



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 08:29 PM UTC

PDB ID : 3VDP / pdb_00003vdp
Title : Structure and biochemical studies of the recombination mediator protein RecR
in RecFOR pathway
Authors : Tang, Q.; Yan, X.X.; Liang, D.C.
Deposited on : 2012-01-05
Resolution : 2.45 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

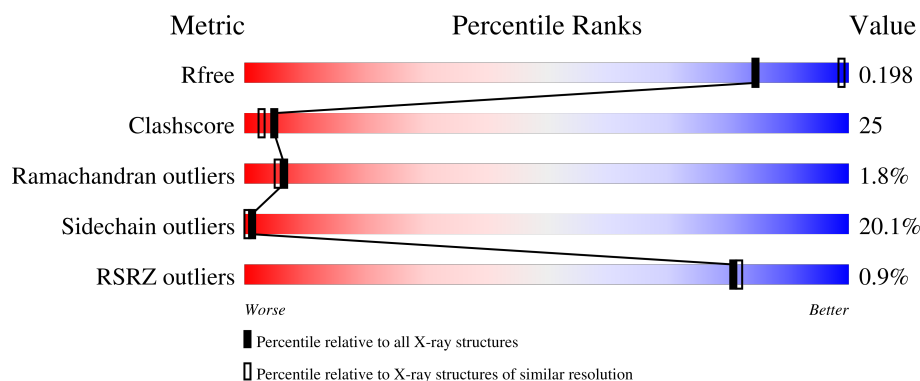
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	212	
1	B	212	
1	C	212	
1	D	212	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	IMD	A	204	-	-	X	-
3	IMD	B	202	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6481 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 Recombination protein recR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	197	Total	C	N	O	S	0	0	0
			1535	976	259	291	9			
1	B	195	Total	C	N	O	S	0	0	0
			1504	955	256	284	9			
1	C	197	Total	C	N	O	S	0	0	0
			1535	976	259	291	9			
1	D	199	Total	C	N	O	S	0	0	0
			1548	984	261	293	10			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	GLY	-	expression tag	UNP Q8RDI4
A	-14	SER	-	expression tag	UNP Q8RDI4
A	-13	SER	-	expression tag	UNP Q8RDI4
A	-12	HIS	-	expression tag	UNP Q8RDI4
A	-11	HIS	-	expression tag	UNP Q8RDI4
A	-10	HIS	-	expression tag	UNP Q8RDI4
A	-9	HIS	-	expression tag	UNP Q8RDI4
A	-8	HIS	-	expression tag	UNP Q8RDI4
A	-7	HIS	-	expression tag	UNP Q8RDI4
A	-6	SER	-	expression tag	UNP Q8RDI4
A	-5	GLN	-	expression tag	UNP Q8RDI4
A	-4	ASP	-	expression tag	UNP Q8RDI4
A	-3	PRO	-	expression tag	UNP Q8RDI4
B	-15	GLY	-	expression tag	UNP Q8RDI4
B	-14	SER	-	expression tag	UNP Q8RDI4
B	-13	SER	-	expression tag	UNP Q8RDI4
B	-12	HIS	-	expression tag	UNP Q8RDI4
B	-11	HIS	-	expression tag	UNP Q8RDI4
B	-10	HIS	-	expression tag	UNP Q8RDI4
B	-9	HIS	-	expression tag	UNP Q8RDI4
B	-8	HIS	-	expression tag	UNP Q8RDI4

Continued on next page...

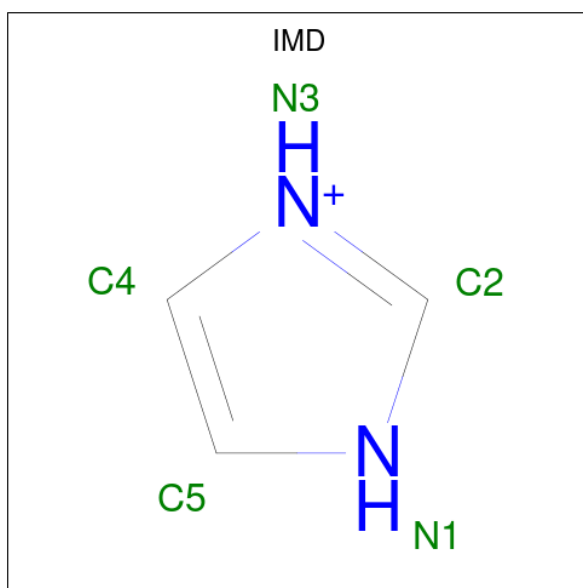
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP Q8RDI4
B	-6	SER	-	expression tag	UNP Q8RDI4
B	-5	GLN	-	expression tag	UNP Q8RDI4
B	-4	ASP	-	expression tag	UNP Q8RDI4
B	-3	PRO	-	expression tag	UNP Q8RDI4
C	-15	GLY	-	expression tag	UNP Q8RDI4
C	-14	SER	-	expression tag	UNP Q8RDI4
C	-13	SER	-	expression tag	UNP Q8RDI4
C	-12	HIS	-	expression tag	UNP Q8RDI4
C	-11	HIS	-	expression tag	UNP Q8RDI4
C	-10	HIS	-	expression tag	UNP Q8RDI4
C	-9	HIS	-	expression tag	UNP Q8RDI4
C	-8	HIS	-	expression tag	UNP Q8RDI4
C	-7	HIS	-	expression tag	UNP Q8RDI4
C	-6	SER	-	expression tag	UNP Q8RDI4
C	-5	GLN	-	expression tag	UNP Q8RDI4
C	-4	ASP	-	expression tag	UNP Q8RDI4
C	-3	PRO	-	expression tag	UNP Q8RDI4
D	-15	GLY	-	expression tag	UNP Q8RDI4
D	-14	SER	-	expression tag	UNP Q8RDI4
D	-13	SER	-	expression tag	UNP Q8RDI4
D	-12	HIS	-	expression tag	UNP Q8RDI4
D	-11	HIS	-	expression tag	UNP Q8RDI4
D	-10	HIS	-	expression tag	UNP Q8RDI4
D	-9	HIS	-	expression tag	UNP Q8RDI4
D	-8	HIS	-	expression tag	UNP Q8RDI4
D	-7	HIS	-	expression tag	UNP Q8RDI4
D	-6	SER	-	expression tag	UNP Q8RDI4
D	-5	GLN	-	expression tag	UNP Q8RDI4
D	-4	ASP	-	expression tag	UNP Q8RDI4
D	-3	PRO	-	expression tag	UNP Q8RDI4

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

- Molecule 3 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			5	3	2		
3	A	1	Total	C	N	0	0
			5	3	2		
3	A	1	Total	C	N	0	0
			5	3	2		
3	B	1	Total	C	N	0	0
			5	3	2		
3	C	1	Total	C	N	0	0
			5	3	2		
3	D	1	Total	C	N	0	0
			5	3	2		

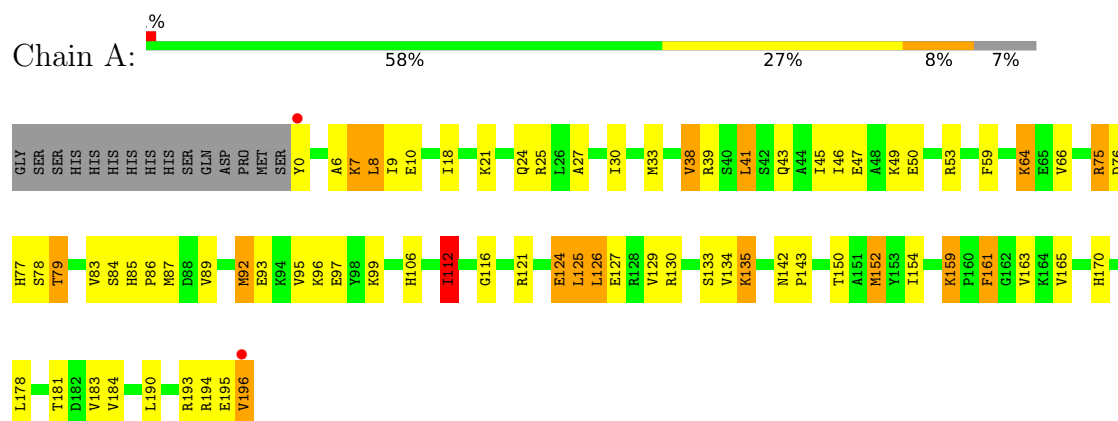
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	92	Total	O	0	0
			92	92		
4	B	74	Total	O	0	0
			74	74		
4	C	74	Total	O	0	0
			74	74		
4	D	85	Total	O	0	0
			85	85		

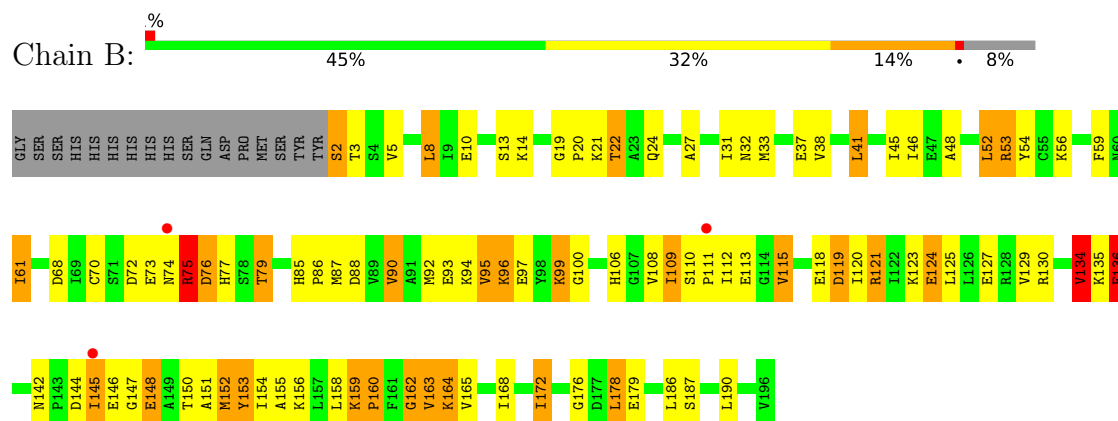
3 Residue-property plots [i](#)

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

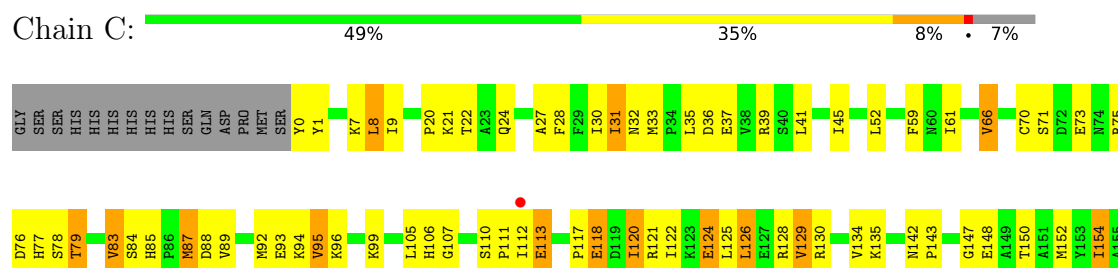
• Molecule 1: Recombination protein recR

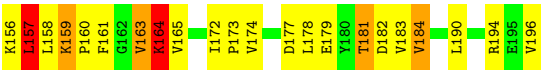


• Molecule 1: Recombination protein recR

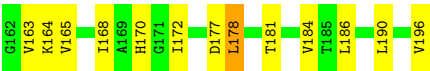


• Molecule 1: Recombination protein recR





• Molecule 1: Recombination protein recR



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.26Å 68.33Å 94.06Å 90.00° 92.95° 90.00°	Depositor
Resolution (Å)	19.96 – 2.45 19.96 – 2.45	Depositor EDS
% Data completeness (in resolution range)	97.6 (19.96-2.45) 97.6 (19.96-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.73 (at 2.45Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.196 , 0.227 0.197 , 0.198	Depositor DCC
R_{free} test set	2096 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.957	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 57.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.000 for h,-k,-l 0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6481	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.8015e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

Bond lengths and bond angles in the following residue types are not validated in this section: IMD, ZN

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	0.77	0/1559	1.10	7/2105 (0.3%)
1	B	0.75	0/1526	1.11	7/2060 (0.3%)
1	C	0.75	1/1559 (0.1%)	1.14	3/2105 (0.1%)
1	D	0.73	0/1572	1.08	6/2122 (0.3%)
All	All	0.75	1/6216 (0.0%)	1.11	23/8392 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	129	VAL	CA-CB	5.28	1.61	1.54

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	GLU	N-CA-C	-8.69	96.60	108.86
1	B	119	ASP	N-CA-C	-7.52	104.14	113.38
1	C	85	HIS	CA-C-N	6.87	127.44	119.47
1	C	85	HIS	C-N-CA	6.87	127.44	119.47
1	A	125	LEU	N-CA-C	-6.33	104.46	111.36
1	A	112	ILE	CB-CA-C	-6.33	103.73	112.02
1	B	172	ILE	CB-CA-C	-6.31	105.07	110.94
1	B	136	GLU	N-CA-C	6.12	123.84	110.80
1	B	53	ARG	N-CA-C	5.50	119.48	112.12
1	A	161	PHE	N-CA-C	-5.41	102.47	110.48
1	D	107	GLY	N-CA-C	5.41	118.19	110.46
1	D	136	GLU	N-CA-C	5.40	117.61	109.23
1	D	116	GLY	CA-C-N	5.40	125.48	119.32
1	D	116	GLY	C-N-CA	5.40	125.48	119.32
1	B	134	VAL	N-CA-C	5.36	116.02	108.36
1	A	75	ARG	N-CA-C	-5.32	100.73	109.40

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	46	ILE	N-CA-C	5.30	116.08	110.72
1	D	29	PHE	N-CA-C	-5.24	105.24	111.69
1	C	84	SER	N-CA-C	5.17	118.74	112.23
1	D	142	ASN	N-CA-C	5.16	116.19	109.64
1	B	115	VAL	N-CA-C	5.15	115.53	108.12
1	A	116	GLY	CA-C-N	5.01	124.79	119.28
1	A	116	GLY	C-N-CA	5.01	124.79	119.28

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	1535	0	1583	76	0
1	B	1504	0	1554	103	0
1	C	1535	0	1583	79	0
1	D	1548	0	1594	91	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	15	0	15	11	0
3	B	5	0	5	7	0
3	C	5	0	5	3	0
3	D	5	0	5	1	0
4	A	92	0	0	4	0
4	B	74	0	0	14	0
4	C	74	0	0	8	0
4	D	85	0	0	13	0
All	All	6481	0	6344	315	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 25.

All (315) 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)
1:A:87:MET:HB2	3:A:204:IMD:H5	1.37	1.06
1:C:163:VAL:O	1:C:164:LYS:HB3	1.49	1.05
1:A:112:ILE:HD13	1:A:112:ILE:H	1.24	1.01
1:B:85:HIS:HD2	1:B:87:MET:H	1.03	1.01
1:D:99:LYS:CD	1:D:99:LYS:H	1.75	0.97
1:B:134:VAL:HG23	1:B:135:LYS:O	1.65	0.96
1:B:20:PRO:HD2	3:B:202:IMD:H5	1.50	0.93
1:C:182:ASP:OD2	1:C:184:VAL:HG23	1.70	0.91
1:A:112:ILE:HD13	1:A:112:ILE:N	1.87	0.90
1:B:14:LYS:HE2	4:C:374:HOH:O	1.72	0.89
1:B:76:ASP:OD2	1:B:79:THR:HG23	1.73	0.88
1:B:146:GLU:HG2	4:B:365:HOH:O	1.73	0.88
1:D:8:LEU:HD13	1:D:27:ALA:HA	1.56	0.87
1:B:85:HIS:CD2	1:B:87:MET:H	1.93	0.85
1:A:30:ILE:HA	1:A:33:MET:HE2	1.57	0.84
1:B:99:LYS:HD3	1:B:99:LYS:H	1.43	0.83
1:C:130:ARG:HD2	4:C:343:HOH:O	1.80	0.80
1:B:52:LEU:HD23	1:B:61:ILE:HG12	1.63	0.80
1:B:22:THR:H	3:B:202:IMD:H4	1.47	0.80
1:B:56:LYS:HE3	4:B:360:HOH:O	1.85	0.77
1:D:99:LYS:H	1:D:99:LYS:HD3	1.49	0.77
1:A:38:VAL:HG11	1:D:49:LYS:HG3	1.67	0.77
1:A:59:PHE:O	1:A:106:HIS:HE1	1.68	0.76
1:C:148:GLU:OE1	1:C:152:MET:HE3	1.85	0.76
1:D:170:HIS:HE1	4:D:369:HOH:O	1.68	0.76
1:C:59:PHE:O	1:C:106:HIS:HE1	1.69	0.76
1:B:85:HIS:HD2	1:B:87:MET:N	1.84	0.75
1:C:0:TYR:HD2	1:C:1:TYR:H	1.32	0.75
1:D:99:LYS:H	1:D:99:LYS:HD2	1.49	0.75
1:A:127:GLU:HG2	1:A:130:ARG:HH22	1.49	0.74
1:D:-2:MET:HG2	4:D:379:HOH:O	1.88	0.74
1:B:41:LEU:HD13	1:C:45:ILE:HD13	1.69	0.73
1:A:92:MET:HE1	1:B:190:LEU:HD22	1.70	0.72
1:D:95:VAL:HG22	4:D:376:HOH:O	1.89	0.72
1:C:112:ILE:HG22	1:C:113:GLU:HG3	1.72	0.72
1:C:76:ASP:OD2	1:C:79:THR:HG23	1.90	0.72
1:A:112:ILE:HD11	4:A:386:HOH:O	1.90	0.71
1:D:112:ILE:O	1:D:113:GLU:HB2	1.91	0.70
1:B:74:ASN:O	1:B:75:ARG:C	2.35	0.69
1:D:75:ARG:NH1	1:D:100:GLY:O	2.26	0.69
1:D:59:PHE:O	1:D:106:HIS:HE1	1.75	0.68
1:C:122:ILE:O	1:C:126:LEU:HD22	1.93	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-2:MET:HE3	4:D:378:HOH:O	1.92	0.68
1:D:85:HIS:HD2	1:D:87:MET:HB2	1.59	0.68
1:C:30:ILE:HA	1:C:33:MET:HE2	1.76	0.67
3:A:204:IMD:H4	1:D:28:PHE:HB3	1.76	0.67
1:D:21:LYS:HE3	1:D:25:ARG:NH2	2.09	0.67
1:A:75:ARG:O	1:A:77:HIS:HD2	1.78	0.67
1:D:190:LEU:HD23	1:D:190:LEU:C	2.20	0.67
1:D:152:MET:HE3	1:D:152:MET:HA	1.75	0.67
1:B:72:ASP:HB3	1:B:75:ARG:HG2	1.78	0.66
1:C:107:GLY:HA3	1:C:120:ILE:HD11	1.77	0.66
1:C:163:VAL:O	1:C:164:LYS:CB	2.33	0.66
1:C:126:LEU:HD11	1:C:157:LEU:HD23	1.78	0.65
1:C:177:ASP:O	1:C:181:THR:HG23	1.97	0.65
1:D:85:HIS:CD2	1:D:87:MET:HB2	2.30	0.65
1:D:85:HIS:CD2	1:D:87:MET:H	2.15	0.65
1:B:152:MET:HA	1:B:152:MET:HE2	1.79	0.65
1:B:121:ARG:HD3	4:B:339:HOH:O	1.97	0.64
1:D:77:HIS:HD2	4:D:361:HOH:O	1.81	0.64
1:B:99:LYS:H	1:B:99:LYS:CD	2.10	0.64
1:C:21:LYS:H	3:C:202:IMD:HN3	1.44	0.64
1:B:59:PHE:O	1:B:106:HIS:HE1	1.79	0.64
1:D:121:ARG:HD2	1:D:121:ARG:N	2.12	0.64
1:D:85:HIS:HD2	1:D:87:MET:H	1.47	0.63
1:B:2:SER:HB3	4:B:362:HOH:O	1.98	0.63
1:A:143:PRO:HD2	1:B:176:GLY:O	1.98	0.62
1:A:95:VAL:HG13	1:B:187:SER:HB2	1.80	0.62
1:C:75:ARG:HD3	1:C:99:LYS:O	1.99	0.62
1:D:145:ILE:HD12	1:D:145:ILE:H	1.64	0.62
1:B:109:ILE:HD11	1:B:150:THR:HA	1.81	0.62
1:D:85:HIS:HD2	1:D:87:MET:CB	2.13	0.62
1:B:111:PRO:HB3	4:B:368:HOH:O	1.98	0.62
1:C:79:THR:HG21	4:C:353:HOH:O	1.99	0.61
1:B:129:VAL:HG12	1:B:129:VAL:O	1.99	0.61
1:A:46:ILE:O	1:A:50:GLU:HG3	2.01	0.61
1:A:87:MET:HE3	1:D:24:GLN:OE1	2.00	0.61
1:A:112:ILE:H	1:A:112:ILE:CD1	1.95	0.60
1:A:85:HIS:HB3	3:A:204:IMD:N1	2.16	0.60
1:C:22:THR:OG1	3:C:202:IMD:H2	2.02	0.60
1:A:59:PHE:CE2	1:A:124:GLU:HG3	2.36	0.60
1:B:85:HIS:CD2	1:B:87:MET:HG3	2.37	0.60
1:D:8:LEU:CD1	1:D:27:ALA:HA	2.28	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:HIS:HB3	3:A:204:IMD:HN1	1.66	0.60
1:B:70:CYS:O	1:B:77:HIS:HE1	1.85	0.60
1:A:152:MET:HA	1:A:152:MET:HE3	1.82	0.60
1:D:76:ASP:OD2	1:D:79:THR:CG2	2.50	0.60
1:A:178:LEU:H	1:B:142:ASN:HD22	1.50	0.59
1:A:87:MET:HE1	1:D:25:ARG:HA	1.83	0.59
1:B:85:HIS:NE2	1:B:87:MET:HG3	2.17	0.59
1:D:121:ARG:HD3	4:D:375:HOH:O	2.01	0.59
1:A:87:MET:HE1	1:D:25:ARG:HG2	1.84	0.59
1:A:87:MET:HE3	1:D:24:GLN:HG2	1.85	0.59
1:D:112:ILE:O	1:D:113:GLU:CB	2.50	0.59
1:D:99:LYS:CD	1:D:99:LYS:N	2.55	0.59
1:B:2:SER:OG	1:B:5:VAL:HG23	2.03	0.58
1:D:134:VAL:O	1:D:163:VAL:HG12	2.03	0.58
1:D:143:PRO:HA	4:D:369:HOH:O	2.02	0.58
1:B:159:LYS:C	1:B:160:PRO:O	2.45	0.58
1:B:75:ARG:NH1	1:B:100:GLY:O	2.37	0.58
1:D:172:ILE:CD1	1:D:181:THR:HG21	2.34	0.58
1:B:52:LEU:CD2	1:B:61:ILE:HG12	2.33	0.58
1:B:19:GLY:HA3	3:B:202:IMD:C5	2.34	0.58
1:A:112:ILE:N	1:A:112:ILE:CD1	2.60	0.57
1:A:134:VAL:H	3:A:202:IMD:H5	1.69	0.57
1:C:194:ARG:HB3	4:C:306:HOH:O	2.04	0.57
1:A:127:GLU:HG2	1:A:130:ARG:NH2	2.19	0.57
1:B:121:ARG:HG3	1:B:124:GLU:CG	2.34	0.57
1:C:142:ASN:O	1:C:147:GLY:HA3	2.05	0.57
1:B:110:SER:HB3	1:B:146:GLU:OE1	2.05	0.57
1:A:152:MET:HA	1:A:152:MET:CE	2.35	0.56
1:B:8:LEU:HD13	1:B:27:ALA:HA	1.87	0.56
1:A:142:ASN:ND2	1:B:178:LEU:H	2.03	0.56
1:B:48:ALA:O	1:B:52:LEU:HB3	2.05	0.56
1:B:52:LEU:C	1:B:52:LEU:HD13	2.30	0.56
1:B:41:LEU:CD1	1:C:45:ILE:HD13	2.35	0.56
1:B:41:LEU:O	1:B:45:ILE:HG13	2.05	0.56
1:D:49:LYS:O	1:D:52:LEU:HD22	2.06	0.56
1:B:41:LEU:O	1:B:41:LEU:HD22	2.05	0.56
1:B:52:LEU:HD13	1:B:52:LEU:O	2.06	0.56
1:D:133:SER:O	1:D:134:VAL:HB	2.06	0.56
1:A:87:MET:HB2	3:A:204:IMD:C5	2.25	0.56
1:B:20:PRO:CD	3:B:202:IMD:H5	2.31	0.56
1:B:14:LYS:CD	4:B:373:HOH:O	2.53	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:MET:HA	1:B:152:MET:CE	2.36	0.55
1:D:79:THR:HG22	1:D:135:LYS:HE3	1.88	0.55
1:C:159:LYS:HG3	1:C:160:PRO:HD3	1.88	0.55
1:C:172:ILE:CD1	1:C:181:THR:HG21	2.36	0.55
1:B:134:VAL:O	1:B:163:VAL:HG12	2.07	0.55
1:B:159:LYS:O	1:B:160:PRO:O	2.24	0.55
1:C:105:LEU:O	1:C:120:ILE:HG13	2.06	0.54
1:A:150:THR:O	1:A:154:ILE:HG13	2.07	0.54
1:B:121:ARG:HG3	1:B:124:GLU:CD	2.32	0.54
1:D:118:GLU:CD	1:D:118:GLU:H	2.15	0.54
1:A:30:ILE:CA	1:A:33:MET:HE2	2.32	0.54
1:B:61:ILE:O	1:B:61:ILE:HG22	2.08	0.54
1:A:43:GLN:O	1:A:47:GLU:HG3	2.07	0.54
1:D:92:MET:O	1:D:95:VAL:HG13	2.08	0.54
1:C:20:PRO:HD2	3:C:202:IMD:C4	2.38	0.54
1:B:14:LYS:HD2	4:B:373:HOH:O	2.07	0.54
1:B:22:THR:N	3:B:202:IMD:H4	2.21	0.54
1:D:21:LYS:HE3	1:D:25:ARG:HH22	1.73	0.54
1:D:22:THR:OG1	3:D:202:IMD:H2	2.08	0.54
1:D:53:ARG:HB2	4:D:365:HOH:O	2.08	0.53
1:B:121:ARG:HG3	1:B:124:GLU:HG3	1.91	0.53
1:B:129:VAL:O	1:B:129:VAL:CG1	2.55	0.53
1:C:9:ILE:HD13	1:C:24:GLN:HG2	1.90	0.53
1:D:112:ILE:O	1:D:112:ILE:HG23	2.08	0.53
1:D:106:HIS:O	1:D:120:ILE:HG22	2.09	0.53
1:A:142:ASN:HD21	1:B:178:LEU:H	1.56	0.53
1:C:83:VAL:HG13	1:C:88:ASP:HB2	1.91	0.53
1:B:136:GLU:HA	1:B:164:LYS:O	2.09	0.53
1:C:52:LEU:HD21	1:C:61:ILE:HG12	1.90	0.53
1:D:83:VAL:HG13	1:D:88:ASP:HB2	1.91	0.53
1:D:98:TYR:CZ	1:D:100:GLY:HA3	2.44	0.53
1:B:85:HIS:HE2	1:B:87:MET:HG3	1.74	0.52
1:A:21:LYS:HB2	3:A:203:IMD:C2	2.39	0.52
1:A:87:MET:HE1	1:D:25:ARG:CA	2.39	0.52
1:A:178:LEU:H	1:B:142:ASN:ND2	2.07	0.52
1:C:126:LEU:CD1	1:C:158:LEU:HD13	2.40	0.52
1:C:142:ASN:OD1	1:D:178:LEU:HD23	2.10	0.52
1:D:94:LYS:O	1:D:96:LYS:HD3	2.10	0.51
1:A:87:MET:HE3	1:D:24:GLN:CG	2.40	0.51
1:B:142:ASN:O	1:B:147:GLY:HA3	2.11	0.51
1:D:152:MET:HA	1:D:152:MET:CE	2.33	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-2:MET:CG	4:D:379:HOH:O	2.52	0.51
1:A:45:ILE:HD13	1:D:41:LEU:HD13	1.93	0.51
1:A:129:VAL:CG1	1:A:163:VAL:HG21	2.40	0.51
1:A:95:VAL:HG12	1:A:95:VAL:O	2.10	0.51
1:B:21:LYS:HB2	3:B:202:IMD:C4	2.40	0.51
1:C:92:MET:O	1:C:95:VAL:HG13	2.11	0.51
1:B:53:ARG:O	1:B:54:TYR:HB2	2.11	0.50
1:A:8:LEU:HD13	1:A:27:ALA:HA	1.93	0.50
1:D:178:LEU:HD23	1:D:178:LEU:H	1.76	0.50
1:A:33:MET:O	1:D:49:LYS:NZ	2.44	0.50
1:C:75:ARG:O	1:C:77:HIS:HD2	1.95	0.50
1:A:39:ARG:NH1	4:A:345:HOH:O	2.44	0.50
1:C:95:VAL:HG22	1:C:95:VAL:O	2.10	0.50
1:B:115:VAL:HG13	1:B:119:ASP:HB3	1.92	0.50
1:D:110:SER:O	1:D:112:ILE:N	2.37	0.50
1:C:172:ILE:HG13	1:C:173:PRO:HD2	1.93	0.50
1:B:144:ASP:O	1:B:148:GLU:HB2	2.12	0.49
1:C:150:THR:O	1:C:154:ILE:HD12	2.11	0.49
1:B:112:ILE:O	1:B:112:ILE:HG13	2.11	0.49
1:B:160:PRO:C	1:B:162:GLY:H	2.20	0.49
1:A:38:VAL:CG1	1:D:49:LYS:HG3	2.41	0.49
1:A:85:HIS:HB3	3:A:204:IMD:C2	2.43	0.49
1:B:74:ASN:O	1:B:75:ARG:O	2.30	0.49
1:C:66:VAL:CG2	1:C:71:SER:HB3	2.43	0.49
1:D:110:SER:C	1:D:112:ILE:H	2.20	0.49
1:C:156:LYS:HD3	4:C:355:HOH:O	2.12	0.49
1:C:30:ILE:HG12	1:C:33:MET:HE2	1.95	0.48
1:A:75:ARG:O	1:A:77:HIS:CD2	2.62	0.48
1:B:24:GLN:HG3	4:B:351:HOH:O	2.12	0.48
1:B:90:VAL:HG23	1:B:94:LYS:HE2	1.96	0.48
1:B:33:MET:HE3	1:B:37:GLU:HB3	1.95	0.48
1:B:41:LEU:HD22	1:B:41:LEU:C	2.38	0.48
1:C:1:TYR:CD2	1:C:28:PHE:HZ	2.32	0.48
1:D:52:LEU:HD23	1:D:52:LEU:O	2.13	0.48
1:D:99:LYS:HD3	1:D:99:LYS:N	2.24	0.48
1:D:76:ASP:OD2	1:D:79:THR:HG23	2.14	0.48
1:A:43:GLN:HA	1:A:46:ILE:HG22	1.96	0.47
1:C:124:GLU:O	1:C:128:ARG:HG3	2.14	0.47
1:A:59:PHE:O	1:A:106:HIS:CE1	2.58	0.47
1:A:142:ASN:HD22	1:B:178:LEU:HD23	1.79	0.47
1:D:62:THR:HG21	1:D:67:CYS:HB2	1.96	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:LYS:HA	1:A:10:GLU:HG2	1.97	0.47
1:B:24:GLN:CD	4:B:351:HOH:O	2.58	0.47
1:C:122:ILE:HG22	1:C:126:LEU:HD21	1.96	0.47
1:B:108:VAL:HG21	1:B:146:GLU:HB3	1.96	0.47
1:C:59:PHE:CZ	1:C:124:GLU:HG3	2.50	0.47
1:A:41:LEU:HD13	1:D:45:ILE:HD13	1.97	0.47
1:D:54:TYR:N	4:D:365:HOH:O	2.47	0.46
1:D:-2:MET:HE1	4:D:377:HOH:O	2.15	0.46
1:C:75:ARG:NH2	1:C:93:GLU:OE1	2.47	0.46
1:D:145:ILE:HD12	1:D:145:ILE:N	2.29	0.46
1:B:145:ILE:HG22	1:B:146:GLU:N	2.29	0.46
1:C:122:ILE:HG22	1:C:126:LEU:CD2	2.45	0.46
1:C:159:LYS:O	1:C:161:PHE:O	2.34	0.46
1:C:0:TYR:HD2	1:C:1:TYR:N	2.07	0.46
1:C:172:ILE:HG12	1:D:143:PRO:HG3	1.97	0.46
1:D:159:LYS:O	1:D:161:PHE:O	2.34	0.46
1:D:10:GLU:O	1:D:14:LYS:HG3	2.16	0.45
1:C:78:SER:OG	1:C:135:LYS:HG2	2.15	0.45
1:A:6:ALA:HB3	4:A:343:HOH:O	2.15	0.45
1:A:89:VAL:O	1:A:93:GLU:HB2	2.16	0.45
1:B:96:LYS:HD3	1:B:96:LYS:N	2.31	0.45
1:B:123:LYS:O	1:B:127:GLU:HG3	2.16	0.45
1:C:143:PRO:HG3	1:D:172:ILE:HG22	1.98	0.45
1:C:196:VAL:CG1	1:D:152:MET:HE1	2.46	0.45
1:A:87:MET:HE3	1:D:24:GLN:CD	2.42	0.45
1:C:36:ASP:OD1	1:C:37:GLU:N	2.50	0.45
1:B:163:VAL:HG23	4:B:361:HOH:O	2.16	0.45
1:B:99:LYS:CD	1:B:99:LYS:N	2.79	0.45
1:A:163:VAL:HG22	3:A:202:IMD:H4	1.98	0.45
1:A:92:MET:HE2	1:A:92:MET:HB3	1.82	0.44
1:B:2:SER:N	4:B:301:HOH:O	2.49	0.44
1:B:85:HIS:CD2	1:B:86:PRO:HD2	2.52	0.44
1:C:31:ILE:HD12	1:C:31:ILE:C	2.42	0.44
1:C:126:LEU:HD12	1:C:158:LEU:HD13	2.00	0.44
1:D:8:LEU:HD23	1:D:8:LEU:HA	1.73	0.44
1:A:78:SER:HB2	1:A:135:LYS:HD2	1.99	0.44
1:A:159:LYS:O	1:A:161:PHE:O	2.34	0.44
1:B:75:ARG:NH2	1:B:93:GLU:OE1	2.50	0.44
1:C:75:ARG:HH22	1:C:93:GLU:CD	2.26	0.44
1:C:179:GLU:HB2	4:C:373:HOH:O	2.17	0.44
1:A:87:MET:HE1	1:D:25:ARG:CG	2.47	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:143:PRO:HG3	1:D:172:ILE:CG2	2.48	0.44
1:A:76:ASP:OD2	1:A:79:THR:CG2	2.66	0.44
1:C:106:HIS:O	1:C:120:ILE:HD11	2.18	0.44
1:C:95:VAL:O	1:C:95:VAL:CG2	2.66	0.43
1:D:8:LEU:HD13	1:D:27:ALA:CA	2.36	0.43
1:A:87:MET:CE	1:D:24:GLN:OE1	2.66	0.43
1:B:53:ARG:O	1:B:54:TYR:CB	2.66	0.43
1:B:110:SER:HB3	1:B:146:GLU:CD	2.43	0.43
1:C:110:SER:HA	1:C:111:PRO:HD2	1.86	0.43
1:A:21:LYS:HB3	1:A:25:ARG:NH2	2.33	0.43
1:C:159:LYS:HE3	1:C:159:LYS:HB2	1.90	0.43
1:D:59:PHE:O	1:D:106:HIS:CE1	2.63	0.43
1:A:18:ILE:O	1:A:18:ILE:HG22	2.17	0.43
1:B:72:ASP:OD1	1:B:73:GLU:N	2.52	0.43
1:C:70:CYS:O	1:C:77:HIS:HE1	2.02	0.43
1:C:76:ASP:OD2	1:C:79:THR:CG2	2.64	0.43
1:C:83:VAL:CG1	1:C:88:ASP:HB2	2.47	0.43
1:D:126:LEU:N	1:D:126:LEU:HD12	2.34	0.43
1:A:195:GLU:O	1:A:196:VAL:OXT	2.36	0.43
1:C:89:VAL:O	1:C:93:GLU:HB2	2.18	0.43
1:C:194:ARG:CB	4:C:306:HOH:O	2.65	0.43
1:C:178:LEU:HD12	1:D:88:ASP:HB3	2.01	0.43
1:C:154:ILE:O	1:C:158:LEU:HB2	2.19	0.43
1:D:85:HIS:CD2	1:D:87:MET:CB	2.96	0.43
1:B:95:VAL:O	1:B:97:GLU:HB2	2.19	0.43
1:A:9:ILE:HD13	1:A:24:GLN:HG2	2.01	0.43
1:A:86:PRO:HD2	3:A:204:IMD:C2	2.48	0.43
1:B:154:ILE:O	1:B:158:LEU:HG	2.18	0.43
1:A:129:VAL:HG13	1:A:163:VAL:HG21	1.99	0.43
1:A:142:ASN:ND2	1:B:178:LEU:HD23	2.33	0.43
1:B:153:TYR:C	1:B:155:ALA:N	2.74	0.42
1:D:68:ASP:HA	4:D:381:HOH:O	2.18	0.42
1:A:95:VAL:O	1:A:95:VAL:CG1	2.67	0.42
1:B:14:LYS:HD3	4:B:373:HOH:O	2.15	0.42
1:C:194:ARG:CG	4:C:306:HOH:O	2.67	0.42
1:D:178:LEU:N	1:D:178:LEU:CD2	2.81	0.42
1:B:153:TYR:C	1:B:153:TYR:CD1	2.97	0.42
1:A:92:MET:HE3	1:A:92:MET:CA	2.49	0.42
1:A:159:LYS:HB3	1:A:159:LYS:NZ	2.33	0.42
1:A:21:LYS:HB2	3:A:203:IMD:H2	2.01	0.42
1:A:92:MET:HE3	1:A:92:MET:HA	2.02	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:ILE:HG12	1:B:150:THR:OG1	2.19	0.42
1:C:117:PRO:HB2	1:C:118:GLU:OE1	2.20	0.42
1:D:66:VAL:HG22	1:D:71:SER:OG	2.20	0.42
4:A:359:HOH:O	1:C:21:LYS:HE2	2.19	0.41
1:C:31:ILE:HD12	1:C:32:ASN:N	2.35	0.41
1:D:129:VAL:HG12	1:D:129:VAL:O	2.19	0.41
1:A:64:LYS:HB2	1:A:64:LYS:HE3	1.67	0.41
1:B:129:VAL:C	4:B:314:HOH:O	2.63	0.41
1:C:94:LYS:O	1:C:96:LYS:HD2	2.19	0.41
1:C:172:ILE:CG1	1:C:173:PRO:HD2	2.50	0.41
1:B:146:GLU:HG2	1:B:146:GLU:H	1.64	0.41
1:A:126:LEU:HD22	1:A:126:LEU:HA	1.93	0.41
1:B:77:HIS:HD2	4:B:349:HOH:O	2.04	0.41
1:B:85:HIS:O	1:B:88:ASP:HB2	2.21	0.41
1:B:22:THR:OG1	3:B:202:IMD:N3	2.54	0.41
1:D:178:LEU:HD23	1:D:178:LEU:N	2.35	0.41
1:D:145:ILE:HA	4:D:371:HOH:O	2.21	0.41
1:A:193:ARG:NH2	1:B:136:GLU:OE1	2.39	0.40
1:C:9:ILE:CD1	1:C:24:GLN:HG2	2.51	0.40
1:B:52:LEU:HD21	1:B:54:TYR:CZ	2.56	0.40
1:C:87:MET:HE2	1:C:87:MET:HB2	1.49	0.40
1:C:174:VAL:HG21	1:D:184:VAL:HG12	2.02	0.40
1:B:99:LYS:HG2	1:B:100:GLY:N	2.37	0.40
1:B:45:ILE:HG12	1:C:8:LEU:HD21	2.04	0.40
1:B:151:ALA:C	1:B:153:TYR:H	2.28	0.40
1:C:8:LEU:HD13	1:C:27:ALA:HA	2.02	0.40
1:D:-2:MET:HE3	1:D:-2:MET:HB2	1.91	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	195/212 (92%)	183 (94%)	11 (6%)	1 (0%)	24	32
1	B	193/212 (91%)	166 (86%)	21 (11%)	6 (3%)	3	1
1	C	195/212 (92%)	177 (91%)	16 (8%)	2 (1%)	12	14
1	D	197/212 (93%)	179 (91%)	13 (7%)	5 (2%)	4	3
All	All	780/848 (92%)	705 (90%)	61 (8%)	14 (2%)	6	5

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	136	GLU
1	B	160	PRO
1	A	96	LYS
1	B	113	GLU
1	B	162	GLY
1	C	164	LYS
1	D	54	TYR
1	D	113	GLU
1	B	75	ARG
1	D	112	ILE
1	B	164	LYS
1	D	134	VAL
1	C	157	LEU
1	D	111	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	172/186 (92%)	141 (82%)	31 (18%)	2	1
1	B	168/186 (90%)	126 (75%)	42 (25%)	0	0
1	C	172/186 (92%)	141 (82%)	31 (18%)	2	1
1	D	173/186 (93%)	139 (80%)	34 (20%)	1	0
All	All	685/744 (92%)	547 (80%)	138 (20%)	1	0

All (138) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	TYR
1	A	7	LYS
1	A	8	LEU
1	A	38	VAL
1	A	41	LEU
1	A	49	LYS
1	A	53	ARG
1	A	64	LYS
1	A	66	VAL
1	A	79	THR
1	A	83	VAL
1	A	84	SER
1	A	92	MET
1	A	99	LYS
1	A	112	ILE
1	A	121	ARG
1	A	124	GLU
1	A	125	LEU
1	A	126	LEU
1	A	133	SER
1	A	135	LYS
1	A	152	MET
1	A	159	LYS
1	A	165	VAL
1	A	170	HIS
1	A	181	THR
1	A	183	VAL
1	A	184	VAL
1	A	190	LEU
1	A	194	ARG
1	A	196	VAL
1	B	2	SER
1	B	3	THR
1	B	8	LEU
1	B	10	GLU
1	B	13	SER
1	B	22	THR
1	B	31	ILE
1	B	32	ASN
1	B	38	VAL
1	B	41	LEU
1	B	52	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	61	ILE
1	B	68	ASP
1	B	75	ARG
1	B	76	ASP
1	B	79	THR
1	B	90	VAL
1	B	92	MET
1	B	95	VAL
1	B	96	LYS
1	B	99	LYS
1	B	109	ILE
1	B	118	GLU
1	B	120	ILE
1	B	121	ARG
1	B	124	GLU
1	B	125	LEU
1	B	130	ARG
1	B	134	VAL
1	B	145	ILE
1	B	148	GLU
1	B	152	MET
1	B	153	TYR
1	B	156	LYS
1	B	159	LYS
1	B	163	VAL
1	B	165	VAL
1	B	168	ILE
1	B	172	ILE
1	B	178	LEU
1	B	179	GLU
1	B	186	LEU
1	C	7	LYS
1	C	8	LEU
1	C	31	ILE
1	C	35	LEU
1	C	39	ARG
1	C	41	LEU
1	C	66	VAL
1	C	73	GLU
1	C	79	THR
1	C	83	VAL
1	C	87	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	95	VAL
1	C	113	GLU
1	C	118	GLU
1	C	120	ILE
1	C	121	ARG
1	C	124	GLU
1	C	125	LEU
1	C	126	LEU
1	C	129	VAL
1	C	134	VAL
1	C	154	ILE
1	C	157	LEU
1	C	159	LYS
1	C	163	VAL
1	C	164	LYS
1	C	165	VAL
1	C	181	THR
1	C	183	VAL
1	C	184	VAL
1	C	190	LEU
1	D	-2	MET
1	D	8	LEU
1	D	10	GLU
1	D	31	ILE
1	D	32	ASN
1	D	35	LEU
1	D	37	GLU
1	D	38	VAL
1	D	41	LEU
1	D	49	LYS
1	D	52	LEU
1	D	63	ASP
1	D	66	VAL
1	D	75	ARG
1	D	79	THR
1	D	83	VAL
1	D	84	SER
1	D	95	VAL
1	D	96	LYS
1	D	99	LYS
1	D	112	ILE
1	D	118	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	121	ARG
1	D	124	GLU
1	D	125	LEU
1	D	146	GLU
1	D	152	MET
1	D	164	LYS
1	D	165	VAL
1	D	168	ILE
1	D	177	ASP
1	D	178	LEU
1	D	186	LEU
1	D	196	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	77	HIS
1	A	85	HIS
1	A	106	HIS
1	A	142	ASN
1	B	43	GLN
1	B	77	HIS
1	B	85	HIS
1	B	106	HIS
1	B	142	ASN
1	C	85	HIS
1	C	106	HIS
1	D	77	HIS
1	D	85	HIS
1	D	106	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	IMD	A	202	-	5,5,5	0.55	0	5,5,5	0.62	0
3	IMD	A	204	-	5,5,5	0.87	0	5,5,5	0.31	0
3	IMD	B	202	-	5,5,5	0.63	0	5,5,5	0.64	0
3	IMD	C	202	-	5,5,5	0.73	0	5,5,5	0.32	0
3	IMD	D	202	-	5,5,5	0.98	0	5,5,5	0.89	0
3	IMD	A	203	-	5,5,5	0.75	0	5,5,5	0.30	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	IMD	A	202	-	-	-	0/1/1/1
3	IMD	A	204	-	-	-	0/1/1/1
3	IMD	B	202	-	-	-	0/1/1/1
3	IMD	C	202	-	-	-	0/1/1/1
3	IMD	D	202	-	-	-	0/1/1/1
3	IMD	A	203	-	-	-	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	202	IMD	2	0
3	A	204	IMD	7	0
3	B	202	IMD	7	0
3	C	202	IMD	3	0
3	D	202	IMD	1	0
3	A	203	IMD	2	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	197/212 (92%)	-0.61	2 (1%) 79 80	24, 46, 73, 109	0
1	B	195/212 (91%)	-0.23	3 (1%) 72 74	30, 54, 100, 127	0
1	C	197/212 (92%)	-0.41	1 (0%) 87 88	29, 50, 85, 114	0
1	D	199/212 (93%)	-0.34	1 (0%) 87 88	26, 51, 93, 141	0
All	All	788/848 (92%)	-0.40	7 (0%) 81 82	24, 50, 92, 141	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	74	ASN	6.8
1	D	112	ILE	3.3
1	C	112	ILE	2.4
1	B	111	PRO	2.3
1	A	196	VAL	2.2
1	B	145	ILE	2.1
1	A	0	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	IMD	B	202	5/5	0.74	0.16	66,70,74,77	0
3	IMD	D	202	5/5	0.78	0.20	52,55,58,63	0
3	IMD	C	202	5/5	0.80	0.16	63,64,69,71	0
3	IMD	A	203	5/5	0.82	0.16	55,56,60,61	0
3	IMD	A	202	5/5	0.82	0.16	82,84,85,87	0
3	IMD	A	204	5/5	0.90	0.13	52,55,59,61	0
2	ZN	C	201	1/1	1.00	0.01	36,36,36,36	0
2	ZN	D	201	1/1	1.00	0.01	40,40,40,40	0
2	ZN	A	201	1/1	1.00	0.02	39,39,39,39	0
2	ZN	B	201	1/1	1.00	0.02	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.