4:H:112:GLN:HE21	4:H:112:GLN:N	1.83	0.75
1:A:72:GLN:NE2	4:H:59:TYR:H	1.85	0.74
1:A:62:ARG:HD3	6:A:297:HOH:O	1.89	0.72
1:A:176:ASN:HB2	6:A:308:HOH:O	1.92	0.69
5:L:121:SER:O	5:L:125:THR:HG23	1.93	0.69
4:H:50:ILE:HG12	4:H:69:LEU:HD23	1.76	0.67
4:H:35:TRP:O	4:H:46:TRP:O	2.12	0.67
5:L:192:THR:HB	5:L:207:SER:OG	1.93	0.66
4:H:5:GLN:H	4:H:112:GLN:NE2	1.93	0.65
2:B:34:HIS:HB2	6:B:104:HOH:O	1.97	0.65
5:R:119:PRO:O	5:R:120:SER:HB2	1.98	0.64
1:A:81:LEU:HD13	3:C:8:LEU:HD13	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:R:186:GLU:HB3	6:R:214:HOH:O	1.99	0.62
1:M:35:ARG:NH1	2:N:53:ASP:HB3	2.15	0.61
5:R:155:GLN:HB2	5:R:158:VAL:HG21	1.83	0.61
4:Q:5:GLN:HB2	4:Q:112:GLN:HE22	1.66	0.61
1:A:35:ARG:NH2	2:B:54:MET:O	2.34	0.60
1:M:44:ARG:HD3	6:M:277:HOH:O	2.01	0.59
6:A:316:HOH:O	5:L:1:ILE:N	2.34	0.59
2:N:12:ARG:NH2	2:N:22:ILE:HD13	2.18	0.59
5:L:194:GLU:HG2	5:L:205:VAL:HG22	1.85	0.59
4:Q:161:TRP:H	4:Q:166:LEU:HD21	1.68	0.58
5:L:5:GLN:HE21	5:L:98:GLY:HA3	1.68	0.56
4:H:16:SER:HB3	4:H:83:ARG:HA	1.86	0.56
5:R:60:ARG:HD3	5:R:76:SER:O	2.05	0.56
1:A:114:GLN:HG2	1:A:156:LEU:HD21	1.86	0.56
1:A:114:GLN:CG	1:A:156:LEU:HD21	2.36	0.56
1:A:236:ALA:HB1	2:B:12:ARG:HG3	1.88	0.56
1:A:10:THR:HG23	1:A:23:MET:HB2	1.87	0.56
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.41	0.56
1:M:162:GLY:O	1:M:166:GLU:HG3	2.06	0.55
2:N:17:ASN:HA	2:N:72:PRO:O	2.07	0.55
4:H:154:PRO:O	4:H:206:HIS:HE1	1.90	0.54
4:Q:128:VAL:HG21	4:Q:204:VAL:HG21	1.89	0.54
1:A:81:LEU:HD23	1:A:118:TYR:CD1	2.43	0.54
5:R:16:ASP:H	5:R:77:LEU:H	1.57	0.54
5:R:135:LEU:HD12	5:R:174:MET:HG2	1.88	0.53
2:N:99:MET:SD	2:N:99:MET:C	2.86	0.53
4:H:50:ILE:HG23	4:H:57:THR:HG22	1.90	0.53
1:A:103:VAL:HG11	1:A:165:VAL:HG22	1.91	0.53
4:H:99:PRO:HG3	4:H:106:TRP:CE2	2.44	0.52
1:M:194:ARG:HB2	1:M:198:LYS:O	2.10	0.52
4:H:134:GLY:CA	6:H:224:HOH:O	2.57	0.52
1:A:172:LEU:O	1:A:176:ASN:HB3	2.10	0.52
4:H:51:ASN:ND2	4:H:53:ASN:H	2.06	0.52
1:M:6:ARG:HH12	2:N:58:LYS:HE2	1.75	0.52
4:Q:166:LEU:H	4:Q:166:LEU:HD22	1.75	0.52
1:M:117:ALA:HB2	2:N:60:TRP:CE2	2.45	0.52
1:M:194:ARG:HG2	1:M:195:PRO:HD2	1.91	0.52
4:Q:143:VAL:HG13	4:Q:190:VAL:HG23	1.92	0.52
5:R:143:ILE:HG23	5:R:197:HIS:ND1	2.25	0.51
1:M:233:THR:HG23	1:M:243:LYS:HB2	1.91	0.51
1:A:145:HIS:HB3	6:A:279:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:R:60:ARG:NE	5:R:78:GLN:HG3	2.25	0.51
4:H:161:TRP:CZ3	4:H:202:CYS:HB2	2.46	0.51
1:M:194:ARG:CB	1:M:198:LYS:O	2.59	0.51
4:H:134:GLY:CA	6:H:233:HOH:O	2.48	0.51
5:R:127:GLY:HA2	5:R:182:LYS:HB2	1.93	0.51
4:H:141:SER:N	4:H:192:SER:HG	2.09	0.50
1:A:13:SER:HB3	1:A:78:LEU:HD13	1.94	0.50
4:H:141:SER:N	6:H:222:HOH:O	2.44	0.50
5:R:124:LEU:O	5:R:182:LYS:HD2	2.12	0.50
4:H:46:TRP:O	4:H:47:ILE:HB	2.13	0.49
1:M:192:HIS:CE1	2:N:98:ASP:HB3	2.48	0.49
4:H:112:GLN:NE2	4:H:112:GLN:H	1.89	0.49
1:A:32:GLU:OE2	1:A:35:ARG:HD2	2.12	0.49
4:H:134:GLY:HA3	6:H:224:HOH:O	2.13	0.48
5:R:53:LEU:HD11	5:R:57:VAL:O	2.12	0.48
4:Q:161:TRP:HB2	4:Q:166:LEU:HD11	1.95	0.48
4:Q:178:GLN:HB3	5:R:159:LEU:HD11	1.96	0.48
4:Q:98:LYS:HD3	4:Q:105:ALA:HB3	1.94	0.48
5:L:37:GLN:NE2	5:L:43:PRO:HD3	2.29	0.48
4:H:43:SER:OG	4:H:44:LEU:N	2.47	0.48
1:A:114:GLN:CB	1:A:156:LEU:HD21	2.43	0.47
1:A:190:THR:HG22	1:A:202:ARG:HB3	1.96	0.47
5:L:5:GLN:HE22	5:L:86:TYR:HA	1.80	0.47
5:R:5:GLN:HE21	5:R:5:GLN:HA	1.79	0.47
5:R:5:GLN:NE2	5:R:87:CYS:SG	2.82	0.46
1:M:30:ASP:OD1	1:M:30:ASP:N	2.48	0.46
4:Q:5:GLN:CB	4:Q:112:GLN:HE22	2.29	0.46
1:M:135:ALA:HB1	1:M:140:ALA:HB3	1.95	0.46
1:M:71:GLU:O	1:M:75:ARG:HG3	2.16	0.46
5:R:184:GLU:HA	5:R:187:ARG:HH21	1.81	0.46
1:A:185:PRO:HD2	1:A:266:LEU:HG	1.97	0.46
1:M:55:GLU:OE1	1:M:170:ARG:NH2	2.49	0.46
1:A:201:LEU:HD21	1:A:254:GLU:HB2	1.98	0.46
4:H:143:VAL:HG13	4:H:143:VAL:O	2.16	0.46
1:A:218:GLN:HG3	1:A:258:THR:OG1	2.16	0.46
5:L:85:TYR:O	5:L:100:GLY:HA2	2.16	0.46
2:N:3:LYS:HE3	2:N:29:GLN:HE21	1.80	0.46
4:H:50:ILE:HG21	4:H:71:VAL:HG22	1.98	0.46
2:N:99:MET:SD	2:N:99:MET:O	2.74	0.46
5:R:160:ASN:HD22	5:R:176:SER:HA	1.81	0.45
1:A:253:LYS:HB3	1:A:256:TYR:CD1	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:32:PRO:HB2	2:N:33:PRO:HD2	1.98	0.45
2:B:7:ILE:HD12	2:B:91:LYS:HD3	1.98	0.45
4:H:46:TRP:O	4:H:47:ILE:CB	2.64	0.45
5:R:162:TRP:NE1	5:R:174:MET:SD	2.90	0.45
1:M:216:THR:OG1	1:M:260:HIS:HB2	2.16	0.45
5:L:22:CYS:SG	5:L:32:LEU:HD11	2.57	0.45
5:R:84:THR:HA	5:R:102:LYS:HA	1.99	0.45
4:Q:128:VAL:HG22	4:Q:149:VAL:HG13	1.98	0.44
5:R:11:SER:HA	5:R:104:GLU:O	2.18	0.44
4:H:178:GLN:HE21	4:H:183:THR:HG1	1.61	0.44
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.99	0.44
1:A:222:GLU:HA	6:A:313:HOH:O	2.17	0.44
4:H:150:LYS:HG3	4:H:178:GLN:NE2	2.32	0.44
1:A:26:GLY:O	1:A:32:GLU:HA	2.18	0.44
1:A:49:ALA:O	1:A:50:ARG:C	2.56	0.44
4:H:5:GLN:N	4:H:112:GLN:HE22	2.04	0.44
4:H:98:LYS:HB2	4:H:100:TYR:O	2.17	0.43
1:A:216:THR:OG1	1:A:260:HIS:HB2	2.18	0.43
1:A:123:TYR:CZ	1:A:140:ALA:HA	2.54	0.43
2:N:38:GLN:HE22	2:N:45:LYS:HE2	1.83	0.43
4:Q:155:GLU:HB3	4:Q:156:PRO:HA	2.01	0.43
4:H:141:SER:CB	6:H:222:HOH:O	2.49	0.43
5:R:45:LEU:CD2	5:R:54:GLU:HG3	2.49	0.43
2:N:29:GLN:HA	2:N:61:SER:HB2	2.01	0.43
1:A:196:GLU:HB2	1:A:197:ASP:H	1.73	0.43
5:L:109:ASP:OD2	5:L:198:LYS:HE3	2.19	0.42
4:H:209:SER:O	4:H:211:THR:HG23	2.19	0.42
2:N:59:ASP:O	2:N:60:TRP:HB2	2.19	0.42
4:H:104:PHE:HB3	5:L:90:TYR:O	2.19	0.42
1:M:76:VAL:HG13	4:Q:56:GLY:HA3	2.01	0.42
1:A:159:TYR:CD1	1:A:163:THR:HB	2.55	0.42
5:R:2:GLN:HE21	5:R:2:GLN:HB3	1.68	0.42
4:H:16:SER:CB	4:H:83:ARG:HA	2.49	0.42
4:H:184:LEU:HD12	4:H:184:LEU:C	2.40	0.42
1:A:157:ARG:HE	1:A:161:GLU:CD	2.23	0.42
3:O:7:LYS:HE2	4:Q:32:ASN:ND2	2.34	0.42
5:R:149:ILE:HD11	5:R:178:LEU:HD21	2.02	0.41
1:A:32:GLU:CD	1:A:35:ARG:HH11	2.22	0.41
2:N:36:GLU:HB3	2:N:83:LYS:HB2	2.03	0.41
4:H:163:SER:H	4:H:203:ASN:ND2	2.09	0.41
1:M:22:TYR:OH	1:M:70:ASN:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:Q:154:PRO:HB2	4:Q:155:GLU:H	1.72	0.41
2:N:40:LEU:HD11	2:N:81:ARG:HD3	2.02	0.41
1:A:16:GLY:O	1:A:17:LEU:HD23	2.21	0.41
2:B:19:LYS:HA	2:B:20:PRO:HD3	1.97	0.41
4:H:159:VAL:HA	4:H:203:ASN:O	2.20	0.41
2:N:96:ASP:O	2:N:99:MET:OXT	2.38	0.41
1:M:114:GLN:HE21	1:M:114:GLN:HA	1.86	0.41
5:R:105:LEU:HD12	5:R:165:GLN:HG2	2.03	0.40
5:L:118:PRO:HB3	5:L:208:PHE:CE1	2.56	0.40
5:R:5:GLN:OE1	5:R:86:TYR:HA	2.21	0.40
1:A:218:GLN:HA	1:A:223:GLU:HA	2.02	0.40
4:Q:37:LYS:HG3	4:Q:93:TYR:CE1	2.56	0.40
4:Q:132:ALA:HA	4:Q:133:PRO:HD3	1.98	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	272/274 (99%)	251 (92%)	15 (6%)	6 (2%)	8	31
1	M	272/274 (99%)	258 (95%)	13 (5%)	1 (0%)	39	74
2	B	97/99 (98%)	93 (96%)	4 (4%)	0	100	100
2	N	97/99 (98%)	90 (93%)	7 (7%)	0	100	100
3	C	6/8 (75%)	6 (100%)	0	0	100	100
3	O	6/8 (75%)	5 (83%)	1 (17%)	0	100	100
4	H	209/219 (95%)	191 (91%)	14 (7%)	4 (2%)	10	35
4	Q	206/219 (94%)	185 (90%)	19 (9%)	2 (1%)	19	54
5	L	207/209 (99%)	199 (96%)	6 (3%)	2 (1%)	19	54
5	R	207/209 (99%)	180 (87%)	19 (9%)	8 (4%)	4	15

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1579/1618 (98%)	1458 (92%)	98 (6%)	23 (2%)	13	42

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	H	47	ILE
4	H	143	VAL
4	H	178	GLN
5	R	120	SER
5	R	136	ASN
1	A	50	ARG
5	R	81	ASP
5	R	119	PRO
5	L	6	SER
5	L	67	ARG
4	Q	84	SER
5	R	6	SER
1	A	54	GLN
1	A	220	ASN
5	R	67	ARG
1	A	195	PRO
1	A	196	GLU
1	A	197	ASP
4	H	46	TRP
5	R	7	SER
5	R	185	TYR
1	M	195	PRO
4	Q	154	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	232/232 (100%)	225 (97%)	7 (3%)	48	83
1	M	232/232 (100%)	218 (94%)	14 (6%)	24	57

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	94/94 (100%)	88 (94%)	6 (6%)	22	53
2	N	94/94 (100%)	86 (92%)	8 (8%)	13	37
3	C	8/8 (100%)	7 (88%)	1 (12%)	6	17
3	O	8/8 (100%)	7 (88%)	1 (12%)	6	17
4	H	182/186 (98%)	171 (94%)	11 (6%)	24	57
4	Q	180/186 (97%)	167 (93%)	13 (7%)	18	46
5	L	186/186 (100%)	172 (92%)	14 (8%)	17	44
5	R	186/186 (100%)	167 (90%)	19 (10%)	9	27
All	All	1402/1412 (99%)	1308 (93%)	94 (7%)	20	50

All (94) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ASP
1	A	45	TYR
1	A	73	SER
1	A	81	LEU
1	A	88	SER
1	A	141	LEU
1	A	157	ARG
2	B	22	ILE
2	B	44	LYS
2	B	69	GLU
2	B	70	PHE
2	B	75	THR
2	B	77	THR
3	C	7	LYS
4	H	50	ILE
4	H	51	ASN
4	H	71	VAL
4	H	112	GLN
4	H	147	CYS
4	H	158	THR
4	H	168	SER
4	H	180	ASP
4	H	193	SER
4	H	202	CYS
4	H	212	LYS
5	L	2	GLN

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Mol	Chain	Res	Type
5	L	10	PHE
5	L	32	LEU
5	L	52	SER
5	L	55	THR
5	L	59	ASP
5	L	66	SER
5	L	95	LEU
5	L	115	SER
5	L	142	ASP
5	L	165	GLN
5	L	168	LYS
5	L	192	THR
5	L	206	LYS
1	M	10	THR
1	M	45	TYR
1	M	62	ARG
1	M	80	THR
1	M	81	LEU
1	M	114	GLN
1	M	144	LYS
1	M	156	LEU
1	M	178	THR
1	M	194	ARG
1	M	212	ASP
1	M	230	LEU
1	M	268	GLU
1	M	272	LEU
2	N	12	ARG
2	N	48	LYS
2	N	55	SER
2	N	57	SER
2	N	73	THR
2	N	87	MET
2	N	89	GLU
2	N	99	MET
3	O	4	ASN
4	Q	12	LYS
4	Q	42	LYS
4	Q	49	ASP
4	Q	66	LYS
4	Q	70	THR
4	Q	112	GLN

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Mol	Chain	Res	Type
4	Q	124	THR
4	Q	145	LEU
4	Q	158	THR
4	Q	166	LEU
4	Q	176	VAL
4	Q	190	VAL
4	Q	194	THR
5	R	2	GLN
5	R	5	GLN
5	R	12	VAL
5	R	32	LEU
5	R	47	ILE
5	R	55	THR
5	R	71	THR
5	R	77	LEU
5	R	78	GLN
5	R	104	GLU
5	R	109	ASP
5	R	131	VAL
5	R	143	ILE
5	R	154	ARG
5	R	174	MET
5	R	183	ASP
5	R	187	ARG
5	R	198	LYS
5	R	209	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	226	GLN
1	A	255	GLN
4	H	38	GLN
4	H	51	ASN
4	H	112	GLN
4	H	203	ASN
4	H	206	HIS
5	L	2	GLN
5	L	5	GLN
5	L	37	GLN
5	L	78	GLN

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Mol	Chain	Res	Type
5	L	123	GLN
5	L	137	ASN
5	L	160	ASN
1	M	87	GLN
1	M	114	GLN
1	M	174	ASN
2	N	38	GLN
4	Q	61	GLN
4	Q	112	GLN
5	R	2	GLN
5	R	36	GLN
5	R	78	GLN
5	R	160	ASN
5	R	189	ASN
5	R	209	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 carbohydrates 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%)	0.14	4 (1%) 76 74	24, 35, 53, 59	0
1	M	274/274 (100%)	0.24	6 (2%) 65 60	35, 46, 54, 62	0
2	B	99/99 (100%)	-0.14	0 100 100	27, 36, 43, 45	0
2	N	99/99 (100%)	0.20	2 (2%) 68 64	34, 45, 53, 63	0
3	C	8/8 (100%)	-0.00	0 100 100	25, 26, 28, 29	0
3	O	8/8 (100%)	0.22	0 100 100	47, 50, 51, 51	0
4	H	213/219 (97%)	-0.04	0 100 100	18, 28, 37, 43	0
4	Q	210/219 (95%)	0.81	28 (13%) 4 2	47, 56, 77, 89	0
5	L	209/209 (100%)	-0.02	0 100 100	17, 30, 41, 51	0
5	R	209/209 (100%)	1.32	51 (24%) 1 0	51, 60, 67, 70	0
All	All	1603/1618 (99%)	0.34	91 (5%) 27 21	17, 43, 65, 89	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	R	112	PRO	7.5
5	R	179	THR	6.7
5	R	191	TYR	6.5
5	R	209	ASN	6.2
4	Q	119	SER	6.0
5	R	178	LEU	5.9
1	M	197	ASP	5.5
5	R	172	TYR	5.4
5	R	144	ASN	5.2
5	R	204	ILE	5.2
4	Q	143	VAL	4.8
2	N	1	ILE	4.4
1	A	226	GLN	4.3

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Mol	Chain	Res	Type	RSRZ
4	Q	216	LYS	4.3
5	R	130	SER	4.1
5	R	151	GLY	4.1
4	Q	189	THR	4.1
5	R	192	THR	4.0
5	R	194	GLU	3.8
5	R	162	TRP	3.8
5	R	188	HIS	3.8
1	A	198	LYS	3.8
5	R	189	ASN	3.6
4	Q	218	VAL	3.5
4	Q	199	THR	3.5
5	R	205	VAL	3.4
5	R	202	SER	3.4
4	Q	211	THR	3.4
4	Q	129	TYR	3.4
5	R	187	ARG	3.4
5	R	105	LEU	3.3
5	R	195	ALA	3.3
5	R	154	ARG	3.2
5	R	134	PHE	3.2
1	M	195	PRO	3.1
2	N	98	ASP	3.1
5	R	177	THR	3.1
5	R	7	SER	3.1
4	Q	120	ALA	3.0
5	R	149	ILE	3.0
5	R	114	VAL	3.0
5	R	150	ASP	2.9
1	M	251	LEU	2.9
4	Q	131	LEU	2.9
5	R	152	SER	2.9
1	M	249	VAL	2.8
5	R	142	ASP	2.8
5	R	200	SER	2.8
5	R	74	ILE	2.8
5	R	86	TYR	2.7
5	R	155	GLN	2.7
4	Q	210	SER	2.7
4	Q	154	PRO	2.7
5	R	153	GLU	2.6
5	R	198	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
4	Q	197	SER	2.6
5	R	145	VAL	2.6
5	R	157	GLY	2.6
4	Q	121	ALA	2.6
5	R	143	ILE	2.5
5	R	109	ASP	2.5
5	R	9	SER	2.4
4	Q	59	TYR	2.4
5	R	197	HIS	2.4
4	Q	122	LYS	2.4
5	R	141	LYS	2.4
5	R	20	ILE	2.4
5	R	186	GLU	2.4
1	A	252	GLY	2.4
4	Q	198	GLU	2.4
5	R	73	ILE	2.4
5	R	10	PHE	2.3
4	Q	145	LEU	2.3
1	M	255	GLN	2.3
5	R	13	SER	2.3
4	Q	177	LEU	2.3
4	Q	190	VAL	2.3
4	Q	153	PHE	2.2
4	Q	175	ALA	2.2
1	A	249	VAL	2.2
5	R	158	VAL	2.2
4	Q	130	PRO	2.2
5	R	77	LEU	2.2
1	M	139	ALA	2.1
4	Q	144	THR	2.1
4	Q	10	LEU	2.1
5	R	18	VAL	2.1
4	Q	209	SER	2.0
4	Q	146	GLY	2.0
4	Q	147	CYS	2.0
5	R	129	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.