



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 29, 2020 – 04:51 PM EST

PDB ID : 1U3H
Title : Crystal structure of mouse TCR 172.10 complexed with MHC class II I-Au molecule at 2.4 Å
Authors : Maynard, J.; Petersson, K.; Wilson, D.H.; Adams, E.J.; Blondelle, S.E.; Boulanger, M.J.; Wilson, D.B.; Garcia, K.C.
Deposited on : 2004-07-21
Resolution : 2.42 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.4
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.4

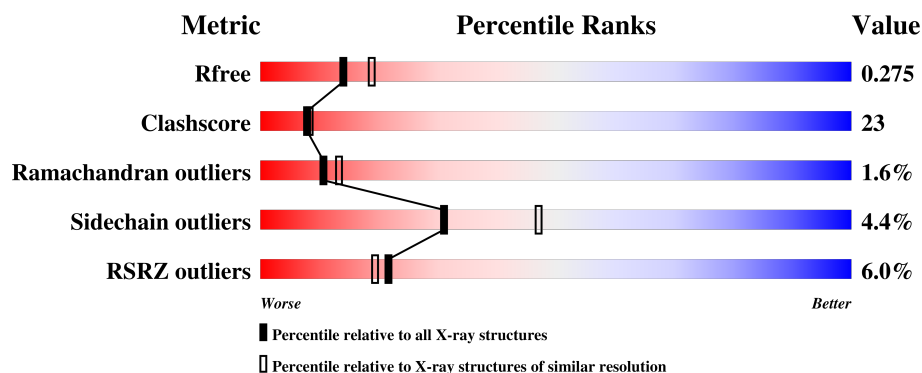
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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}	111664	4090 (2.44-2.40)
Clashscore	122126	4587 (2.44-2.40)
Ramachandran outliers	120053	4522 (2.44-2.40)
Sidechain outliers	120020	4523 (2.44-2.40)
RSRZ outliers	108989	3987 (2.44-2.40)

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 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	110	<div> <div>5%</div> <div>49%</div> <div>47%</div> <div>.</div> </div>
1	E	110	<div> <div>7%</div> <div>50%</div> <div>44%</div> <div>6%</div> </div>
2	B	111	<div> <div>10%</div> <div>45%</div> <div>49%</div> <div>6%</div> </div>
2	F	111	<div> <div>10%</div> <div>45%</div> <div>47%</div> <div>8%</div> </div>
3	C	182	<div> <div>%</div> <div>72%</div> <div>26%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
3	G	182	<div><div>%</div><div><div></div><div>78%</div><div>20%</div><div></div></div><div></div></div>
4	D	189	<div><div>8%</div><div><div></div><div>66%</div><div>30%</div><div></div></div><div></div></div>
4	H	189	<div><div>9%</div><div><div></div><div>63%</div><div>34%</div><div></div></div><div></div></div>
5	I	12	<div><div></div><div><div></div><div>67%</div><div>33%</div><div></div></div><div></div></div>
5	P	12	<div><div></div><div><div></div><div>58%</div><div>42%</div><div></div></div><div></div></div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9909 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 T-cell receptor alpha-chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	110	Total	C	N	O	S	0	0	0
			869	554	144	169	2			
1	E	110	Total	C	N	O	S	0	0	0
			869	554	144	169	2			

- Molecule 2 is a protein called Mouse TCRVbeta 172.10, extracellular variable domain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	111	Total	C	N	O	S	0	0	0
			853	528	148	174	3			
2	F	111	Total	C	N	O	S	0	0	0
			853	528	148	174	3			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	182	Total	C	N	O	S	0	0	0
			1464	946	233	282	3			
3	G	182	Total	C	N	O	S	0	0	0
			1464	946	233	282	3			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	ILE	-	cloning artifact	UNP P14438
C	2	GLU	-	cloning artifact	UNP P14438
C	3	ALA	-	cloning artifact	UNP P14438
G	1	ILE	-	cloning artifact	UNP P14438
G	2	GLU	-	cloning artifact	UNP P14438
G	3	ALA	-	cloning artifact	UNP P14438

- Molecule 4 is a protein called H-2 class II histocompatibility antigen, A-U beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	189	Total	C	N	O	S	0	0	0
			1586	998	283	299	6			
4	H	189	Total	C	N	O	S	0	0	0
			1586	998	283	299	6			

- Molecule 5 is a protein called Myelin basic protein (MBP)-peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	P	12	Total	C	N	O	0	0	0
			91	52	20	19			
5	I	12	Total	C	N	O	0	0	0
			91	52	20	19			

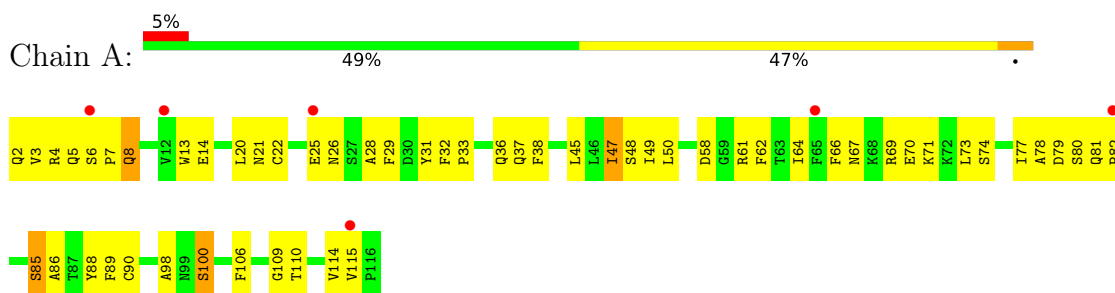
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	4	Total	O	0	0
			4	4		
6	B	4	Total	O	0	0
			4	4		
6	C	49	Total	O	0	0
			49	49		
6	D	42	Total	O	0	0
			42	42		
6	P	3	Total	O	0	0
			3	3		
6	E	4	Total	O	0	0
			4	4		
6	F	3	Total	O	0	0
			3	3		
6	G	42	Total	O	0	0
			42	42		
6	H	30	Total	O	0	0
			30	30		
6	I	2	Total	O	0	0
			2	2		

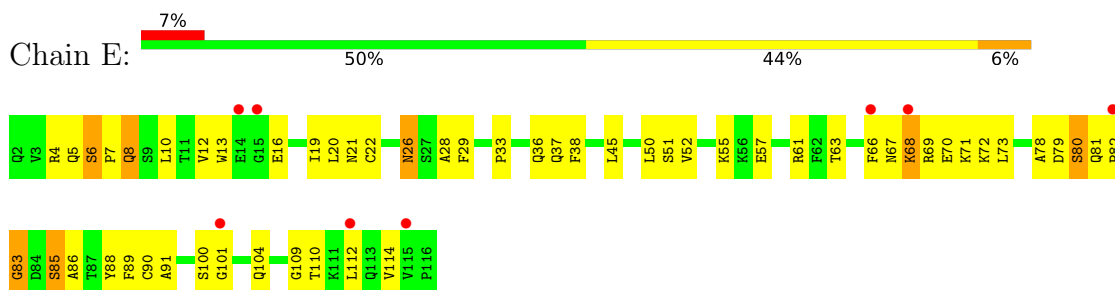
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

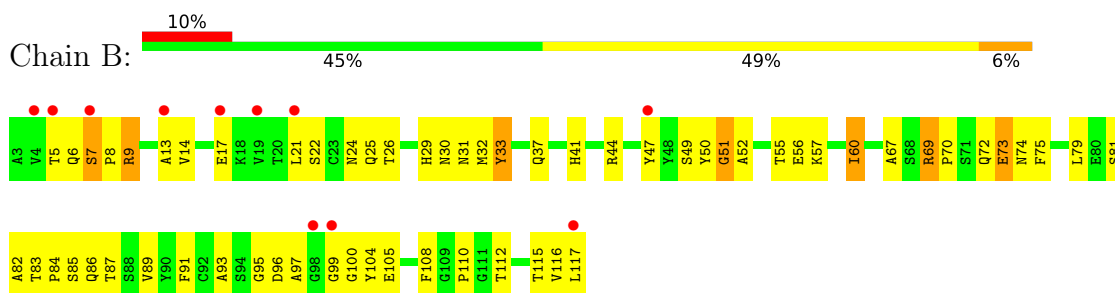
- Molecule 1: T-cell receptor alpha-chain



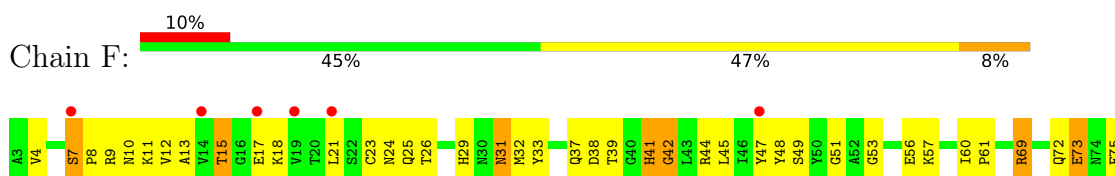
- Molecule 1: T-cell receptor alpha-chain

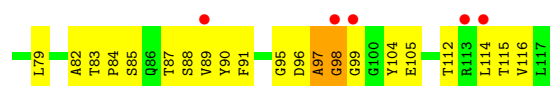


- Molecule 2: Mouse TCRVbeta 172.10, extracellular variable domain

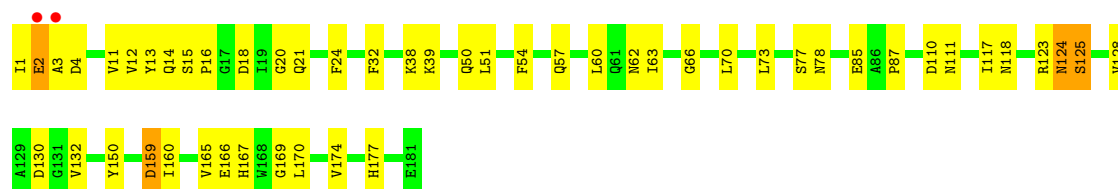


- Molecule 2: Mouse TCRVbeta 172.10, extracellular variable domain

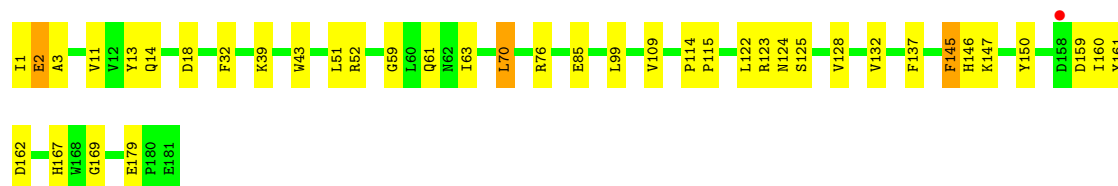
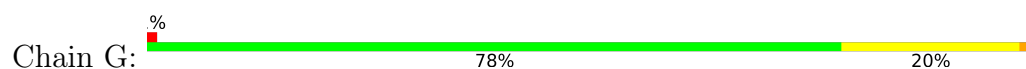




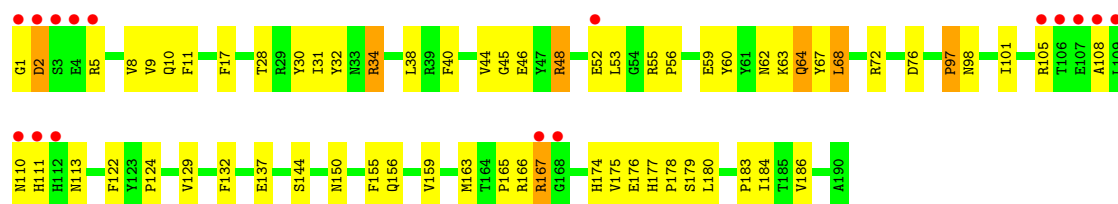
- Molecule 3: H-2 class II histocompatibility antigen, A-U alpha chain



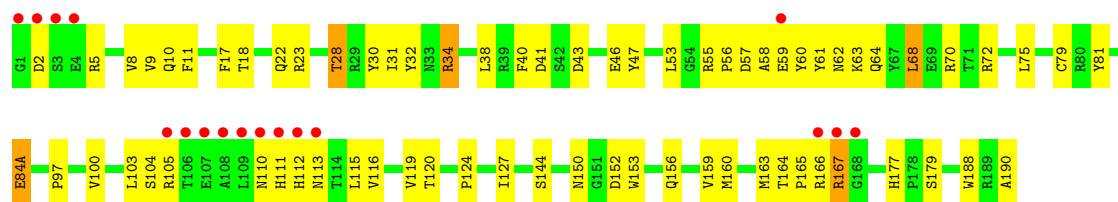
- Molecule 3: H-2 class II histocompatibility antigen, A-U alpha chain



- Molecule 4: H-2 class II histocompatibility antigen, A-U beta chain



- Molecule 4: H-2 class II histocompatibility antigen, A-U beta chain



- Molecule 5: Myelin basic protein (MBP)-peptide





- Molecule 5: Myelin basic protein (MBP)-peptide

Chain I:  67% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	87.83Å 327.16Å 127.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.51 – 2.42 41.51 – 2.42	Depositor EDS
% Data completeness (in resolution range)	89.1 (41.51-2.42) 89.2 (41.51-2.42)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.42Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.232 , 0.274 0.235 , 0.275	Depositor DCC
R_{free} test set	3489 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	42.1	Xtriage
Anisotropy	0.898	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9909	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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	0.35	0/892	0.65	2/1208 (0.2%)
1	E	0.37	0/892	0.68	3/1208 (0.2%)
2	B	0.39	0/873	0.76	4/1184 (0.3%)
2	F	0.37	0/873	0.67	1/1184 (0.1%)
3	C	0.40	0/1509	0.69	1/2059 (0.0%)
3	G	0.42	0/1509	0.68	1/2059 (0.0%)
4	D	0.41	0/1625	0.70	1/2206 (0.0%)
4	H	0.42	0/1625	0.69	0/2206
5	I	0.41	0/92	0.69	0/120
5	P	0.43	0/92	0.68	0/120
All	All	0.40	0/9982	0.69	13/13554 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	100	SER	C-N-CA	-7.21	107.17	122.30
2	F	53	GLY	N-CA-C	-6.52	96.79	113.10
2	B	52	ALA	CA-C-N	-6.26	103.68	116.20
1	E	100	SER	CA-C-N	6.21	128.61	116.20
2	B	52	ALA	C-N-CA	6.13	135.18	122.30
2	B	51	GLY	N-CA-C	5.89	127.83	113.10
1	E	100	SER	O-C-N	-5.69	113.53	123.20
1	A	100	SER	N-CA-C	-5.57	95.96	111.00
1	A	85	SER	N-CA-C	5.51	125.88	111.00
3	C	124	ASN	CB-CA-C	-5.29	99.82	110.40
2	B	52	ALA	O-C-N	5.25	132.13	123.20
3	G	70	LEU	N-CA-C	-5.17	97.03	111.00
4	D	108	ALA	C-N-CA	-5.13	108.87	121.70

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	869	0	828	49	0
1	E	869	0	828	57	0
2	B	853	0	795	54	0
2	F	853	0	795	74	0
3	C	1464	0	1392	56	0
3	G	1464	0	1392	39	0
4	D	1586	0	1521	65	0
4	H	1586	0	1521	66	0
5	I	91	0	86	5	0
5	P	91	0	86	9	0
6	A	4	0	0	0	0
6	B	4	0	0	1	0
6	C	49	0	0	6	0
6	D	42	0	0	2	0
6	E	4	0	0	3	0
6	F	3	0	0	2	0
6	G	42	0	0	1	0
6	H	30	0	0	1	0
6	I	2	0	0	0	0
6	P	3	0	0	0	0
All	All	9909	0	9244	437	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 23.

All (437) 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)
3:C:1:ILE:CG2	6:C:230:HOH:O	1.78	1.26
2:F:60:ILE:HG22	6:F:120:HOH:O	1.41	1.17
3:C:124:ASN:O	3:C:125:SER:HB2	1.30	1.09
1:E:37:GLN:O	1:E:86:ALA:HB1	1.54	1.08
3:C:124:ASN:O	3:C:125:SER:CB	2.01	1.05
3:G:99:LEU:H	3:G:99:LEU:HD12	1.26	0.98
1:A:37:GLN:O	1:A:86:ALA:HB1	1.63	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:2:ASP:H	4:H:5:ARG:HD3	1.33	0.93
2:B:13:ALA:HB1	2:B:17:GLU:OE1	1.69	0.91
2:F:13:ALA:HB1	2:F:17:GLU:OE1	1.68	0.91
1:E:6:SER:HB3	1:E:7:PRO:HD3	1.52	0.90
2:B:41:HIS:HB3	2:B:44:ARG:HD2	1.52	0.90
3:C:62:ASN:HD21	5:P:3:GLN:NE2	1.70	0.89
1:A:6:SER:HB3	1:A:21:ASN:HB2	1.54	0.88
2:F:47:TYR:CE1	2:F:61:PRO:HB2	2.08	0.88
4:H:167:ARG:HB3	4:H:167:ARG:HH11	1.40	0.87
2:F:7:SER:HB3	2:F:8:PRO:HD3	1.58	0.86
1:E:37:GLN:HE22	2:F:37:GLN:HE22	1.25	0.85
4:H:59:GLU:HG2	4:H:63:LYS:HE2	1.60	0.84
4:H:68:LEU:HD22	4:H:72:ARG:HE	1.44	0.82
4:H:150:ASN:HD21	4:H:156:GLN:HE21	1.23	0.82
3:C:62:ASN:HD21	5:P:3:GLN:HE22	1.23	0.82
2:B:84:PRO:HA	2:B:116:VAL:HG21	1.61	0.81
3:C:1:ILE:HG22	6:C:230:HOH:O	1.60	0.80
1:A:47:ILE:HD11	1:A:58:ASP:HB3	1.64	0.79
4:D:34:ARG:HG3	4:D:34:ARG:O	1.81	0.79
2:B:84:PRO:HA	2:B:116:VAL:CG2	2.12	0.79
4:D:150:ASN:HD21	4:D:156:GLN:HE21	1.28	0.79
1:A:6:SER:HB3	1:A:7:PRO:HD3	1.66	0.78
1:E:85:SER:HB2	6:E:118:HOH:O	1.83	0.78
1:A:25:GLU:HA	1:A:71:LYS:HD3	1.65	0.77
2:B:7:SER:OG	2:B:8:PRO:HD2	1.84	0.77
2:F:47:TYR:CE1	2:F:61:PRO:CB	2.69	0.75
1:E:6:SER:HB3	1:E:21:ASN:HB2	1.67	0.75
4:D:28:THR:HG21	6:D:194:HOH:O	1.86	0.74
2:F:31:ASN:HB2	2:F:49:SER:O	1.86	0.74
1:E:5:GLN:HE22	1:E:89:PHE:HA	1.52	0.74
4:H:2:ASP:H	4:H:5:ARG:CD	1.99	0.74
1:A:47:ILE:HG22	1:A:64:ILE:HD11	1.70	0.74
2:F:37:GLN:HE21	2:F:91:PHE:HE2	1.34	0.73
2:F:13:ALA:CB	2:F:17:GLU:OE1	2.37	0.73
2:B:37:GLN:HE21	2:B:91:PHE:HE2	1.35	0.73
4:D:59:GLU:HG2	4:D:63:LYS:HE2	1.69	0.73
2:F:41:HIS:CD2	2:F:41:HIS:H	2.08	0.71
3:G:14:GLN:HB3	4:H:8:VAL:HG23	1.71	0.71
4:H:28:THR:CG2	4:H:30:TYR:HE1	2.03	0.71
4:D:105:ARG:HB2	4:D:113:ASN:OD1	1.91	0.71
4:H:10:GLN:HB2	4:H:31:ILE:HB	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:99:LEU:N	3:G:99:LEU:HD12	2.01	0.70
4:H:28:THR:HG21	6:H:199:HOH:O	1.92	0.70
2:F:41:HIS:HD2	2:F:41:HIS:H	1.38	0.69
4:H:68:LEU:HD22	4:H:72:ARG:NE	2.06	0.69
2:F:87:THR:HG23	2:F:114:LEU:O	1.92	0.69
4:H:104:SER:O	4:H:105:ARG:HG2	1.92	0.69
2:F:11:LYS:NZ	2:F:13:ALA:HB2	2.08	0.68
3:G:99:LEU:CD1	3:G:99:LEU:H	2.03	0.68
4:D:124:PRO:O	4:D:177:HIS:HE1	1.77	0.68
2:F:41:HIS:N	2:F:41:HIS:CD2	2.58	0.68
4:H:8:VAL:O	4:H:32:TYR:O	2.10	0.68
3:C:160:ILE:HG23	3:C:177:HIS:CE1	2.29	0.67
4:H:18:THR:HB	4:H:23:ARG:HB3	1.76	0.67
4:D:167:ARG:HB3	4:D:167:ARG:NH1	2.09	0.67
2:B:50:TYR:OH	3:C:57:GLN:HG3	1.95	0.67
2:B:73:GLU:CD	2:B:73:GLU:H	1.97	0.67
3:C:1:ILE:HG22	3:C:2:GLU:N	2.10	0.67
1:A:37:GLN:O	1:A:86:ALA:CB	2.39	0.66
2:F:11:LYS:HZ2	2:F:13:ALA:HB2	1.61	0.66
3:C:160:ILE:HG23	3:C:177:HIS:HE1	1.58	0.65
2:F:51:GLY:O	2:F:69:ARG:HD2	1.97	0.65
4:D:10:GLN:HB2	4:D:31:ILE:HB	1.77	0.65
2:B:69:ARG:HH11	2:B:69:ARG:HG3	1.60	0.65
2:B:9:ARG:NH2	2:B:110:PRO:HB2	2.11	0.65
2:B:32:MET:HG3	2:B:69:ARG:NH2	2.11	0.65
2:B:97:ALA:O	2:B:104:TYR:CE2	2.50	0.65
1:E:13:TRP:HB2	1:E:16:GLU:OE2	1.97	0.64
2:F:7:SER:CB	2:F:8:PRO:HD3	2.26	0.64
2:B:21:LEU:H	2:B:21:LEU:HD12	1.63	0.64
2:F:88:SER:O	2:F:90:TYR:CD1	2.50	0.64
1:E:6:SER:HB3	1:E:7:PRO:CD	2.27	0.64
2:F:24:ASN:HA	2:F:73:GLU:O	1.97	0.63
3:C:117:ILE:HD11	3:C:165:VAL:CG1	2.28	0.63
4:D:11:PHE:CE1	4:D:28:THR:HG23	2.33	0.63
1:A:69:ARG:O	1:A:70:GLU:HB2	1.97	0.63
4:D:2:ASP:H	4:D:5:ARG:HD3	1.64	0.63
4:D:177:HIS:HD2	4:D:179:SER:OG	1.82	0.63
2:B:13:ALA:CB	2:B:17:GLU:OE1	2.45	0.62
2:B:84:PRO:CA	2:B:116:VAL:HG21	2.28	0.62
3:C:160:ILE:CG2	3:C:177:HIS:HE1	2.12	0.62
2:F:11:LYS:HD3	2:F:12:VAL:N	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:38:ASP:OD1	2:F:88:SER:CB	2.48	0.62
2:B:51:GLY:HA2	2:B:69:ARG:NH1	2.14	0.62
2:B:79:LEU:N	2:B:79:LEU:HD12	2.14	0.62
2:B:21:LEU:HD12	2:B:21:LEU:N	2.14	0.62
2:F:11:LYS:O	2:F:114:LEU:HD12	1.99	0.62
2:F:7:SER:HB3	2:F:8:PRO:CD	2.27	0.62
4:D:8:VAL:O	4:D:32:TYR:O	2.16	0.62
1:E:6:SER:CB	1:E:7:PRO:HD3	2.26	0.62
1:A:47:ILE:HG22	1:A:64:ILE:CD1	2.29	0.61
3:C:1:ILE:N	3:C:1:ILE:HD12	2.15	0.61
3:C:167:HIS:CD2	3:C:169:GLY:H	2.18	0.61
1:E:66:PHE:HD1	1:E:73:LEU:HD13	1.65	0.61
4:H:81:TYR:O	4:H:84(A):GLU:HB2	2.00	0.61
2:B:41:HIS:HB3	2:B:44:ARG:CD	2.28	0.61
4:D:97:PRO:HD3	4:D:122:PHE:CB	2.31	0.61
4:D:167:ARG:HB3	4:D:167:ARG:HH11	1.64	0.60
1:E:7:PRO:HD3	1:E:21:ASN:HD22	1.67	0.60
3:G:1:ILE:HG22	3:G:2:GLU:N	2.16	0.60
4:H:177:HIS:HD2	4:H:179:SER:OG	1.84	0.60
3:C:1:ILE:HG23	6:C:230:HOH:O	1.68	0.60
3:G:14:GLN:HB3	4:H:8:VAL:CG2	2.31	0.60
1:A:66:PHE:HD1	1:A:73:LEU:HD13	1.67	0.60
2:F:83:THR:HG23	2:F:84:PRO:HD2	1.84	0.60
2:B:7:SER:CB	2:B:8:PRO:CD	2.80	0.59
2:F:88:SER:O	2:F:90:TYR:CE1	2.56	0.59
2:B:51:GLY:CA	2:B:69:ARG:NH1	2.66	0.59
2:B:30:ASN:HA	2:B:72:GLN:HE22	1.68	0.58
1:A:25:GLU:HA	1:A:71:LYS:CD	2.33	0.58
2:F:15:THR:HG23	2:F:84:PRO:HD3	1.85	0.58
3:C:167:HIS:HD2	3:C:169:GLY:H	1.51	0.58
4:D:62:ASN:O	4:D:68:LEU:HB2	2.03	0.58
2:F:26:THR:O	2:F:26:THR:HG22	2.03	0.58
2:F:87:THR:HG23	2:F:115:THR:HA	1.85	0.58
4:H:116:VAL:HG22	4:H:160:MET:HG3	1.86	0.58
2:F:21:LEU:HD12	2:F:21:LEU:N	2.18	0.58
3:C:123:ARG:HB2	3:C:128:VAL:HG21	1.86	0.58
1:A:45:LEU:HB3	2:B:105:GLU:OE2	2.03	0.58
3:C:123:ARG:HD3	3:C:159:ASP:OD2	2.04	0.58
2:F:84:PRO:HA	2:F:116:VAL:CG2	2.34	0.58
4:D:55:ARG:HB3	4:D:56:PRO:HD3	1.86	0.58
1:E:8:GLN:CD	1:E:8:GLN:H	2.07	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:87:THR:OG1	2:F:116:VAL:HG22	2.03	0.57
1:E:26:ASN:HD22	1:E:26:ASN:C	2.07	0.57
5:P:3:GLN:HA	5:P:3:GLN:NE2	2.19	0.57
2:F:89:VAL:HA	2:F:112:THR:O	2.05	0.57
1:E:38:PHE:CD1	1:E:86:ALA:HB2	2.40	0.56
3:G:114:PRO:O	3:G:167:HIS:HE1	1.88	0.56
4:H:28:THR:CG2	4:H:30:TYR:CE1	2.87	0.56
2:F:60:ILE:CG2	6:F:120:HOH:O	2.18	0.56
3:C:3:ALA:HB1	4:D:17:PHE:O	2.05	0.56
2:B:7:SER:HG	2:B:22:SER:H	1.54	0.56
1:E:19:ILE:O	1:E:20:LEU:HD23	2.06	0.56
3:G:1:ILE:HD12	3:G:1:ILE:N	2.21	0.56
3:G:76:ARG:HD2	4:H:53:LEU:HD23	1.88	0.56
2:F:32:MET:HG3	2:F:69:ARG:NH2	2.20	0.56
1:E:20:LEU:HD22	1:E:110:THR:HG21	1.88	0.56
2:F:79:LEU:N	2:F:79:LEU:HD12	2.20	0.56
3:G:85:GLU:O	3:G:169:GLY:HA3	2.06	0.56
4:D:111:HIS:O	4:D:165:PRO:HD2	2.05	0.55
1:A:13:TRP:HD1	1:A:115:VAL:HB	1.70	0.55
4:D:52:GLU:HG2	4:D:55:ARG:HH21	1.71	0.55
2:F:73:GLU:CD	2:F:73:GLU:H	2.09	0.55
4:H:30:TYR:HB2	4:H:38:LEU:HB3	1.88	0.55
3:C:170:LEU:HD13	3:C:174:VAL:HG23	1.89	0.55
4:D:9:VAL:HA	4:D:32:TYR:O	2.06	0.55
1:A:7:PRO:HD3	1:A:21:ASN:HB2	1.89	0.55
1:E:69:ARG:C	1:E:71:LYS:H	2.10	0.55
1:A:89:PHE:CE1	1:A:109:GLY:HA3	2.41	0.55
2:B:83:THR:HG22	2:B:85:SER:H	1.71	0.54
4:H:167:ARG:NH1	4:H:167:ARG:HB3	2.17	0.54
4:H:2:ASP:N	4:H:5:ARG:HD3	2.13	0.54
2:B:69:ARG:NH1	2:B:69:ARG:HG3	2.23	0.54
3:G:159:ASP:O	3:G:160:ILE:HD12	2.07	0.54
4:H:60:TYR:CZ	4:H:64:GLN:HG3	2.43	0.54
2:F:56:GLU:OE1	3:G:39:LYS:HE2	2.08	0.53
2:B:93:ALA:HB2	2:B:108:PHE:CD1	2.43	0.53
2:B:73:GLU:CD	2:B:73:GLU:N	2.62	0.53
1:E:101:GLY:HA3	5:I:3:GLN:HE22	1.73	0.53
1:E:81:GLN:N	1:E:114:VAL:HG11	2.23	0.53
3:G:167:HIS:CD2	3:G:169:GLY:H	2.27	0.53
4:H:55:ARG:HB3	4:H:56:PRO:HD3	1.91	0.53
3:C:1:ILE:HG22	3:C:2:GLU:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:166:ARG:HG2	4:D:166:ARG:HH11	1.73	0.53
4:D:11:PHE:HE1	4:D:28:THR:HG23	1.72	0.53
3:G:3:ALA:HB1	4:H:17:PHE:O	2.09	0.53
4:H:111:HIS:O	4:H:164:THR:HA	2.09	0.53
3:C:57:GLN:NE2	3:C:60:LEU:HD12	2.23	0.53
3:C:70:LEU:HD13	4:D:9:VAL:HG13	1.90	0.53
3:G:123:ARG:HG3	3:G:161:TYR:CE2	2.43	0.53
4:H:150:ASN:ND2	4:H:156:GLN:HE21	2.02	0.52
5:P:7:SER:O	5:P:8:GLN:HB2	2.08	0.52
2:F:37:GLN:NE2	2:F:91:PHE:HE2	2.06	0.52
2:B:97:ALA:O	2:B:104:TYR:HE2	1.92	0.52
1:E:80:SER:O	1:E:81:GLN:HB2	2.10	0.52
2:F:11:LYS:O	2:F:114:LEU:HA	2.09	0.52
1:E:55:LYS:HE2	1:E:57:GLU:OE2	2.09	0.52
2:F:47:TYR:CZ	2:F:57:LYS:HG2	2.45	0.52
4:H:41:ASP:OD1	4:H:43:ASP:HB2	2.09	0.52
1:E:57:GLU:HB3	1:E:63:THR:HG23	1.90	0.52
2:F:32:MET:HE2	2:F:72:GLN:O	2.09	0.52
4:H:55:ARG:CB	4:H:56:PRO:HD3	2.39	0.52
4:D:11:PHE:HB2	5:P:4:TYR:CE1	2.45	0.52
2:F:41:HIS:O	2:F:42:GLY:O	2.27	0.52
3:G:115:PRO:HG3	3:G:145:PHE:CE1	2.44	0.52
4:D:11:PHE:HE1	4:D:28:THR:CG2	2.23	0.51
4:D:28:THR:HG22	4:D:30:TYR:CE1	2.45	0.51
2:F:29:HIS:HB3	2:F:95:GLY:O	2.10	0.51
1:A:31:TYR:CD2	2:B:100:GLY:HA2	2.46	0.51
3:C:51:LEU:N	3:C:51:LEU:HD12	2.25	0.51
4:H:53:LEU:O	4:H:53:LEU:HD23	2.11	0.51
1:A:8:GLN:CD	1:A:8:GLN:H	2.13	0.51
4:D:97:PRO:HD3	4:D:122:PHE:HB2	1.92	0.51
1:E:38:PHE:HD1	1:E:86:ALA:HB2	1.75	0.51
3:C:24:PHE:HZ	5:P:1:ALA:HB2	1.76	0.51
1:A:85:SER:OG	1:A:114:VAL:N	2.37	0.51
1:E:101:GLY:HA3	5:I:3:GLN:NE2	2.25	0.51
4:D:45:GLY:O	4:D:72:ARG:NH1	2.44	0.51
4:H:61:TYR:CE2	5:I:5:ARG:HB2	2.46	0.51
2:B:89:VAL:HA	2:B:112:THR:O	2.11	0.50
1:E:66:PHE:CE2	1:E:68:LYS:HA	2.46	0.50
2:F:97:ALA:HB1	3:G:61:GLN:NE2	2.26	0.50
3:G:137:PHE:CE1	3:G:147:LYS:HE3	2.46	0.50
2:B:56:GLU:CD	3:C:39:LYS:HZ2	2.14	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:31:ASN:OD1	2:B:97:ALA:HA	2.11	0.50
4:D:111:HIS:HB2	4:D:113:ASN:HD21	1.77	0.50
2:B:24:ASN:HA	2:B:73:GLU:O	2.12	0.50
4:D:97:PRO:HD3	4:D:122:PHE:HB3	1.92	0.50
2:B:37:GLN:NE2	2:B:91:PHE:CE2	2.78	0.50
1:E:10:LEU:O	1:E:112:LEU:HA	2.12	0.50
1:A:4:ARG:HG2	1:A:4:ARG:HH11	1.77	0.50
4:D:177:HIS:CG	4:D:178:PRO:HD2	2.46	0.49
1:A:14:GLU:HB2	1:A:82:PRO:HD3	1.94	0.49
4:D:174:HIS:HD2	6:D:225:HOH:O	1.95	0.49
1:E:81:GLN:O	1:E:83:GLY:N	2.41	0.49
2:B:60:ILE:HG22	6:B:121:HOH:O	2.11	0.49
4:D:167:ARG:HH11	4:D:167:ARG:CB	2.26	0.49
2:F:41:HIS:HB2	2:F:44:ARG:HD2	1.94	0.49
4:H:40:PHE:HB2	4:H:47:TYR:CE1	2.47	0.49
3:C:62:ASN:ND2	5:P:3:GLN:HE22	2.00	0.49
1:E:45:LEU:HD22	2:F:105:GLU:HB2	1.95	0.49
4:H:110:ASN:OD1	4:H:166:ARG:NH1	2.47	0.48
1:A:38:PHE:HD1	1:A:86:ALA:HB2	1.79	0.48
1:A:61:ARG:HD2	1:A:79:ASP:O	2.13	0.48
1:E:81:GLN:C	1:E:114:VAL:HG11	2.33	0.48
2:F:31:ASN:C	2:F:31:ASN:HD22	2.15	0.48
2:F:7:SER:CB	2:F:8:PRO:CD	2.87	0.48
3:G:1:ILE:CG2	3:G:2:GLU:N	2.77	0.48
4:H:167:ARG:CB	4:H:167:ARG:HH11	2.20	0.48
4:D:8:VAL:HG22	4:D:9:VAL:N	2.29	0.48
1:A:26:ASN:HB3	1:A:29:PHE:CD2	2.48	0.48
3:C:77:SER:O	3:C:78:ASN:HB2	2.13	0.48
2:B:49:SER:HB3	2:B:69:ARG:NH1	2.29	0.48
3:G:122:LEU:HB2	3:G:162:ASP:HB2	1.96	0.48
1:A:69:ARG:C	1:A:71:LYS:H	2.18	0.48
4:D:28:THR:HB	4:D:40:PHE:HB3	1.95	0.48
1:E:61:ARG:HB3	1:E:78:ALA:O	2.13	0.48
4:H:144:SER:CB	4:H:159:VAL:HG22	2.44	0.48
2:B:87:THR:HG23	2:B:115:THR:HA	1.95	0.47
1:A:26:ASN:HD21	1:A:28:ALA:HB3	1.79	0.47
1:A:31:TYR:CE1	1:A:48:SER:HB2	2.48	0.47
1:A:38:PHE:CD1	1:A:86:ALA:HB2	2.49	0.47
2:B:29:HIS:HB3	2:B:95:GLY:O	2.14	0.47
2:B:33:TYR:N	2:B:33:TYR:CD1	2.81	0.47
3:G:123:ARG:HH11	3:G:123:ARG:HG3	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1:ILE:CG2	3:C:2:GLU:N	2.75	0.47
4:D:97:PRO:CD	4:D:122:PHE:HB3	2.44	0.47
1:E:37:GLN:HA	6:E:119:HOH:O	2.14	0.47
4:D:60:TYR:CZ	4:D:64:GLN:HG3	2.50	0.47
1:E:26:ASN:HB3	1:E:29:PHE:CE2	2.49	0.47
4:H:9:VAL:HA	4:H:32:TYR:O	2.15	0.47
3:C:70:LEU:CD1	4:D:9:VAL:HG13	2.44	0.47
2:F:47:TYR:HE1	2:F:61:PRO:HB2	1.72	0.47
1:A:26:ASN:HB3	1:A:29:PHE:CE2	2.50	0.47
2:B:9:ARG:CZ	2:B:110:PRO:HB2	2.44	0.47
3:C:123:ARG:HG3	3:C:123:ARG:HH11	1.78	0.47
1:E:5:GLN:NE2	1:E:109:GLY:HA2	2.30	0.47
4:D:46:GLU:HG3	4:D:62:ASN:OD1	2.15	0.47
4:H:104:SER:C	4:H:105:ARG:HG2	2.36	0.47
2:F:83:THR:HG22	2:F:85:SER:H	1.79	0.46
3:G:70:LEU:HD13	4:H:9:VAL:HG13	1.96	0.46
1:E:37:GLN:NE2	2:F:37:GLN:HE22	2.03	0.46
4:D:111:HIS:HB2	4:D:113:ASN:ND2	2.30	0.46
2:F:84:PRO:HA	2:F:116:VAL:HG21	1.97	0.46
3:C:50:GLN:HB2	3:C:51:LEU:HD12	1.97	0.46
1:A:5:GLN:NE2	1:A:90:CYS:H	2.13	0.46
4:D:166:ARG:HG2	4:D:166:ARG:NH1	2.30	0.46
4:D:101:ILE:HD12	4:D:186:VAL:HG12	1.98	0.46
4:D:1:GLY:N	4:D:5:ARG:HD3	2.30	0.46
1:E:4:ARG:HH11	1:E:4:ARG:HG2	1.79	0.46
2:F:11:LYS:HB3	2:F:114:LEU:HD13	1.97	0.46
2:F:4:VAL:HG13	2:F:25:GLN:HB3	1.96	0.46
1:A:81:GLN:C	1:A:114:VAL:HG11	2.35	0.46
3:C:117:ILE:HD11	3:C:165:VAL:HG13	1.98	0.46
4:H:11:PHE:HB2	5:I:4:TYR:CE1	2.51	0.46
4:D:132:PHE:CE2	4:D:137:GLU:HB2	2.51	0.46
1:E:70:GLU:HG2	1:E:72:LYS:HE2	1.98	0.46
4:H:62:ASN:O	4:H:68:LEU:HB2	2.15	0.46
1:A:38:PHE:HA	1:A:86:ALA:HB2	1.97	0.45
1:E:26:ASN:HB3	1:E:29:PHE:CD2	2.51	0.45
4:H:144:SER:HB2	4:H:159:VAL:HG22	1.97	0.45
1:A:62:PHE:HD1	1:A:77:ILE:HG12	1.81	0.45
3:C:14:GLN:HB3	4:D:8:VAL:HB	1.99	0.45
1:E:5:GLN:NE2	1:E:90:CYS:H	2.15	0.45
2:F:33:TYR:N	2:F:33:TYR:CD1	2.84	0.45
2:F:96:ASP:HB3	2:F:104:TYR:CD2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:119:VAL:HG11	4:H:127:ILE:HD11	1.99	0.45
2:B:7:SER:HG	2:B:22:SER:N	2.14	0.45
3:C:15:SER:HB2	3:C:16:PRO:HA	1.98	0.45
1:E:61:ARG:HD2	1:E:79:ASP:O	2.17	0.45
2:F:23:CYS:SG	2:F:24:ASN:N	2.90	0.45
4:D:31:ILE:HD12	4:D:31:ILE:N	2.32	0.45
1:E:26:ASN:ND2	1:E:28:ALA:H	2.15	0.45
1:A:4:ARG:O	1:A:22:CYS:HA	2.16	0.45
2:F:31:ASN:CB	2:F:49:SER:O	2.62	0.45
4:D:144:SER:HB3	4:D:159:VAL:HG22	1.99	0.45
1:A:6:SER:CB	1:A:7:PRO:HD3	2.40	0.44
3:C:167:HIS:HD2	3:C:169:GLY:N	2.14	0.44
4:H:177:HIS:CD2	4:H:179:SER:H	2.35	0.44
3:C:85:GLU:HG3	4:D:34:ARG:NH2	2.32	0.44
4:D:67:TYR:O	4:D:68:LEU:C	2.54	0.44
4:H:40:PHE:HB2	4:H:47:TYR:CD1	2.52	0.44
3:C:32:PHE:C	3:C:32:PHE:CD1	2.90	0.44
1:E:50:LEU:C	1:E:52:VAL:H	2.20	0.44
1:E:66:PHE:CD1	1:E:73:LEU:HD13	2.50	0.44
3:G:43:TRP:CH2	3:G:52:ARG:HG3	2.52	0.44
1:E:81:GLN:CA	1:E:114:VAL:HG11	2.47	0.44
3:G:160:ILE:HG13	3:G:179:GLU:HB3	1.99	0.44
4:H:8:VAL:HG22	4:H:9:VAL:N	2.32	0.44
4:D:97:PRO:HG3	4:D:122:PHE:HB3	1.99	0.44
4:D:2:ASP:N	4:D:5:ARG:HD3	2.30	0.44
2:B:14:VAL:HA	2:B:116:VAL:O	2.18	0.44
3:C:110:ASP:OD1	3:C:111:ASN:N	2.47	0.43
3:C:132:VAL:HA	3:C:150:TYR:O	2.18	0.43
2:F:7:SER:O	2:F:8:PRO:C	2.54	0.43
4:H:31:ILE:HD12	4:H:31:ILE:N	2.33	0.43
4:H:5:ARG:HG2	4:H:5:ARG:HH11	1.83	0.43
1:A:7:PRO:O	1:A:110:THR:HG23	2.18	0.43
1:A:2:GLN:HG2	1:A:3:VAL:N	2.32	0.43
1:E:10:LEU:HA	6:E:117:HOH:O	2.18	0.43
3:G:132:VAL:HA	3:G:150:TYR:O	2.19	0.43
3:G:137:PHE:CZ	3:G:147:LYS:HE3	2.53	0.43
4:H:58:ALA:O	4:H:62:ASN:ND2	2.51	0.43
3:C:1:ILE:CG2	3:C:2:GLU:H	2.30	0.43
1:E:104:GLN:HG3	2:F:99:GLY:O	2.18	0.43
4:H:105:ARG:HG3	4:H:113:ASN:HA	1.99	0.43
1:A:85:SER:HG	1:A:114:VAL:H	1.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:81:GLN:H	1:E:114:VAL:HG11	1.84	0.43
3:C:118:ASN:HB2	3:C:166:GLU:HB2	2.00	0.43
2:F:47:TYR:OH	2:F:57:LYS:HE2	2.19	0.43
4:D:1:GLY:H2	4:D:5:ARG:HD3	1.83	0.43
1:E:26:ASN:ND2	1:E:26:ASN:C	2.72	0.43
3:G:85:GLU:CD	4:H:34:ARG:HH21	2.22	0.43
3:C:13:TYR:OH	3:C:18:ASP:HB3	2.19	0.43
3:G:11:VAL:HB	4:H:11:PHE:HB3	1.99	0.43
2:B:67:ALA:HB1	2:B:75:PHE:CE1	2.53	0.43
4:D:44:VAL:HG13	4:D:48:ARG:NH2	2.34	0.43
4:H:5:ARG:H	4:H:5:ARG:HG3	1.57	0.43
1:A:32:PHE:N	1:A:32:PHE:CD1	2.87	0.43
4:H:57:ASP:OD1	5:I:7:SER:HB2	2.19	0.43
1:A:38:PHE:HA	1:A:86:ALA:CB	2.48	0.42
2:B:70:PRO:HD2	2:B:74:ASN:O	2.18	0.42
4:H:188:TRP:CH2	4:H:190:ALA:HA	2.54	0.42
1:A:98:ALA:O	1:A:100:SER:O	2.37	0.42
3:C:21:GLN:HG2	6:C:204:HOH:O	2.18	0.42
4:D:163:MET:HE1	4:D:165:PRO:HB3	2.00	0.42
3:G:59:GLY:O	3:G:63:ILE:HG12	2.18	0.42
2:B:47:TYR:CZ	2:B:57:LYS:HG2	2.54	0.42
4:D:122:PHE:CE1	4:D:155:PHE:HB2	2.54	0.42
1:E:83:GLY:C	1:E:85:SER:H	2.22	0.42
2:F:87:THR:O	2:F:87:THR:HG22	2.19	0.42
1:A:20:LEU:O	1:A:74:SER:HB2	2.19	0.42
1:A:61:ARG:HB3	1:A:78:ALA:O	2.19	0.42
3:C:12:VAL:O	3:C:20:GLY:HA2	2.19	0.42
2:F:15:THR:O	2:F:15:THR:HG22	2.19	0.42
2:F:69:ARG:HH11	2:F:69:ARG:HG3	1.83	0.42
3:C:87:PRO:HG3	6:C:196:HOH:O	2.18	0.42
4:D:177:HIS:ND1	4:D:178:PRO:HD2	2.35	0.42
1:E:22:CYS:N	1:E:73:LEU:O	2.52	0.42
2:F:45:LEU:HD21	2:F:48:TYR:CD2	2.55	0.42
1:E:6:SER:CB	1:E:21:ASN:HB2	2.43	0.42
4:H:28:THR:HG21	4:H:30:TYR:HE1	1.83	0.42
3:C:73:LEU:HB3	4:D:32:TYR:CD1	2.54	0.42
1:A:31:TYR:HE1	1:A:48:SER:HB2	1.85	0.42
1:A:61:ARG:HD2	1:A:78:ALA:O	2.19	0.42
4:D:55:ARG:CB	4:D:56:PRO:HD3	2.50	0.42
1:A:33:PRO:HB2	1:A:45:LEU:CD1	2.50	0.42
4:D:46:GLU:OE1	4:D:48:ARG:NH1	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:32:MET:CE	2:F:69:ARG:NE	2.82	0.42
3:G:123:ARG:HB2	3:G:128:VAL:CG2	2.50	0.42
1:A:49:ILE:HG12	1:A:50:LEU:N	2.35	0.42
4:D:30:TYR:HB2	4:D:38:LEU:HB3	2.01	0.42
2:F:25:GLN:O	2:F:73:GLU:HB2	2.20	0.42
3:G:1:ILE:HD12	3:G:1:ILE:H1	1.84	0.42
4:H:152:ASP:O	4:H:153:TRP:HB2	2.19	0.42
2:B:26:THR:HG22	2:B:26:THR:O	2.20	0.41
1:E:19:ILE:HG22	1:E:20:LEU:N	2.35	0.41
3:G:76:ARG:HD2	4:H:53:LEU:CD2	2.49	0.41
1:A:36:GLN:HB2	1:A:88:TYR:HE1	1.85	0.41
2:F:17:GLU:HG2	2:F:18:LYS:N	2.35	0.41
3:G:124:ASN:O	3:G:125:SER:HB2	2.21	0.41
4:H:75:LEU:O	4:H:79:CYS:HB2	2.19	0.41
2:B:7:SER:HB2	2:B:8:PRO:HD3	2.01	0.41
3:C:123:ARG:CD	3:C:159:ASP:OD2	2.68	0.41
1:E:82:PRO:HA	1:E:114:VAL:HB	2.03	0.41
3:C:38:LYS:O	3:C:39:LYS:HB2	2.20	0.41
4:D:97:PRO:CG	4:D:122:PHE:HB3	2.50	0.41
2:F:96:ASP:O	2:F:98:GLY:N	2.53	0.41
1:E:12:VAL:HG13	1:E:16:GLU:CD	2.40	0.41
2:F:21:LEU:CD1	2:F:21:LEU:N	2.82	0.41
4:H:34:ARG:HG3	4:H:34:ARG:O	2.20	0.41
3:C:15:SER:HA	3:C:16:PRO:C	2.41	0.41
4:H:124:PRO:O	4:H:177:HIS:HE1	2.03	0.41
4:D:129:VAL:HG13	4:D:175:VAL:HG22	2.03	0.41
4:D:59:GLU:O	4:D:63:LYS:HG3	2.21	0.41
1:E:69:ARG:C	1:E:71:LYS:N	2.74	0.41
2:F:82:ALA:HB1	2:F:116:VAL:CG1	2.51	0.41
4:H:103:LEU:HD13	4:H:115:LEU:HD23	2.02	0.41
4:D:180:LEU:HD13	4:D:184:ILE:HG13	2.02	0.41
1:E:6:SER:CB	1:E:7:PRO:CD	2.91	0.41
2:B:7:SER:HB2	2:B:8:PRO:CD	2.51	0.41
3:C:1:ILE:N	3:C:1:ILE:CD1	2.82	0.41
1:E:33:PRO:HG2	1:E:91:ALA:HB3	2.03	0.41
4:H:163:MET:SD	4:H:165:PRO:HD3	2.61	0.41
1:A:106:PHE:CD1	1:A:106:PHE:N	2.89	0.41
1:A:36:GLN:HB2	1:A:88:TYR:CE1	2.55	0.41
3:G:109:VAL:O	3:G:146:HIS:HA	2.20	0.41
3:C:50:GLN:C	3:C:51:LEU:HD12	2.42	0.41
1:E:36:GLN:HB2	1:E:88:TYR:HE1	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:13:ALA:O	2:F:116:VAL:HA	2.21	0.41
4:H:111:HIS:O	4:H:164:THR:HG23	2.20	0.41
3:C:11:VAL:HG13	3:C:63:ILE:HD13	2.03	0.40
3:C:66:GLY:HA2	5:P:4:TYR:HE2	1.86	0.40
3:G:32:PHE:C	3:G:32:PHE:CD1	2.94	0.40
6:G:183:HOH:O	4:H:120:THR:HB	2.21	0.40
2:B:25:GLN:HG2	2:B:32:MET:SD	2.62	0.40
2:B:32:MET:HE3	2:B:69:ARG:NE	2.37	0.40
2:B:5:THR:HG22	2:B:6:GLN:N	2.37	0.40
2:F:32:MET:CE	2:F:75:PHE:HB2	2.52	0.40
3:G:70:LEU:CD1	4:H:9:VAL:HG13	2.50	0.40
3:C:62:ASN:ND2	5:P:3:GLN:NE2	2.53	0.40
2:B:116:VAL:HG23	2:B:117:LEU:N	2.37	0.40
3:C:54:PHE:HA	6:C:220:HOH:O	2.21	0.40
2:B:79:LEU:HD23	2:B:86:GLN:OE1	2.22	0.40
4:D:176:GLU:HG2	4:D:183:PRO:HG3	2.04	0.40
2:F:8:PRO:O	2:F:112:THR:HG23	2.21	0.40
3:G:13:TYR:OH	3:G:18:ASP:HB3	2.21	0.40
3:G:145:PHE:CD1	3:G:145:PHE:N	2.89	0.40
3:G:63:ILE:HA	3:G:63:ILE:HD13	1.95	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	108/110 (98%)	90 (83%)	17 (16%)	1 (1%)	19	26
1	E	108/110 (98%)	89 (82%)	14 (13%)	5 (5%)	2	1
2	B	109/111 (98%)	94 (86%)	11 (10%)	4 (4%)	4	2
2	F	109/111 (98%)	89 (82%)	15 (14%)	5 (5%)	2	1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	180/182 (99%)	169 (94%)	10 (6%)	1 (1%)	27	38
3	G	180/182 (99%)	174 (97%)	6 (3%)	0	100	100
4	D	187/189 (99%)	171 (91%)	14 (8%)	2 (1%)	16	21
4	H	187/189 (99%)	161 (86%)	25 (13%)	1 (0%)	31	43
5	I	10/12 (83%)	10 (100%)	0	0	100	100
5	P	10/12 (83%)	9 (90%)	1 (10%)	0	100	100
All	All	1188/1208 (98%)	1056 (89%)	113 (10%)	19 (2%)	11	13

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	80	SER
2	F	7	SER
2	F	42	GLY
2	F	97	ALA
1	A	80	SER
2	B	82	ALA
4	D	2	ASP
4	D	110	ASN
1	E	51	SER
1	E	68	LYS
2	F	15	THR
2	B	81	SER
2	B	99	GLY
1	E	6	SER
1	E	83	GLY
3	C	125	SER
4	H	22	GLN
2	B	7	SER
2	F	98	GLY

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	94/94 (100%)	91 (97%)	3 (3%)	42	61
1	E	94/94 (100%)	90 (96%)	4 (4%)	32	48
2	B	91/91 (100%)	84 (92%)	7 (8%)	14	21
2	F	91/91 (100%)	84 (92%)	7 (8%)	14	21
3	C	163/163 (100%)	159 (98%)	4 (2%)	50	69
3	G	163/163 (100%)	160 (98%)	3 (2%)	62	78
4	D	175/175 (100%)	166 (95%)	9 (5%)	26	41
4	H	175/175 (100%)	165 (94%)	10 (6%)	23	34
5	I	9/9 (100%)	9 (100%)	0	100	100
5	P	9/9 (100%)	9 (100%)	0	100	100
All	All	1064/1064 (100%)	1017 (96%)	47 (4%)	31	47

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	47	ILE
1	A	67	ASN
2	B	9	ARG
2	B	33	TYR
2	B	55	THR
2	B	60	ILE
2	B	69	ARG
2	B	73	GLU
2	B	96	ASP
3	C	2	GLU
3	C	4	ASP
3	C	130	ASP
3	C	159	ASP
4	D	34	ARG
4	D	48	ARG
4	D	53	LEU
4	D	64	GLN
4	D	68	LEU
4	D	76	ASP
4	D	97	PRO
4	D	98	ASN
4	D	167	ARG
1	E	8	GLN

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Mol	Chain	Res	Type
1	E	26	ASN
1	E	67	ASN
1	E	85	SER
2	F	9	ARG
2	F	10	ASN
2	F	31	ASN
2	F	39	THR
2	F	41	HIS
2	F	69	ARG
2	F	73	GLU
3	G	2	GLU
3	G	51	LEU
3	G	145	PHE
4	H	28	THR
4	H	34	ARG
4	H	46	GLU
4	H	68	LEU
4	H	70	ARG
4	H	84(A)	GLU
4	H	97	PRO
4	H	100	VAL
4	H	112	HIS
4	H	167	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	26	ASN
1	A	67	ASN
1	A	81	GLN
1	A	104	GLN
2	B	24	ASN
2	B	28	ASN
2	B	37	GLN
2	B	72	GLN
2	B	74	ASN
2	B	106	GLN
3	C	57	GLN
3	C	84	ASN
3	C	118	ASN
3	C	167	HIS

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Mol	Chain	Res	Type
4	D	64	GLN
4	D	111	HIS
4	D	146	GLN
4	D	150	ASN
4	D	177	HIS
5	P	3	GLN
1	E	5	GLN
1	E	21	ASN
1	E	26	ASN
1	E	36	GLN
1	E	67	ASN
1	E	76	HIS
1	E	81	GLN
1	E	104	GLN
2	F	24	ASN
2	F	28	ASN
2	F	31	ASN
2	F	37	GLN
2	F	41	HIS
2	F	72	GLN
2	F	74	ASN
3	G	14	GLN
3	G	101	GLN
3	G	118	ASN
3	G	167	HIS
4	H	64	GLN
4	H	111	HIS
4	H	150	ASN
4	H	174	HIS
4	H	177	HIS
5	I	3	GLN
5	I	8	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	110/110 (100%)	0.53	6 (5%) 25 23	47, 78, 102, 112	0
1	E	110/110 (100%)	0.66	8 (7%) 15 13	51, 80, 100, 110	0
2	B	111/111 (100%)	0.51	11 (9%) 7 6	48, 74, 92, 101	0
2	F	111/111 (100%)	0.61	11 (9%) 7 6	52, 78, 97, 104	0
3	C	182/182 (100%)	0.14	2 (1%) 80 78	27, 46, 79, 107	0
3	G	182/182 (100%)	-0.02	1 (0%) 90 89	28, 45, 74, 102	0
4	D	189/189 (100%)	0.57	16 (8%) 11 9	25, 49, 119, 154	0
4	H	189/189 (100%)	0.69	17 (8%) 9 8	24, 50, 120, 155	0
5	I	12/12 (100%)	0.81	0 100 100	45, 52, 83, 96	0
5	P	12/12 (100%)	0.44	0 100 100	38, 48, 71, 84	0
All	All	1208/1208 (100%)	0.44	72 (5%) 22 19	24, 60, 99, 155	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	H	111	HIS	14.6
4	H	110	ASN	12.9
4	H	106	THR	12.6
4	D	108	ALA	12.5
4	D	106	THR	12.4
4	H	109	LEU	12.3
4	H	107	GLU	11.1
4	D	110	ASN	10.8
4	D	111	HIS	10.0
4	D	109	LEU	9.1
4	H	108	ALA	9.0
4	D	3	SER	7.1
4	H	105	ARG	5.4

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Mol	Chain	Res	Type	RSRZ
4	D	107	GLU	5.2
4	H	3	SER	5.0
3	C	2	GLU	4.7
1	E	82	PRO	4.7
4	H	112	HIS	4.5
4	D	168	GLY	4.0
3	C	3	ALA	4.0
4	D	1	GLY	3.7
4	H	2	ASP	3.6
1	E	115	VAL	3.6
4	D	5	ARG	3.5
4	D	167	ARG	3.5
4	D	105	ARG	3.3
1	E	68	LYS	3.3
2	B	4	VAL	3.3
1	A	82	PRO	3.3
1	E	101	GLY	3.3
2	F	47	TYR	3.3
2	B	98	GLY	3.3
4	D	2	ASP	3.2
1	A	6	SER	3.2
4	H	167	ARG	3.1
2	F	98	GLY	3.1
3	G	158	ASP	3.1
4	H	4	GLU	3.0
4	D	4	GLU	2.9
4	H	113	ASN	2.9
2	B	21	LEU	2.8
1	E	112	LEU	2.8
4	H	1	GLY	2.8
2	F	17	GLU	2.7
1	A	115	VAL	2.7
2	B	7	SER	2.6
2	F	19	VAL	2.6
2	F	114	LEU	2.5
2	F	99	GLY	2.5
2	F	7	SER	2.5
4	H	166	ARG	2.4
2	B	117	LEU	2.4
2	B	47	TYR	2.4
2	B	17	GLU	2.4
2	B	99	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	12	VAL	2.4
1	A	25	GLU	2.4
1	E	14	GLU	2.4
2	F	89	VAL	2.4
1	A	65	PHE	2.4
4	D	52	GLU	2.4
1	E	15	GLY	2.4
4	H	168	GLY	2.3
2	B	19	VAL	2.3
2	B	13	ALA	2.3
2	F	21	LEU	2.2
4	D	112	HIS	2.1
2	F	113	ARG	2.1
2	B	5	THR	2.1
4	H	59	GLU	2.1
1	E	66	PHE	2.1
2	F	14	VAL	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 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.