



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 07:44 pm BST

PDB ID : 1P7Q
Title : Crystal Structure of HLA-A2 Bound to LIR-1, a Host and Viral MHC Receptor
Authors : Willcox, B.E.; Thomas, L.M.; Bjorkman, P.J.
Deposited on : 2003-05-05
Resolution : 3.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

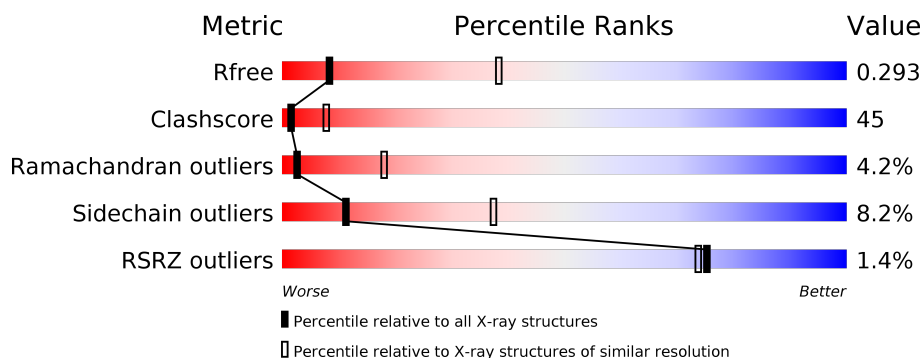
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

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 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	276	
2	B	99	
3	C	9	
4	D	197	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4498 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-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2237	1399	403	426	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	93	Total	C	N	O	S	0	0	0
			775	496	129	147	3			

- Molecule 3 is a protein called POL polyprotein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			69	46	12	11			

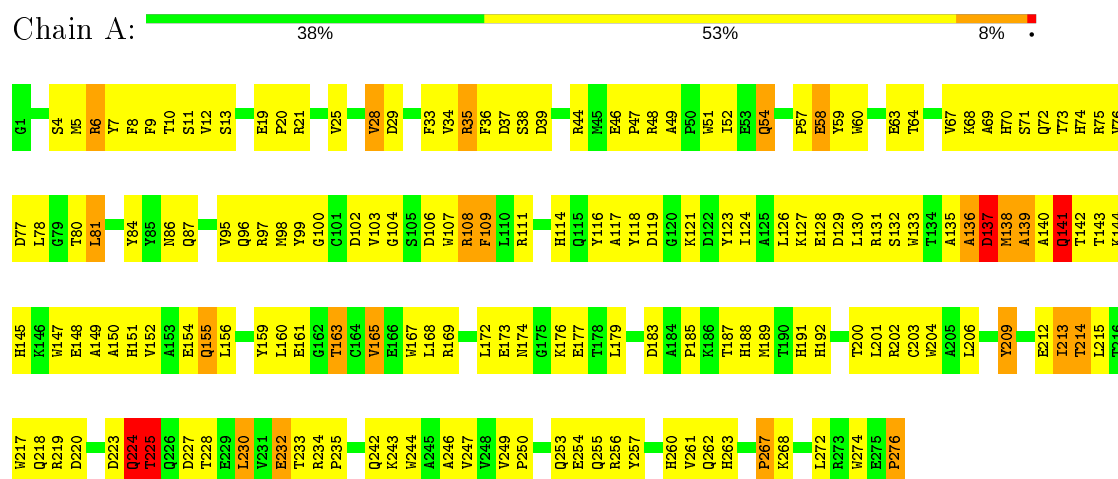
- Molecule 4 is a protein called leukocyte immunoglobulin-like receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	183	Total	C	N	O	S	0	0	0
			1417	907	244	260	6			

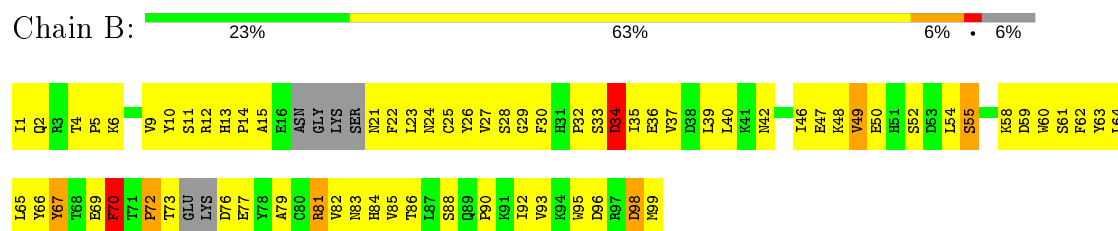
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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-2 alpha chain



- Molecule 2: Beta-2-microglobulin

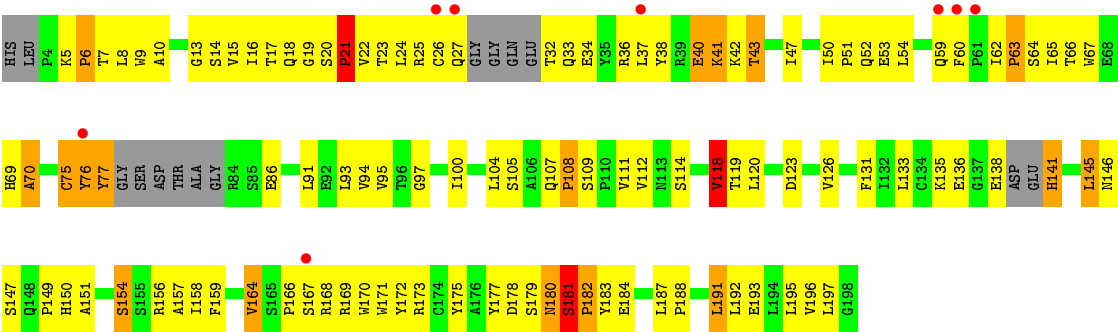


- Molecule 3: POL polyprotein



- Molecule 4: leukocyte immunoglobulin-like receptor 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	113.74 Å 113.74 Å 89.46 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.11 – 3.40 33.11 – 3.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (33.11-3.40) 99.7 (33.11-3.20)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 3.18 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.218 , 0.309 0.211 , 0.293	Depositor DCC
R_{free} test set	489 reflections (4.31%)	wwPDB-VP
Wilson B-factor (Å ²)	57.1	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4498	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.49	1/2303 (0.0%)	1.36	14/3131 (0.4%)
2	B	0.43	0/796	0.69	0/1080
3	C	0.56	0/70	0.82	0/94
4	D	0.44	1/1460 (0.1%)	0.78	3/1995 (0.2%)
All	All	0.46	2/4629 (0.0%)	1.10	17/6300 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	3
4	D	0	1
All	All	1	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	225	THR	C-N	-10.90	1.08	1.34
4	D	52	GLN	C-N	6.28	1.48	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	GLN	O-C-N	-35.73	65.53	122.70
1	A	225	THR	O-C-N	-30.64	73.67	122.70
1	A	225	THR	CA-C-N	22.10	165.82	117.20
1	A	224	GLN	CA-C-N	20.79	162.94	117.20
1	A	139	ALA	N-CA-CB	-16.48	87.02	110.10
1	A	141	GLN	CB-CG-CD	15.79	152.64	111.60
1	A	109	PHE	N-CA-CB	-12.36	88.36	110.60
1	A	108	ARG	CB-CA-C	-12.26	85.89	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	THR	N-CA-C	11.44	141.88	111.00
1	A	225	THR	C-N-CA	9.97	146.62	121.70
1	A	109	PHE	N-CA-C	9.21	135.88	111.00
4	D	118	VAL	O-C-N	-8.21	109.56	122.70
4	D	182	PRO	CA-N-CD	-7.08	101.59	111.50
1	A	138	MET	N-CA-C	7.05	130.05	111.00
1	A	108	ARG	N-CA-C	-6.60	93.17	111.00
1	A	28	VAL	N-CA-C	-6.56	93.29	111.00
4	D	181	SER	C-N-CD	-5.99	107.42	120.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	225	THR	CA

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	224	GLN	Mainchain,Peptide
1	A	225	THR	Mainchain
4	D	118	VAL	Mainchain

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	2237	0	2069	205	0
2	B	775	0	724	78	0
3	C	69	0	78	15	0
4	D	1417	0	1348	129	0
All	All	4498	0	4219	394	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 45.

All (394) 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:118:VAL:HG22	4:D:119:THR:H	0.90	1.04
4:D:118:VAL:CG2	4:D:119:THR:H	1.72	1.02
4:D:118:VAL:HG22	4:D:119:THR:N	1.74	1.01
4:D:6:PRO:HB2	4:D:27:GLN:H	1.26	1.00
4:D:37:LEU:HG	4:D:47:ILE:HD13	1.46	0.98
1:A:141:GLN:HA	1:A:144:LYS:HB3	1.55	0.88
1:A:130:LEU:HD23	1:A:160:LEU:HD12	1.58	0.84
1:A:9:PHE:HE2	1:A:99:TYR:CE2	1.95	0.83
4:D:70:ALA:HB2	4:D:95:VAL:HG23	1.61	0.83
4:D:16:ILE:HG13	4:D:22:VAL:HB	1.61	0.83
1:A:253:GLN:HA	1:A:255:GLN:NE2	1.95	0.80
1:A:219:ARG:HH21	1:A:256:ARG:HH11	1.28	0.79
1:A:84:TYR:CE2	1:A:142:THR:HB	2.18	0.79
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.63	0.79
1:A:121:LYS:HE2	2:B:1:ILE:HG13	1.64	0.79
4:D:50:ILE:HG23	4:D:54:LEU:HD12	1.66	0.78
1:A:152:VAL:HA	1:A:155:GLN:HE21	1.50	0.77
1:A:73:THR:HG21	3:C:6:VAL:HG13	1.64	0.77
4:D:6:PRO:HB2	4:D:27:GLN:N	2.00	0.77
1:A:218:GLN:HG2	1:A:223:ASP:OD2	1.84	0.76
4:D:181:SER:H	4:D:182:PRO:HD3	1.51	0.73
4:D:150:HIS:HE1	4:D:158:ILE:HG12	1.53	0.72
4:D:70:ALA:HB1	4:D:183:TYR:HB2	1.70	0.72
1:A:84:TYR:HE2	1:A:142:THR:HB	1.54	0.72
1:A:225:THR:O	1:A:227:ASP:N	2.21	0.72
4:D:107:GLN:HA	4:D:107:GLN:HE21	1.55	0.72
4:D:66:THR:H	4:D:69:HIS:HD2	1.35	0.72
1:A:13:SER:HA	1:A:20:PRO:HB3	1.70	0.72
1:A:219:ARG:HH21	1:A:256:ARG:NH1	1.87	0.71
4:D:70:ALA:HB2	4:D:95:VAL:CG2	2.21	0.70
4:D:37:LEU:HG	4:D:47:ILE:CD1	2.22	0.70
1:A:267:PRO:HG2	1:A:268:LYS:H	1.56	0.70
1:A:189:MET:HE1	1:A:274:TRP:HE3	1.55	0.70
4:D:181:SER:H	4:D:182:PRO:CD	2.05	0.70
4:D:169:ARG:HB2	4:D:171:TRP:NE1	2.07	0.69
1:A:173:GLU:O	1:A:176:LYS:HG3	1.93	0.69
1:A:163:THR:HG22	1:A:167:TRP:CD1	2.27	0.69
2:B:83:ASN:HB2	2:B:90:PRO:HG3	1.75	0.69
1:A:5:MET:CB	1:A:168:LEU:HD13	2.23	0.69
2:B:11:SER:HB2	2:B:21:ASN:HD21	1.56	0.69
4:D:32:THR:HG22	4:D:33:GLN:N	2.08	0.68
1:A:9:PHE:CE2	1:A:99:TYR:CE2	2.81	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2:GLN:HG2	2:B:32:PRO:HD3	1.75	0.67
1:A:74:HIS:NE2	1:A:97:ARG:HD2	2.11	0.66
2:B:40:LEU:HD11	2:B:79:ALA:HB3	1.75	0.66
4:D:41:LYS:HA	4:D:41:LYS:HE2	1.77	0.66
1:A:133:TRP:HB2	1:A:144:LYS:HD3	1.76	0.66
1:A:54:GLN:NE2	1:A:174:ASN:HB3	2.11	0.66
1:A:117:ALA:HB1	1:A:121:LYS:O	1.96	0.66
1:A:165:VAL:O	1:A:169:ARG:HG3	1.95	0.66
1:A:202:ARG:HE	1:A:246:ALA:HB2	1.61	0.66
1:A:8:PHE:HB2	1:A:25:VAL:CG1	2.26	0.65
1:A:19:GLU:HB2	1:A:75:ARG:HH12	1.59	0.65
2:B:50:GLU:HB3	2:B:67:TYR:CE1	2.31	0.65
4:D:7:THR:HB	4:D:9:TRP:CZ3	2.31	0.65
4:D:107:GLN:HA	4:D:107:GLN:NE2	2.11	0.65
1:A:133:TRP:H	1:A:144:LYS:NZ	1.95	0.65
4:D:32:THR:HG22	4:D:33:GLN:H	1.62	0.64
4:D:112:VAL:CG2	4:D:164:VAL:HG21	2.28	0.64
4:D:34:GLU:O	4:D:77:TYR:HB3	1.98	0.64
2:B:30:PHE:HE2	2:B:35:ILE:HB	1.62	0.63
1:A:253:GLN:HA	1:A:255:GLN:HE22	1.61	0.63
4:D:66:THR:H	4:D:69:HIS:CD2	2.15	0.63
1:A:244:TRP:HE1	2:B:99:MET:HG3	1.63	0.63
1:A:54:GLN:HE22	1:A:174:ASN:HB3	1.64	0.62
1:A:149:ALA:O	1:A:151:HIS:N	2.33	0.62
1:A:187:THR:OG1	1:A:272:LEU:HD21	1.98	0.62
1:A:214:THR:OG1	1:A:262:GLN:HB2	1.99	0.62
2:B:4:THR:HG23	2:B:5:PRO:HD2	1.82	0.62
2:B:5:PRO:HB3	2:B:30:PHE:CB	2.30	0.62
4:D:173:ARG:NH1	4:D:191:LEU:HB2	2.15	0.62
4:D:191:LEU:HD23	4:D:193:GLU:HG3	1.80	0.62
4:D:135:LYS:HG2	4:D:172:TYR:CE2	2.35	0.62
2:B:93:VAL:HG22	4:D:67:TRP:CZ2	2.35	0.62
2:B:29:GLY:HA2	2:B:61:SER:OG	2.01	0.61
1:A:219:ARG:NH2	1:A:256:ARG:NH1	2.49	0.61
1:A:233:THR:N	1:A:243:LYS:NZ	2.49	0.61
2:B:37:VAL:HG22	2:B:82:VAL:HG22	1.83	0.61
1:A:133:TRP:HB2	1:A:144:LYS:CD	2.31	0.61
4:D:24:LEU:HD22	4:D:91:LEU:HD21	1.82	0.61
1:A:235:PRO:HG3	2:B:26:TYR:CE1	2.35	0.60
1:A:103:VAL:CG1	1:A:168:LEU:HD23	2.32	0.60
2:B:5:PRO:CB	2:B:30:PHE:HB3	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:131:PHE:CD1	4:D:157:ALA:HB2	2.37	0.60
3:C:9:VAL:HG23	3:C:9:VAL:O	2.00	0.60
1:A:70:HIS:HA	3:C:6:VAL:HG11	1.84	0.59
3:C:3:LYS:HG2	3:C:4:GLU:N	2.17	0.59
4:D:181:SER:N	4:D:182:PRO:CD	2.63	0.59
1:A:137:ASP:O	1:A:141:GLN:NE2	2.36	0.59
1:A:8:PHE:HB2	1:A:25:VAL:HG12	1.83	0.59
1:A:111:ARG:NH1	1:A:128:GLU:OE1	2.35	0.59
1:A:212:GLU:O	1:A:213:ILE:HB	2.01	0.59
1:A:47:PRO:HB3	1:A:60:TRP:CZ2	2.38	0.59
1:A:19:GLU:CB	1:A:75:ARG:HH12	2.16	0.59
2:B:30:PHE:CE2	2:B:35:ILE:HD12	2.37	0.59
4:D:70:ALA:HB1	4:D:183:TYR:CB	2.32	0.59
4:D:7:THR:HB	4:D:9:TRP:HZ3	1.67	0.59
1:A:35:ARG:HG2	1:A:48:ARG:HD3	1.85	0.58
2:B:96:ASP:HB3	2:B:99:MET:H	1.67	0.58
4:D:37:LEU:HD12	4:D:38:TYR:H	1.68	0.58
2:B:55:SER:HB3	2:B:63:TYR:CZ	2.38	0.58
4:D:104:LEU:HG	4:D:192:LEU:HD22	1.86	0.58
4:D:136:GLU:HG3	4:D:191:LEU:HD12	1.84	0.58
1:A:156:LEU:O	1:A:160:LEU:HG	2.03	0.58
1:A:35:ARG:HD3	1:A:48:ARG:CZ	2.34	0.57
4:D:112:VAL:HG21	4:D:164:VAL:HG21	1.86	0.57
4:D:37:LEU:HD12	4:D:38:TYR:N	2.19	0.57
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.87	0.57
1:A:10:THR:HG21	2:B:62:PHE:CE1	2.40	0.57
1:A:138:MET:HA	1:A:141:GLN:NE2	2.20	0.56
1:A:37:ASP:O	1:A:39:ASP:N	2.38	0.56
1:A:218:GLN:OE1	1:A:260:HIS:HB2	2.05	0.56
1:A:12:VAL:O	1:A:12:VAL:HG13	2.05	0.56
1:A:143:THR:HG23	3:C:9:VAL:C	2.26	0.56
4:D:131:PHE:HE1	4:D:149:PRO:HG3	1.70	0.56
4:D:7:THR:O	4:D:26:CYS:HA	2.05	0.56
1:A:230:LEU:CD1	1:A:243:LYS:HE3	2.35	0.56
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.39	0.56
1:A:255:GLN:CD	1:A:255:GLN:H	2.09	0.56
4:D:8:LEU:HA	4:D:25:ARG:O	2.06	0.56
1:A:121:LYS:HE2	2:B:1:ILE:CG1	2.35	0.56
1:A:228:THR:HG23	1:A:247:VAL:HG13	1.87	0.56
1:A:124:ILE:O	1:A:124:ILE:HG23	2.06	0.56
1:A:52:ILE:HD11	1:A:59:TYR:CZ	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:52:SER:OG	2:B:64:LEU:HD23	2.05	0.56
2:B:54:LEU:HD11	2:B:62:PHE:HB3	1.88	0.55
2:B:67:TYR:N	2:B:67:TYR:CD1	2.74	0.55
1:A:154:GLU:N	1:A:154:GLU:OE2	2.40	0.55
1:A:73:THR:O	1:A:76:VAL:HG22	2.07	0.55
4:D:50:ILE:CG2	4:D:54:LEU:HD12	2.34	0.55
1:A:202:ARG:NH1	2:B:98:ASP:O	2.40	0.55
1:A:274:TRP:O	1:A:276:PRO:HD3	2.07	0.55
4:D:133:LEU:HD22	4:D:159:PHE:CG	2.42	0.55
4:D:182:PRO:C	4:D:184:GLU:H	2.10	0.55
1:A:253:GLN:CA	1:A:255:GLN:NE2	2.68	0.55
1:A:219:ARG:HB2	1:A:257:TYR:CE2	2.42	0.55
4:D:42:LYS:HD3	4:D:43:THR:O	2.06	0.55
2:B:63:TYR:O	2:B:63:TYR:CD1	2.60	0.54
4:D:138:GLU:OE1	4:D:141:HIS:N	2.40	0.54
4:D:16:ILE:CG2	4:D:65:ILE:HG21	2.37	0.54
4:D:10:ALA:HA	4:D:23:THR:O	2.08	0.54
1:A:191:HIS:HB2	1:A:274:TRP:CH2	2.43	0.54
2:B:50:GLU:O	2:B:66:TYR:HA	2.08	0.54
1:A:244:TRP:NE1	2:B:99:MET:HG3	2.23	0.54
1:A:68:LYS:O	1:A:71:SER:HB3	2.07	0.54
1:A:136:ALA:O	1:A:137:ASP:CB	2.56	0.53
2:B:96:ASP:HB3	2:B:99:MET:N	2.23	0.53
2:B:47:GLU:HB2	2:B:48:LYS:HD2	1.89	0.53
1:A:103:VAL:HG11	1:A:168:LEU:HD23	1.90	0.53
1:A:230:LEU:HD12	1:A:243:LYS:HE3	1.90	0.53
1:A:69:ALA:O	1:A:73:THR:HG23	2.09	0.53
1:A:137:ASP:OD1	1:A:140:ALA:HB2	2.07	0.53
2:B:84:HIS:CG	2:B:85:VAL:H	2.26	0.53
1:A:189:MET:HE1	1:A:274:TRP:CE3	2.40	0.53
4:D:22:VAL:HG11	4:D:65:ILE:HD12	1.90	0.53
2:B:69:GLU:O	2:B:70:PHE:HB3	2.09	0.53
1:A:109:PHE:CZ	1:A:161:GLU:HG2	2.44	0.53
1:A:64:THR:HA	1:A:67:VAL:HG12	1.91	0.53
4:D:175:TYR:CZ	4:D:188:PRO:HB3	2.44	0.53
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.44	0.52
1:A:6:ARG:HH11	1:A:6:ARG:HG3	1.73	0.52
2:B:27:VAL:HG23	2:B:27:VAL:O	2.08	0.52
1:A:149:ALA:C	1:A:151:HIS:H	2.13	0.52
4:D:16:ILE:HG22	4:D:65:ILE:HG21	1.92	0.52
1:A:49:ALA:HB1	1:A:51:TRP:NE1	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:TRP:O	1:A:223:ASP:HA	2.09	0.52
4:D:171:TRP:CD2	4:D:193:GLU:HG2	2.45	0.52
1:A:219:ARG:HB3	1:A:224:GLN:NE2	2.25	0.52
2:B:23:LEU:HD12	2:B:24:ASN:N	2.25	0.52
2:B:13:HIS:HB3	2:B:14:PRO:HD2	1.91	0.52
1:A:219:ARG:HE	1:A:256:ARG:HD3	1.75	0.52
1:A:206:LEU:HD23	1:A:242:GLN:CD	2.29	0.52
4:D:150:HIS:CE1	4:D:158:ILE:HG12	2.40	0.52
4:D:9:TRP:CE2	4:D:25:ARG:HB2	2.45	0.52
1:A:204:TRP:CH2	2:B:99:MET:HG2	2.45	0.51
1:A:114:HIS:CD2	1:A:156:LEU:HD11	2.44	0.51
4:D:41:LYS:CA	4:D:41:LYS:HE2	2.40	0.51
1:A:67:VAL:HG13	1:A:68:LYS:N	2.26	0.51
2:B:40:LEU:O	2:B:40:LEU:HD12	2.11	0.51
1:A:13:SER:HB3	1:A:78:LEU:HD13	1.93	0.51
1:A:159:TYR:CE2	3:C:3:LYS:HB2	2.46	0.51
1:A:152:VAL:C	1:A:154:GLU:H	2.14	0.50
4:D:107:GLN:HE21	4:D:108:PRO:HA	1.75	0.50
4:D:114:SER:OG	4:D:166:PRO:HD3	2.11	0.50
1:A:254:GLU:N	1:A:254:GLU:OE1	2.40	0.50
1:A:159:TYR:CD2	3:C:3:LYS:HB2	2.46	0.50
4:D:112:VAL:HG23	4:D:164:VAL:HG21	1.92	0.50
2:B:15:ALA:HA	2:B:72:PRO:HG2	1.93	0.50
1:A:28:VAL:HG11	1:A:179:LEU:HD13	1.92	0.50
1:A:253:GLN:CA	1:A:255:GLN:HE22	2.23	0.50
1:A:76:VAL:HG23	1:A:77:ASP:N	2.27	0.50
4:D:51:PRO:C	4:D:53:GLU:N	2.65	0.50
1:A:188:HIS:O	1:A:189:MET:HG3	2.11	0.50
4:D:191:LEU:CD2	4:D:193:GLU:HG3	2.40	0.50
1:A:119:ASP:O	1:A:121:LYS:HG3	2.11	0.50
2:B:86:THR:C	4:D:100:ILE:HG13	2.31	0.50
1:A:7:TYR:O	1:A:98:MET:HG2	2.12	0.50
4:D:51:PRO:C	4:D:53:GLU:H	2.15	0.50
1:A:54:GLN:HE22	1:A:174:ASN:CB	2.25	0.49
1:A:77:ASP:O	1:A:81:LEU:HB2	2.11	0.49
1:A:104:GLY:C	1:A:106:ASP:H	2.16	0.49
1:A:10:THR:HG23	1:A:96:GLN:HG3	1.94	0.49
1:A:6:ARG:NH1	1:A:6:ARG:HG3	2.27	0.49
4:D:170:TRP:C	4:D:171:TRP:HD1	2.16	0.49
4:D:171:TRP:N	4:D:171:TRP:CD1	2.78	0.49
1:A:108:ARG:HH11	1:A:108:ARG:HG2	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:ILE:HG13	1:A:262:GLN:O	2.13	0.49
4:D:37:LEU:HD23	4:D:60:PHE:CD2	2.47	0.49
1:A:130:LEU:HD23	1:A:160:LEU:CD1	2.37	0.49
2:B:42:ASN:HD21	2:B:76:ASP:HB2	1.78	0.49
4:D:69:HIS:O	4:D:70:ALA:C	2.50	0.49
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.47	0.49
4:D:123:ASP:HA	4:D:156:ARG:HG3	1.94	0.49
4:D:15:VAL:HA	4:D:94:VAL:O	2.13	0.49
4:D:20:SER:O	4:D:65:ILE:HG22	2.12	0.49
2:B:23:LEU:O	2:B:67:TYR:HA	2.12	0.48
1:A:167:TRP:CD1	3:C:1:ILE:HG13	2.48	0.48
1:A:152:VAL:C	1:A:154:GLU:N	2.67	0.48
1:A:127:LYS:HD3	1:A:132:SER:HB2	1.95	0.48
1:A:209:TYR:CD1	1:A:209:TYR:C	2.87	0.48
1:A:217:TRP:CD1	1:A:247:VAL:HG13	2.48	0.48
4:D:60:PHE:N	4:D:60:PHE:CD1	2.80	0.48
1:A:213:ILE:HD12	1:A:263:HIS:HD2	1.79	0.48
4:D:9:TRP:NE1	4:D:25:ARG:HB2	2.29	0.48
2:B:67:TYR:N	2:B:67:TYR:HD1	2.10	0.48
1:A:106:ASP:OD1	1:A:106:ASP:N	2.47	0.47
1:A:152:VAL:O	1:A:156:LEU:HB2	2.14	0.47
4:D:107:GLN:HA	4:D:108:PRO:HA	1.69	0.47
4:D:172:TYR:O	4:D:191:LEU:HA	2.14	0.47
1:A:121:LYS:HG2	2:B:1:ILE:HD12	1.97	0.47
2:B:29:GLY:C	2:B:61:SER:HB2	2.34	0.47
1:A:137:ASP:O	1:A:140:ALA:HB3	2.15	0.47
1:A:202:ARG:HE	1:A:246:ALA:CB	2.26	0.47
4:D:120:LEU:HD22	4:D:120:LEU:N	2.29	0.47
4:D:5:LYS:HB3	4:D:86:GLU:O	2.15	0.47
1:A:147:TRP:C	1:A:149:ALA:N	2.67	0.47
4:D:8:LEU:HD22	4:D:75:CYS:HB2	1.96	0.47
1:A:103:VAL:CG2	1:A:107:TRP:HA	2.45	0.47
1:A:219:ARG:HB3	1:A:224:GLN:HE22	1.80	0.47
1:A:234:ARG:HD2	2:B:10:TYR:CE1	2.49	0.47
1:A:6:ARG:HB3	1:A:100:GLY:HA3	1.96	0.47
1:A:117:ALA:HA	1:A:123:TYR:HB2	1.97	0.46
1:A:44:ARG:HD3	1:A:64:THR:HG21	1.96	0.46
4:D:32:THR:CG2	4:D:33:GLN:H	2.24	0.46
1:A:84:TYR:HD2	1:A:139:ALA:O	1.97	0.46
1:A:72:GLN:O	1:A:75:ARG:HB3	2.15	0.46
4:D:97:GLY:HA2	4:D:187:LEU:HG	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:ARG:CB	1:A:100:GLY:HA3	2.44	0.46
1:A:219:ARG:O	1:A:220:ASP:HB2	2.16	0.46
2:B:9:VAL:CG1	2:B:95:TRP:HB2	2.45	0.46
1:A:109:PHE:HD1	1:A:165:VAL:HG21	1.81	0.46
1:A:189:MET:HE1	1:A:274:TRP:HB2	1.96	0.46
1:A:33:PHE:O	1:A:52:ILE:HG21	2.15	0.46
1:A:107:TRP:CD1	1:A:107:TRP:N	2.82	0.46
1:A:6:ARG:N	1:A:6:ARG:HD3	2.31	0.46
1:A:80:THR:HG22	1:A:84:TYR:CE1	2.50	0.46
2:B:42:ASN:OD1	2:B:77:GLU:N	2.44	0.46
2:B:86:THR:O	4:D:100:ILE:HG13	2.16	0.46
2:B:85:VAL:O	4:D:100:ILE:CD1	2.63	0.46
4:D:177:TYR:C	4:D:177:TYR:CD1	2.88	0.46
1:A:5:MET:HB3	1:A:168:LEU:HD13	1.98	0.46
1:A:103:VAL:HG13	1:A:168:LEU:HD23	1.97	0.46
1:A:168:LEU:O	1:A:172:LEU:HG	2.15	0.46
4:D:40:GLU:O	4:D:42:LYS:N	2.46	0.46
1:A:202:ARG:HG3	1:A:246:ALA:HB2	1.98	0.46
1:A:217:TRP:NE1	1:A:247:VAL:HG22	2.30	0.46
2:B:4:THR:HG22	2:B:5:PRO:N	2.31	0.46
1:A:10:THR:HG21	2:B:62:PHE:HE1	1.81	0.45
3:C:1:ILE:C	3:C:1:ILE:CD1	2.84	0.45
4:D:109:SER:C	4:D:111:VAL:N	2.68	0.45
4:D:111:VAL:HG22	4:D:195:LEU:HB3	1.99	0.45
1:A:78:LEU:HD23	1:A:95:VAL:HG13	1.99	0.45
1:A:144:LYS:O	1:A:148:GLU:HG2	2.16	0.45
1:A:235:PRO:CG	2:B:26:TYR:CE1	2.99	0.45
1:A:149:ALA:C	1:A:151:HIS:N	2.69	0.45
4:D:109:SER:C	4:D:111:VAL:H	2.19	0.45
2:B:5:PRO:CA	2:B:30:PHE:HB3	2.47	0.45
4:D:54:LEU:HB3	4:D:59:GLN:O	2.17	0.45
4:D:22:VAL:CG1	4:D:65:ILE:HD12	2.47	0.45
4:D:24:LEU:HD22	4:D:91:LEU:CD2	2.45	0.45
1:A:116:TYR:HD1	1:A:117:ALA:N	2.15	0.45
1:A:155:GLN:H	1:A:155:GLN:HG3	1.55	0.45
4:D:97:GLY:HA2	4:D:187:LEU:HD23	1.98	0.45
4:D:178:ASP:O	4:D:180:ASN:N	2.49	0.45
4:D:22:VAL:CG2	4:D:23:THR:N	2.80	0.45
1:A:34:VAL:HA	1:A:46:GLU:O	2.16	0.44
4:D:22:VAL:HG22	4:D:23:THR:N	2.33	0.44
4:D:22:VAL:O	4:D:62:ILE:HB	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:SER:HA	1:A:21:ARG:O	2.17	0.44
1:A:145:HIS:O	1:A:149:ALA:CB	2.65	0.44
4:D:34:GLU:O	4:D:77:TYR:CB	2.63	0.44
4:D:40:GLU:C	4:D:42:LYS:H	2.21	0.44
1:A:49:ALA:HB1	1:A:51:TRP:CE2	2.53	0.44
1:A:185:PRO:O	1:A:187:THR:HG23	2.17	0.44
2:B:33:SER:O	2:B:34:ASP:C	2.56	0.44
4:D:93:LEU:O	4:D:183:TYR:HB3	2.18	0.44
1:A:116:TYR:CD1	1:A:117:ALA:N	2.85	0.44
2:B:24:ASN:HB3	2:B:65:LEU:HD11	2.00	0.44
1:A:81:LEU:HG	1:A:118:TYR:CG	2.53	0.44
2:B:25:CYS:HB2	2:B:39:LEU:HD21	2.00	0.44
4:D:118:VAL:CG2	4:D:119:THR:N	2.46	0.44
1:A:81:LEU:HD12	1:A:118:TYR:CD2	2.53	0.43
2:B:88:SER:HB3	4:D:187:LEU:HD11	2.00	0.43
1:A:59:TYR:O	1:A:63:GLU:HG2	2.18	0.43
2:B:72:PRO:O	2:B:73:THR:C	2.56	0.43
1:A:242:GLN:NE2	2:B:12:ARG:O	2.51	0.43
2:B:58:LYS:C	2:B:60:TRP:H	2.21	0.43
1:A:54:GLN:HE22	1:A:174:ASN:CG	2.22	0.43
1:A:97:ARG:NH1	3:C:6:VAL:CG2	2.82	0.43
1:A:64:THR:O	1:A:68:LYS:HG2	2.17	0.43
4:D:63:PRO:HB2	4:D:64:SER:H	1.66	0.43
4:D:146:ASN:OD1	4:D:147:SER:N	2.52	0.43
4:D:182:PRO:C	4:D:184:GLU:N	2.70	0.43
4:D:97:GLY:HA2	4:D:187:LEU:CG	2.49	0.43
1:A:13:SER:HA	1:A:20:PRO:CB	2.45	0.43
1:A:9:PHE:HZ	3:C:2:LEU:CD2	2.32	0.43
2:B:12:ARG:NH2	2:B:22:PHE:CD1	2.87	0.43
4:D:107:GLN:HE21	4:D:107:GLN:CA	2.22	0.43
1:A:4:SER:HB2	1:A:102:ASP:OD1	2.19	0.42
1:A:84:TYR:H	1:A:84:TYR:HD1	1.66	0.42
4:D:10:ALA:O	4:D:13:GLY:HA2	2.19	0.42
4:D:97:GLY:CA	4:D:187:LEU:HD23	2.50	0.42
1:A:192:HIS:NE2	2:B:98:ASP:HB2	2.34	0.42
4:D:10:ALA:HB3	4:D:13:GLY:HA2	2.02	0.42
1:A:249:VAL:HG13	1:A:250:PRO:HD2	2.00	0.42
2:B:35:ILE:HG12	2:B:36:GLU:N	2.34	0.42
4:D:18:GLN:HA	4:D:65:ILE:HG23	2.02	0.42
1:A:28:VAL:O	1:A:29:ASP:HB2	2.19	0.42
4:D:151:ALA:HB3	4:D:154:SER:OG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:9:VAL:HG12	2:B:95:TRP:HB2	2.02	0.42
1:A:126:LEU:HD12	1:A:132:SER:O	2.19	0.42
1:A:172:LEU:HA	1:A:179:LEU:HD12	2.02	0.42
1:A:253:GLN:OE1	1:A:256:ARG:HD2	2.20	0.42
1:A:57:PRO:O	1:A:58:GLU:C	2.58	0.42
2:B:49:VAL:HG12	2:B:66:TYR:HD2	1.85	0.42
1:A:99:TYR:OH	3:C:2:LEU:HD22	2.20	0.42
4:D:167:SER:C	4:D:168:ARG:HG2	2.39	0.42
4:D:18:GLN:HG2	4:D:19:GLY:N	2.35	0.42
1:A:200:THR:HG22	1:A:201:LEU:N	2.33	0.42
2:B:84:HIS:CG	2:B:85:VAL:N	2.87	0.42
4:D:76:TYR:CD1	4:D:76:TYR:C	2.92	0.42
1:A:117:ALA:HA	1:A:123:TYR:CB	2.50	0.42
1:A:138:MET:HA	1:A:141:GLN:HE22	1.84	0.42
1:A:189:MET:CE	1:A:217:TRP:HH2	2.33	0.42
2:B:59:ASP:O	2:B:60:TRP:HB2	2.19	0.42
1:A:215:LEU:HD12	1:A:243:LYS:HG2	2.02	0.42
1:A:47:PRO:HB3	1:A:60:TRP:CH2	2.55	0.42
1:A:74:HIS:HA	1:A:77:ASP:OD2	2.20	0.42
2:B:81:ARG:HG3	2:B:82:VAL:N	2.33	0.42
1:A:201:LEU:HD12	1:A:249:VAL:HG21	2.02	0.41
1:A:9:PHE:CE2	1:A:99:TYR:HE2	2.32	0.41
2:B:85:VAL:O	4:D:100:ILE:HD12	2.20	0.41
2:B:96:ASP:HB3	2:B:99:MET:CA	2.50	0.41
4:D:151:ALA:HB3	4:D:154:SER:CB	2.50	0.41
4:D:145:LEU:HD12	4:D:146:ASN:N	2.35	0.41
2:B:12:ARG:NH2	2:B:22:PHE:CG	2.88	0.41
3:C:3:LYS:HE2	3:C:5:PRO:O	2.20	0.41
4:D:26:CYS:O	4:D:27:GLN:C	2.57	0.41
1:A:145:HIS:O	1:A:149:ALA:HB3	2.20	0.41
1:A:233:THR:N	1:A:243:LYS:HZ3	2.18	0.41
1:A:87:GLN:HB2	1:A:87:GLN:HE21	1.67	0.41
4:D:178:ASP:C	4:D:180:ASN:N	2.72	0.41
1:A:232:GLU:OE2	2:B:28:SER:OG	2.37	0.41
2:B:92:ILE:HG22	2:B:93:VAL:N	2.36	0.41
4:D:136:GLU:CD	4:D:173:ARG:HD2	2.41	0.41
4:D:97:GLY:HA2	4:D:187:LEU:CD2	2.51	0.41
1:A:215:LEU:HD23	1:A:261:VAL:HG22	2.02	0.41
4:D:36:ARG:HD2	4:D:76:TYR:HE1	1.85	0.41
1:A:36:PHE:CG	1:A:67:VAL:HG21	2.56	0.41
2:B:6:LYS:HB2	2:B:28:SER:OG	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1:ILE:C	3:C:1:ILE:HD12	2.41	0.41
1:A:202:ARG:HD3	1:A:244:TRP:CD2	2.56	0.41
1:A:76:VAL:CG2	1:A:77:ASP:N	2.83	0.41
4:D:62:ILE:HG22	4:D:62:ILE:O	2.20	0.41
2:B:46:ILE:HG22	2:B:47:GLU:O	2.21	0.41
1:A:129:ASP:OD1	1:A:131:ARG:N	2.54	0.41
1:A:135:ALA:O	1:A:137:ASP:N	2.54	0.41
1:A:167:TRP:CG	3:C:1:ILE:HG13	2.55	0.41
1:A:104:GLY:C	1:A:106:ASP:N	2.74	0.40
4:D:107:GLN:O	4:D:119:THR:HB	2.21	0.40
4:D:20:SER:HB3	4:D:21:PRO:CD	2.51	0.40
1:A:19:GLU:HB2	1:A:20:PRO:HD2	2.04	0.40
1:A:249:VAL:HG13	1:A:257:TYR:CE1	2.57	0.40
2:B:4:THR:CG2	2:B:5:PRO:N	2.84	0.40
1:A:173:GLU:CD	1:A:176:LYS:HE2	2.42	0.40
2:B:93:VAL:HG22	4:D:67:TRP:HZ2	1.86	0.40
4:D:120:LEU:N	4:D:120:LEU:CD2	2.83	0.40
1:A:8:PHE:HB2	1:A:25:VAL:HG13	2.02	0.40
1:A:8:PHE:CD1	1:A:8:PHE:N	2.88	0.40
2:B:86:THR:HG23	4:D:126:VAL:CG2	2.51	0.40
4:D:166:PRO:C	4:D:168:ARG:H	2.24	0.40
4:D:67:TRP:HA	4:D:95:VAL:HG21	2.03	0.40
1:A:138:MET:SD	1:A:138:MET:O	2.79	0.40
1:A:44:ARG:HG3	1:A:64:THR:OG1	2.22	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	274/276 (99%)	224 (82%)	42 (15%)	8 (3%)	4 24

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	87/99 (88%)	71 (82%)	12 (14%)	4 (5%)	2	15
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
4	D	175/197 (89%)	145 (83%)	19 (11%)	11 (6%)	1	9
All	All	543/581 (94%)	446 (82%)	74 (14%)	23 (4%)	3	18

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	SER
1	A	137	ASP
4	D	63	PRO
4	D	181	SER
1	A	136	ALA
1	A	150	ALA
1	A	213	ILE
2	B	34	ASP
4	D	21	PRO
1	A	267	PRO
2	B	70	PHE
4	D	14	SER
4	D	43	THR
4	D	70	ALA
4	D	179	SER
1	A	86	ASN
4	D	41	LYS
4	D	145	LEU
1	A	58	GLU
2	B	49	VAL
4	D	40	GLU
2	B	72	PRO
4	D	6	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	229/232 (99%)	212 (93%)	17 (7%)	13	42
2	B	87/94 (93%)	81 (93%)	6 (7%)	15	45
3	C	8/8 (100%)	7 (88%)	1 (12%)	4	17
4	D	152/171 (89%)	137 (90%)	15 (10%)	8	27
All	All	476/505 (94%)	437 (92%)	39 (8%)	11	37

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	35	ARG
1	A	54	GLN
1	A	81	LEU
1	A	137	ASP
1	A	141	GLN
1	A	155	GLN
1	A	163	THR
1	A	165	VAL
1	A	177	GLU
1	A	183	ASP
1	A	209	TYR
1	A	214	THR
1	A	225	THR
1	A	230	LEU
1	A	232	GLU
1	A	276	PRO
2	B	34	ASP
2	B	55	SER
2	B	67	TYR
2	B	70	PHE
2	B	81	ARG
2	B	98	ASP
3	C	1	ILE
4	D	17	THR
4	D	21	PRO
4	D	75	CYS
4	D	76	TYR
4	D	77	TYR
4	D	105	SER
4	D	108	PRO
4	D	141	HIS
4	D	154	SER

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Mol	Chain	Res	Type
4	D	164	VAL
4	D	180	ASN
4	D	181	SER
4	D	191	LEU
4	D	196	VAL
4	D	197	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	54	GLN
1	A	87	GLN
1	A	141	GLN
1	A	151	HIS
1	A	155	GLN
1	A	224	GLN
1	A	255	GLN
2	B	21	ASN
2	B	83	ASN
3	C	7	HIS
4	D	18	GLN
4	D	69	HIS
4	D	107	GLN
4	D	141	HIS
4	D	148	GLN
4	D	150	HIS

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 ⓘ

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	225:THR	C	226:GLN	N	1.09

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	276/276 (100%)	-0.55	0 100 100	2, 26, 71, 117	0
2	B	93/99 (93%)	-0.39	0 100 100	2, 30, 82, 108	0
3	C	9/9 (100%)	-0.70	0 100 100	5, 16, 29, 46	0
4	D	183/197 (92%)	-0.02	8 (4%) 34 34	8, 49, 111, 153	0
All	All	561/581 (96%)	-0.36	8 (1%) 75 74	2, 32, 89, 153	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	61	PRO	3.6
4	D	27	GLN	3.2
4	D	37	LEU	3.2
4	D	26	CYS	2.8
4	D	60	PHE	2.7
4	D	76	TYR	2.4
4	D	59	GLN	2.3
4	D	167	SER	2.1

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 [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.