



# Full wwPDB X-ray Structure Validation Report

Jun 20, 2014 – 07:49 AM EDT

PDB ID : 4P5T  
Title : 14.C6 TCR complexed with MHC class II I-Ab/3K peptide  
Authors : Trenh, P.; Stadinski, B.; Huseby, E.S.; Stern, L.J.  
Deposited on : 2014-03-19  
Resolution : 3.26 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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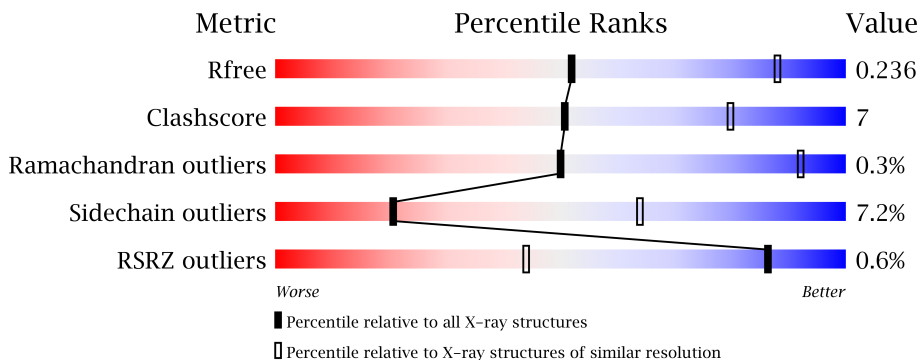
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : **FAILED**  
Xtriage (Phenix) : dev-1439  
EDS : stable23161  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable23161

# 1 Overall quality at a glance

The reported resolution of this entry is 3.26 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))
$R_{free}$	66092	1085 (3.32-3.20)
Clashscore	79885	1374 (3.32-3.20)
Ramachandran outliers	78287	1348 (3.32-3.20)
Sidechain outliers	78261	1346 (3.32-3.20)
RSRZ outliers	66119	1086 (3.32-3.20)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	206	
1	E	206	
2	B	241	
2	F	241	
3	C	182	
3	G	182	
4	D	217	
4	H	217	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12366 atoms, of which 0 are hydrogen and 0 are deuterium.

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 TRA protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	197	Total	C	N	O	S	0	0	0
			1488	948	241	292	7			
1	E	188	Total	C	N	O	S	0	0	0
			1403	902	223	272	6			

- Molecule 2 is a protein called Human nkt tcr beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	238	Total	C	N	O	S	0	0	0
			1815	1146	318	345	6			
2	F	229	Total	C	N	O	S	0	0	0
			1725	1091	300	328	6			

- Molecule 3 is a protein called H-2 class II histocompatibility antigen, A-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	179	Total	C	N	O	S	0	0	0
			1381	897	215	266	3			
3	G	179	Total	C	N	O	S	0	0	0
			1377	895	214	265	3			

- Molecule 4 is a protein called protein of 3K peptide (FEAQKAKANKAVD), Linkerregion - GGGGSLVPRGSGGGG, H-2 class II histocompatibility antigen, A beta chain, H-2 class II histocompatibility antigen, A beta chain.

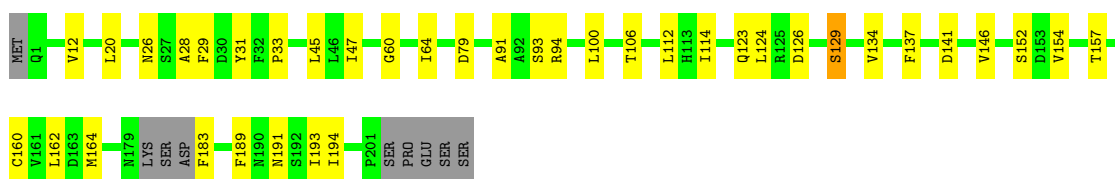
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	199	Total	C	N	O	S	0	0	0
			1593	1001	279	307	6			
4	H	199	Total	C	N	O	S	0	0	0
			1584	997	275	306	6			

### 3 Residue-property plots

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.

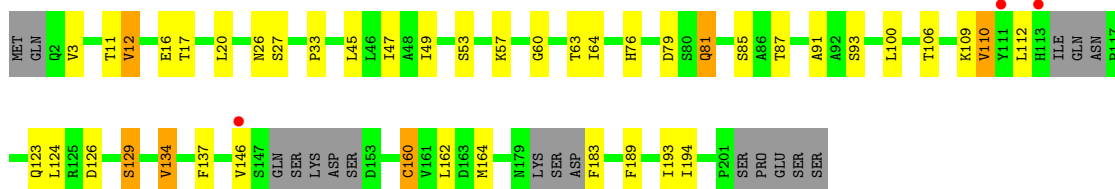
- Molecule 1: TRA protein

Chain A:



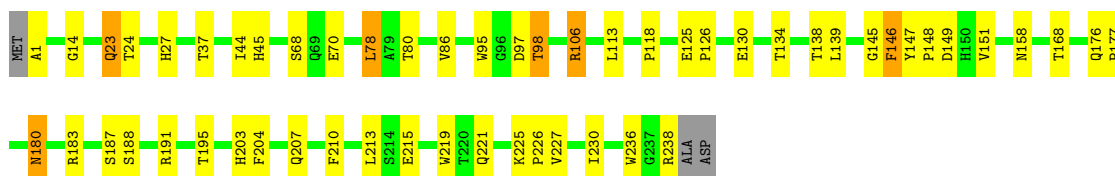
- Molecule 1: TRA protein

Chain E:



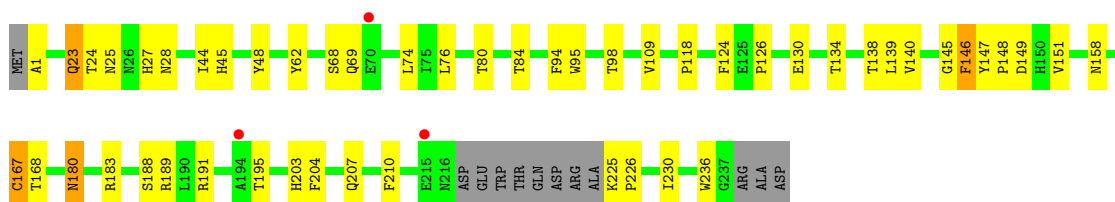
- Molecule 2: Human nkt tcr beta chain

Chain B:



- Molecule 2: Human nkt tcr beta chain

Chain F:



- Molecule 3: H-2 class II histocompatibility antigen, A-B alpha chain

GLU

- |    |    |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |      |      |      |      |      |      |      |      |      |      |      |      |      |      |     |     |     |
|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|------|------|------|------|------|-----|-----|-----|
| IO | G9 | Y13 | Q14 | D18 | I19 | D29 | V34 | D36 | L35 | T41 | V42 | F54 | V65 | V66 | L70 | A86 | P96 | L99 | G100 | Q101 | P102 | N103 | T104 | T110 | N111 | V116 | T120 | P155 | K164 | P173 | H177 | V178 | GLU | PRO | CTD |
|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|------|------|------|------|------|-----|-----|-----|

- |      |      |      |      |      |      |      |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |    |    |    |    |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |
|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| F-25 | Q-22 | K-21 | A-20 | K-19 | V-14 | D-13 | GLY | GLY | GLY | GLY | GLY | LEU | VAL | PRO | ARG | ARG | GLY | GLY | GLY | GLY | GLY | GLY | SER | GLU | ARG | H6 | F7 | V8 | Y9 | M12 | Y16 | R25 | T28 | V38 | R39 | Y40 | E46 | H47 | R48 | E74 | D75 | L76 | R80 | Y83 | T88 | L92 |
|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|----|----|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|

F-25	K-21	A-20	K-19	D-13	GLY	GLY	GLY	GLY	SER	LEU	VAL	PRO	ARG	GLY	SER	GLY	GLY	GLY	SER	GLU	ARG	H6	F7	V8	Y9	Q10	F11	M12	G13	E14	T28	R29	Y30	I31	Y37	V38	R39	Y40	R70	T71	R72	A73	E74	T77	Y83	T88	H89	T90	S91	L92
------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	----	----	----	----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

L95
S105
R106
T107
S108
A109
L110
M111
H112
H113
M114
T115
T121
D122
A126
V130
F133
R134
M135
L148
M154
T155
M164
T165
P166
R167
E170
E171
Y172
T173
S180
L181
L185
E188
L193

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.01Å 74.14Å 259.29Å 90.00° 92.00° 90.00°	Depositor
Resolution (Å)	47.05 – 3.26 48.78 – 3.26	Depositor EDS
% Data completeness (in resolution range)	99.0 (47.05-3.26) 99.0 (48.78-3.26)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.42 (at 3.25Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.186 , 0.236 0.186 , 0.236	Depositor DCC
$R_{free}$ test set	1957 reflections (5.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.4	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 43.1	EDS
Estimated twinning fraction	0.024 for h,-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 39092 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	12366	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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	0.26	0/1525	0.40	0/2078
1	E	0.24	0/1438	0.38	0/1960
2	B	0.25	0/1868	0.44	0/2558
2	F	0.23	0/1775	0.41	0/2432
3	C	0.25	0/1425	0.40	0/1957
3	G	0.22	0/1420	0.38	0/1949
4	D	0.24	0/1632	0.40	0/2225
4	H	0.24	0/1623	0.40	0/2214
All	All	0.24	0/12706	0.40	0/17373

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1488	0	1353	18	0
1	E	1403	0	1255	17	0
2	B	1815	0	1654	36	0
2	F	1725	0	1561	41	0
3	C	1381	0	1263	18	0
3	G	1377	0	1256	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1593	0	1471	30	0
4	H	1584	0	1456	28	0
All	All	12366	0	11269	174	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (174) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:84:THR:HG22	2:F:109:VAL:H	1.46	0.80
1:E:45:LEU:HD22	2:F:98:THR:HG22	1.69	0.74
2:B:130:GLU:OE2	2:B:191:ARG:NH1	2.20	0.74
2:F:130:GLU:OE2	2:F:191:ARG:NH1	2.20	0.74
2:F:145:GLY:HA2	2:F:183:ARG:HD3	1.71	0.73
2:B:145:GLY:HA2	2:B:183:ARG:HD3	1.72	0.71
3:G:34:VAL:HG22	3:G:41:THR:HG22	1.74	0.69
3:C:139:VAL:HG11	4:D:12:MET:HE1	1.75	0.68
1:E:49:ILE:HD13	1:E:57:LYS:HB2	1.75	0.68
4:H:14:GLU:OE1	4:H:29:ARG:NH1	2.26	0.68
3:C:118:ASN:HB2	3:C:166:GLU:HB2	1.79	0.64
4:D:46:GLU:OE2	4:D:48:ARG:NH1	2.30	0.64
4:D:114:ASN:HB3	4:D:164:MET:HE1	1.80	0.62
1:E:93:SER:HB2	1:E:100:LEU:HD23	1.82	0.61
3:C:14:GLN:HG2	4:D:8:VAL:HG22	1.82	0.61
4:H:113:HIS:H	4:H:165:THR:HG23	1.66	0.61
4:H:111:ASN:OD1	4:H:167:ARG:NH1	2.34	0.60
4:D:95:LEU:HG	4:D:180:SER:HA	1.83	0.60
4:D:100:VAL:HG21	4:D:176:VAL:HG11	1.84	0.60
3:G:103:ASN:OD1	3:G:104:THR:N	2.34	0.59
3:G:13:TYR:OH	3:G:18:ASP:OD1	2.20	0.59
2:B:1:ALA:H3	2:B:24:THR:HB	1.68	0.58
2:F:180:ASN:N	2:F:180:ASN:OD1	2.34	0.57
1:A:20:LEU:HD22	1:A:106:THR:HG21	1.87	0.56
1:A:94:ARG:NH2	3:C:55:ASP:OD2	2.38	0.56
4:H:37:TYR:CD2	4:H:38:VAL:HG23	2.40	0.56
3:C:70:LEU:HD13	4:D:9:TYR:HB2	1.87	0.56
4:D:76:ASP:OD1	4:D:80:ARG:NH1	2.39	0.56
4:H:114:ASN:HB3	4:H:164:MET:HE1	1.87	0.55
1:A:114:ILE:HG13	1:A:141:ASP:HA	1.88	0.55
2:F:23:GLN:HG2	2:F:25:ASN:H	1.71	0.55
2:B:68:SER:OG	2:B:70:GLU:OE1	2.25	0.55
1:E:49:ILE:HD11	1:E:53:SER:HB2	1.88	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:225:LYS:HG2	2:B:227:VAL:HG13	1.88	0.55
2:F:158:ASN:HA	2:F:203:HIS:CE1	2.42	0.55
2:F:28:ASN:ND2	2:F:48:TYR:O	2.40	0.54
1:A:28:ALA:O	4:D:-22:GLN:NE2	2.41	0.54
1:E:134:VAL:HB	2:F:124:PHE:CD2	2.41	0.54
1:A:33:PRO:HG2	1:A:91:ALA:HB3	1.89	0.53
2:B:158:ASN:HA	2:B:203:HIS:CE1	2.42	0.53
2:B:95:TRP:CE2	4:D:-19:LYS:HB2	2.43	0.53
2:F:1:ALA:N	2:F:25:ASN:OD1	2.43	0.52
2:B:180:ASN:N	2:B:180:ASN:OD1	2.36	0.52
4:H:10:GLN:HB3	4:H:31:ILE:HB	1.91	0.52
3:G:110:ASP:OD1	3:G:111:ASN:N	2.39	0.51
3:G:41:THR:HG21	3:G:54:PHE:HD2	1.75	0.51
1:E:47:ILE:HG23	1:E:64:ILE:HD11	1.92	0.51
2:B:23:GLN:OE1	2:B:27:HIS:N	2.42	0.51
2:B:147:TYR:HD2	2:B:148:PRO:HA	1.76	0.51
2:F:147:TYR:CD2	2:F:148:PRO:HA	2.46	0.51
4:D:28:THR:HB	4:D:40:TYR:HB3	1.93	0.51
4:H:173:THR:HG23	4:H:188:GLU:HG2	1.92	0.51
2:F:147:TYR:HD2	2:F:148:PRO:HA	1.76	0.50
2:F:68:SER:OG	2:F:69:GLN:N	2.44	0.50
1:A:152:SER:C	1:A:154:VAL:H	2.14	0.50
2:B:147:TYR:CD2	2:B:148:PRO:HA	2.46	0.50
3:C:110:ASP:OD1	3:C:111:ASN:N	2.38	0.50
4:D:121:THR:HG22	4:D:157:GLN:HG3	1.93	0.50
4:H:122:ASP:HA	4:H:155:THR:HB	1.93	0.50
4:D:122:ASP:HA	4:D:155:THR:HB	1.93	0.50
1:E:162:LEU:HB3	2:F:167:CYS:HB3	1.94	0.49
1:E:126:ASP:HA	2:F:124:PHE:HD1	1.77	0.49
2:B:146:PHE:HE1	2:B:151:VAL:HG21	1.78	0.49
2:B:146:PHE:H	2:B:146:PHE:HD2	1.60	0.49
2:F:94:PHE:CG	3:G:65:VAL:HG22	2.47	0.49
2:F:146:PHE:HE1	2:F:151:VAL:HG21	1.78	0.48
2:F:203:HIS:HB3	2:F:236:TRP:CE3	2.47	0.48
2:B:203:HIS:HB3	2:B:236:TRP:CE3	2.47	0.48
3:C:165:VAL:HB	3:C:174:VAL:HG12	1.95	0.48
2:F:27:HIS:O	2:F:69:GLN:NE2	2.46	0.48
2:B:86:VAL:HG22	2:B:106:ARG:HD3	1.95	0.48
4:D:16:TYR:HB2	4:D:25:ARG:HB3	1.96	0.48
2:F:146:PHE:H	2:F:146:PHE:HD2	1.60	0.48
1:A:47:ILE:HG23	1:A:64:ILE:HD11	1.96	0.48
1:E:126:ASP:HB3	1:E:129:SER:O	2.14	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:93:SER:HB2	1:A:100:LEU:HD23	1.95	0.47
3:C:135:THR:O	3:C:147:LYS:HE3	2.14	0.47
4:D:176:VAL:HG13	4:D:185:ILE:HB	1.95	0.47
2:B:44:ILE:HG22	2:B:45:HIS:CD2	2.48	0.47
2:F:23:GLN:HG2	2:F:24:THR:N	2.30	0.47
2:B:168:THR:HG23	2:B:188:SER:HB2	1.97	0.47
2:F:168:THR:HG23	2:F:188:SER:HB2	1.96	0.47
2:F:158:ASN:HA	2:F:203:HIS:HE1	1.81	0.46
4:H:167:ARG:O	4:H:170:GLU:HG2	2.15	0.46
1:E:20:LEU:HD22	1:E:106:THR:HG21	1.97	0.46
2:F:23:GLN:OE1	2:F:27:HIS:N	2.40	0.46
3:C:66:VAL:HG13	4:D:9:TYR:CD2	2.51	0.46
1:A:126:ASP:HB3	1:A:129:SER:O	2.16	0.46
2:B:158:ASN:HA	2:B:203:HIS:HE1	1.81	0.46
3:C:96:PRO:HD3	4:D:121:THR:HG21	1.98	0.46
3:C:82:ALA:HB1	3:C:113:PHE:CE2	2.51	0.46
4:D:-19:LYS:HG2	4:D:74:GLU:OE1	2.16	0.46
4:D:181:LEU:HD22	4:D:185:ILE:HG13	1.97	0.45
4:H:105:SER:OG	4:H:115:THR:HG23	2.15	0.45
2:F:23:GLN:HB3	2:F:23:GLN:HE21	1.52	0.45
1:A:137:PHE:HB2	1:A:189:PHE:CE2	2.50	0.45
1:A:29:PHE:HD1	1:A:94:ARG:HG2	1.81	0.45
1:E:137:PHE:HB2	1:E:189:PHE:CE2	2.51	0.45
2:F:28:ASN:HA	2:F:69:GLN:HE22	1.82	0.45
2:B:207:GLN:HG3	2:B:230:ILE:HG23	1.99	0.45
4:H:-19:LYS:HG2	4:H:74:GLU:OE1	2.17	0.45
3:G:66:VAL:HG13	4:H:9:TYR:CD2	2.52	0.45
3:G:14:GLN:HG2	3:G:19:ILE:HD12	1.99	0.45
3:C:57:GLN:HA	3:C:60:LEU:HB2	1.98	0.45
2:F:138:THR:HG1	2:F:191:ARG:HH11	1.64	0.45
1:A:31:TYR:OH	2:B:97:ASP:HB2	2.17	0.44
2:F:151:VAL:HG12	2:F:210:PHE:HA	1.99	0.44
4:H:83:TYR:O	4:H:88:THR:HG23	2.17	0.44
2:B:118:PRO:HD3	2:B:226:PRO:HB3	2.00	0.44
2:F:118:PRO:HD3	2:F:226:PRO:HB3	1.99	0.44
2:F:207:GLN:HG3	2:F:230:ILE:HG23	1.99	0.44
1:A:29:PHE:CD1	1:A:94:ARG:HG2	2.53	0.44
2:B:219:TRP:CE2	2:B:221:GLN:HB2	2.52	0.44
2:F:149:ASP:N	2:F:149:ASP:OD1	2.50	0.44
2:B:151:VAL:HG12	2:B:210:PHE:HA	2.00	0.44
2:F:225:LYS:HA	2:F:226:PRO:HD3	1.77	0.43
4:D:95:LEU:HA	4:D:95:LEU:HD12	1.83	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:H:73:ALA:O	4:H:77:THR:OG1	2.37	0.43
2:B:23:GLN:HE21	2:B:23:GLN:HB3	1.62	0.43
3:C:103:ASN:OD1	3:C:104:THR:N	2.45	0.43
2:B:149:ASP:OD1	2:B:149:ASP:N	2.50	0.43
2:B:14:GLY:HA2	2:B:78:LEU:HD23	2.01	0.43
3:G:99:LEU:HA	3:G:155:PRO:HB2	1.99	0.43
3:C:76:ARG:NH2	4:D:-14:VAL:O	2.52	0.43
4:H:28:THR:HG22	4:H:40:TYR:HB3	2.01	0.43
1:A:162:LEU:HD21	2:B:191:ARG:HD3	2.01	0.43
1:E:81:GLN:O	1:E:110:VAL:HG21	2.19	0.43
3:G:70:LEU:HD13	4:H:9:TYR:HB2	2.01	0.42
3:G:29:ASP:HB3	4:H:154:TRP:CE2	2.54	0.42
4:H:135:ASN:ND2	4:H:171:VAL:H	2.17	0.42
4:D:98:PRO:HB3	4:D:123:PHE:HB3	2.01	0.42
1:E:126:ASP:HA	2:F:124:PHE:CD1	2.54	0.42
2:F:62:TYR:HD1	2:F:74:LEU:HD21	1.84	0.42
3:C:29:ASP:HB3	4:D:154:TRP:CE2	2.55	0.42
4:H:181:LEU:HD22	4:H:185:ILE:HG13	2.01	0.42
4:D:183:SER:HA	4:D:184:PRO:HD3	1.89	0.42
3:C:59:GLY:O	3:C:63:ILE:HG12	2.20	0.42
2:F:44:ILE:HG22	2:F:45:HIS:CD2	2.55	0.42
2:B:125:GLU:HA	2:B:126:PRO:HD3	1.96	0.41
1:A:45:LEU:HD22	2:B:98:THR:HG22	2.01	0.41
2:F:140:VAL:HG22	2:F:189:ARG:HG3	2.02	0.41
3:G:96:PRO:HD3	4:H:121:THR:HG21	2.02	0.41
1:A:157:THR:HG21	2:B:187:SER:OG	2.20	0.41
2:B:95:TRP:CD2	4:D:-19:LYS:HB2	2.55	0.41
3:G:9:GLY:O	4:H:12:MET:HA	2.20	0.41
4:H:88:THR:HA	4:H:92:LEU:HB2	2.02	0.41
1:A:152:SER:O	1:A:154:VAL:N	2.51	0.41
3:G:101:GLN:HA	3:G:102:PRO:HD2	1.83	0.41
2:F:146:PHE:N	2:F:146:PHE:CD2	2.88	0.41
2:F:76:LEU:HA	2:F:76:LEU:HD12	1.83	0.41
1:E:26:ASN:OD1	1:E:27:SER:N	2.54	0.41
4:D:12:MET:HB3	4:D:12:MET:HE3	1.92	0.41
4:D:83:TYR:O	4:D:88:THR:HG23	2.21	0.41
3:C:13:TYR:OH	3:C:18:ASP:OD1	2.22	0.41
2:F:126:PRO:HD3	2:F:139:LEU:HG	2.02	0.41
1:A:26:ASN:HB3	1:A:29:PHE:CE2	2.56	0.41
2:B:146:PHE:N	2:B:146:PHE:CD2	2.88	0.41
4:D:92:LEU:HA	4:D:92:LEU:HD12	1.88	0.41
4:H:109:ALA:O	4:H:112:HIS:HB2	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:H:126:ALA:HB1	4:H:148:LEU:HD21	2.03	0.41
2:B:138:THR:OG1	2:B:191:ARG:HD2	2.21	0.41
1:E:12:VAL:HG13	1:E:16:GLU:HB2	2.03	0.41
1:E:33:PRO:HG2	1:E:91:ALA:HB3	2.03	0.41
2:B:126:PRO:HD3	2:B:139:LEU:HG	2.03	0.40
2:B:176:GLN:HA	2:B:177:PRO:HD2	1.94	0.40
4:D:104:LEU:HD21	4:D:108:GLU:HB2	2.03	0.40
4:D:167:ARG:HB2	4:D:170:GLU:HG3	2.03	0.40
2:B:113:LEU:HD21	2:B:213:LEU:HD21	2.02	0.40
4:H:133:PHE:HB2	4:H:173:THR:HB	2.02	0.40
1:E:160:CYS:HB3	2:F:189:ARG:NH2	2.36	0.40
2:F:95:TRP:CE2	4:H:-19:LYS:HB2	2.57	0.40
3:C:14:GLN:OE1	3:C:116:VAL:HG23	2.21	0.40
4:H:95:LEU:HD23	4:H:180:SER:HA	2.03	0.40
3:G:14:GLN:HB2	4:H:8:VAL:HG22	2.02	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	193/206 (94%)	182 (94%)	9 (5%)	2 (1%)	22	75
1	E	180/206 (87%)	172 (96%)	6 (3%)	2 (1%)	21	74
2	B	236/241 (98%)	225 (95%)	11 (5%)	0	100	100
2	F	225/241 (93%)	213 (95%)	12 (5%)	0	100	100
3	C	177/182 (97%)	172 (97%)	5 (3%)	0	100	100
3	G	175/182 (96%)	168 (96%)	7 (4%)	0	100	100
4	D	195/217 (90%)	185 (95%)	9 (5%)	1 (0%)	38	86
4	H	195/217 (90%)	189 (97%)	6 (3%)	0	100	100
All	All	1576/1692 (93%)	1506 (96%)	65 (4%)	5 (0%)	50	92

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	GLY
1	E	60	GLY
1	E	79	ASP
1	A	79	ASP
4	D	105	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	157/185 (85%)	144 (92%)	13 (8%)	16	56
1	E	143/185 (77%)	121 (85%)	22 (15%)	4	20
2	B	185/206 (90%)	172 (93%)	13 (7%)	21	66
2	F	175/206 (85%)	167 (95%)	8 (5%)	37	81
3	C	147/163 (90%)	142 (97%)	5 (3%)	49	87
3	G	146/163 (90%)	141 (97%)	5 (3%)	49	87
4	D	168/189 (89%)	156 (93%)	12 (7%)	21	64
4	H	166/189 (88%)	151 (91%)	15 (9%)	14	50
All	All	1287/1486 (87%)	1194 (93%)	93 (7%)	21	64

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	112	LEU
1	A	123	GLN
1	A	124	LEU
1	A	129	SER
1	A	134	VAL
1	A	146	VAL
1	A	160	CYS
1	A	164	MET
1	A	183	PHE
1	A	191	ASN
1	A	193	ILE
1	A	194	ILE

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Mol	Chain	Res	Type
2	B	23	GLN
2	B	37	THR
2	B	78	LEU
2	B	80	THR
2	B	98	THR
2	B	106	ARG
2	B	134	THR
2	B	146	PHE
2	B	180	ASN
2	B	195	THR
2	B	204	PHE
2	B	215	GLU
2	B	238	ARG
3	C	19	ILE
3	C	36	LEU
3	C	60	LEU
3	C	88	GLN
3	C	95	SER
4	D	-21	LYS
4	D	25	ARG
4	D	38	VAL
4	D	39	ARG
4	D	92	LEU
4	D	95	LEU
4	D	115	THR
4	D	148	LEU
4	D	164	MET
4	D	173	THR
4	D	176	VAL
4	D	181	LEU
1	E	3	VAL
1	E	11	THR
1	E	12	VAL
1	E	17	THR
1	E	63	THR
1	E	76	HIS
1	E	81	GLN
1	E	85	SER
1	E	87	THR
1	E	109	LYS
1	E	110	VAL
1	E	112	LEU

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Mol	Chain	Res	Type
1	E	123	GLN
1	E	124	LEU
1	E	129	SER
1	E	134	VAL
1	E	146	VAL
1	E	160	CYS
1	E	164	MET
1	E	183	PHE
1	E	193	ILE
1	E	194	ILE
2	F	23	GLN
2	F	80	THR
2	F	134	THR
2	F	146	PHE
2	F	167	CYS
2	F	180	ASN
2	F	195	THR
2	F	204	PHE
3	G	14	GLN
3	G	36	LEU
3	G	42	VAL
3	G	116	VAL
3	G	120	THR
4	H	-21	LYS
4	H	-19	LYS
4	H	28	THR
4	H	70	ARG
4	H	71	THR
4	H	77	THR
4	H	90	THR
4	H	92	LEU
4	H	105	SER
4	H	107	THR
4	H	112	HIS
4	H	115	THR
4	H	130	VAL
4	H	164	MET
4	H	181	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	69	GLN
4	D	10	GLN
2	F	27	HIS
2	F	69	GLN
4	H	157	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

Mogul failed to run properly - this section will therefore be empty.

## 5.5 Carbohydrates ⓘ

Mogul failed to run properly - this section will therefore be empty.

## 5.6 Ligand geometry ⓘ

Mogul failed to run properly - this section will therefore be empty.

## 5.7 Other polymers ⓘ

Mogul failed to run properly - this section will therefore be empty.

## 5.8 Polymer linkage issues

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	G	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	G	85:GLU	C	86:ALA	N	3.12

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	197/206 (95%)	-0.07	0 100 100	22, 40, 83, 109	0
1	E	188/206 (91%)	0.57	3 (1%) 68 22	62, 99, 133, 142	0
2	B	238/241 (98%)	-0.06	0 100 100	20, 38, 70, 107	0
2	F	229/241 (95%)	0.47	3 (1%) 74 27	60, 92, 123, 153	0
3	C	179/182 (98%)	-0.03	0 100 100	23, 41, 72, 95	0
3	G	179/182 (98%)	0.40	4 (2%) 59 16	53, 79, 113, 132	0
4	D	199/217 (91%)	-0.10	0 100 100	27, 44, 71, 100	0
4	H	199/217 (91%)	0.07	0 100 100	54, 73, 96, 114	0
All	All	1608/1692 (95%)	0.15	10 (0%) 86 44	20, 65, 113, 153	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	86	ALA	3.0
1	E	146	VAL	2.4
2	F	194	ALA	2.4
1	E	113	HIS	2.3
1	E	111	TYR	2.3
2	F	70	GLU	2.2
2	F	215	GLU	2.1
3	G	164	LYS	2.1
3	G	173	PRO	2.1
3	G	177	HIS	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 ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

There are no ligands in this entry.

### 6.5 Other polymers ⓘ

There are no such residues in this entry.