

Full wwPDB X-ray Structure Validation Report (i)

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PDB ID	:	3VJJ
Title	:	Crystal Structure Analysis of the P9-1
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		Τ.
Deposited on	:	2011-10-24
Resolution	:	3.00 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 (i) 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 (1)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
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)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\;DIFFRACTION$

The reported resolution of this entry is 3.00 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$		
R _{free}	130704	2092 (3.00-3.00)		
Clashscore	141614	2416 (3.00-3.00)		
Ramachandran outliers	138981	2333 (3.00-3.00)		
Sidechain outliers	138945	2336 (3.00-3.00)		
RSRZ outliers	127900	1990 (3.00-3.00)		

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 <=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	А	368	25%	38%	11%	•	25%		
1	В	368	25%	39%	10%		27%		



2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 4475 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.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Λ	277	Total	С	Ν	0	S	0	0	0
	1 A	211	2265	1459	372	427	7	0		
1	В	270	Total	С	Ν	0	S	0	0	0
	ГВ	210	2210	1428	364	411	7	0	0	0

• Molecule 1 is a protein called P9-1.

Chain	Residue	Modelled	Actual	Comment	Reference				
А	1	MET	-	expression tag	UNP Q913E4				
А	2	SER	-	expression tag	UNP Q913E4				
А	3	TYR	-	expression tag	UNP Q913E4				
А	4	TYR	-	expression tag	UNP Q913E4				
А	5	HIS	-	expression tag	UNP Q913E4				
А	6	HIS	-	expression tag	UNP Q913E4				
А	7	HIS	-	expression tag	UNP Q913E4				
А	8	HIS	-	expression tag	UNP Q913E4				
А	9	HIS	-	expression tag	UNP Q913E4				
А	10	HIS	-	expression tag	UNP Q913E4				
А	11	LEU	-	expression tag	UNP Q913E4				
А	12	GLY	-	expression tag	UNP Q913E4				
А	13	SER	-	expression tag	UNP Q913E4				
А	14	THR	-	expression tag	UNP Q913E4				
А	15	SER	-	expression tag	UNP Q913E4				
А	16	LEU	-	expression tag	UNP Q913E4				
А	17	TYR	-	expression tag	UNP Q913E4				
А	18	LYS	-	expression tag	UNP Q913E4				
А	19	LYS	-	expression tag	UNP Q913E4				
А	20	ALA	-	expression tag	UNP Q913E4				
А	21	GLY	-	expression tag	UNP Q913E4				
А	346	LYS	ARG	conflict	UNP Q913E4				
В	1	MET	-	expression tag	UNP Q913E4				
В	2	SER	-	expression tag	UNP Q913E4				
В	3	TYR	-	expression tag	UNP Q913E4				
	Continued on next page								

There are 44 discrepancies between the modelled and reference sequences:



Chain	Residue	Modelled	Actual	Comment	Reference
В	4	TYR	_	expression tag	UNP Q913E4
В	5	HIS	-	expression tag	UNP Q913E4
В	6	HIS	-	expression tag	UNP Q913E4
В	7	HIS	-	expression tag	UNP Q913E4
В	8	HIS	-	expression tag	UNP Q913E4
В	9	HIS	-	expression tag	UNP Q913E4
В	10	HIS	-	expression tag	UNP Q913E4
В	11	LEU	-	expression tag	UNP Q913E4
В	12	GLY	-	expression tag	UNP Q913E4
В	13	SER	-	expression tag	UNP Q913E4
В	14	THR	-	expression tag	UNP Q913E4
В	15	SER	-	expression tag	UNP Q913E4
В	16	LEU	-	expression tag	UNP Q913E4
В	17	TYR	-	expression tag	UNP Q913E4
В	18	LYS	-	expression tag	UNP Q913E4
В	19	LYS	-	expression tag	UNP Q913E4
В	20	ALA	-	expression tag	UNP Q913E4
В	21	GLY	-	expression tag	UNP Q913E4
В	346	LYS	ARG	conflict	UNP Q913E4



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: P9-1







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants	127.28Å 127.28Å 143.23Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Bosolution(A)	42.18 - 3.00	Depositor
Resolution (A)	42.18 - 3.00	EDS
% Data completeness	94.6 (42.18-3.00)	Depositor
(in resolution range)	93.6~(42.18-3.00)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$5.33 (at 3.01 \text{\AA})$	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
P. P.	0.243 , 0.293	Depositor
II, II free	0.239 , 0.285	DCC
R_{free} test set	974 reflections (4.25%)	wwPDB-VP
Wilson B-factor $(Å^2)$	94.8	Xtriage
Anisotropy	0.580	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.33, 89.3	EDS
L-test for $twinning^2$	$ L > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4475	wwPDB-VP
Average B, all atoms $(Å^2)$	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.08% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

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

Mol	Chain	Bo	nd lengths	Bond angles		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.70	1/2306~(0.0%)	0.87	7/3106~(0.2%)	
1	В	0.56	0/2247	0.84	3/3021~(0.1%)	
All	All	0.63	1/4553~(0.0%)	0.86	10/6127~(0.2%)	

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	А	0	2
1	В	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
1	А	41	LYS	C-N	18.06	1.75	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	261	ARG	NE-CZ-NH1	13.60	127.10	120.30
1	В	261	ARG	NE-CZ-NH2	12.38	126.49	120.30
1	А	261	ARG	NE-CZ-NH2	-12.34	114.13	120.30
1	В	261	ARG	NE-CZ-NH1	-11.62	114.49	120.30
1	А	41	LYS	C-N-CA	6.88	138.89	121.70
1	А	261	ARG	CD-NE-CZ	6.24	132.34	123.60
1	В	261	ARG	CD-NE-CZ	5.57	131.40	123.60
1	А	245	ARG	NE-CZ-NH1	-5.26	117.67	120.30
1	А	40	ILE	O-C-N	-5.21	114.36	122.70
1	А	196	ARG	NE-CZ-NH2	-5.03	117.78	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	182	GLU	Peptide
1	А	201	SER	Peptide
1	В	182	GLU	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	2265	0	2296	231	0
1	В	2210	0	2251	215	2
All	All	4475	0	4547	438	2

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 49.

All (438) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:41:LYS:C	1:A:63:THR:N	1.75	1.38
1:A:201:SER:H	1:A:202:LEU:HD22	1.16	1.05
1:A:97:SER:HB3	1:A:99:LEU:H	1.22	1.03
1:B:97:SER:HB3	1:B:99:LEU:H	1.19	1.02
1:A:353:VAL:O	1:A:356:VAL:HG12	1.62	0.98
1:A:329:GLY:HA2	1:A:332:GLY:H	1.31	0.93
1:A:233:ASN:ND2	1:A:235:TYR:H	1.68	0.90
1:B:218:LYS:HG2	1:B:251:ILE:HD13	1.54	0.90
1:A:201:SER:N	1:A:202:LEU:HD22	1.87	0.88
1:B:233:ASN:ND2	1:B:235:TYR:H	1.71	0.88
1:B:329:GLY:HA2	1:B:332:GLY:H	1.37	0.88
1:B:183:LYS:HG3	1:B:184:LYS:HE3	1.57	0.87
1:A:183:LYS:HG3	1:A:184:LYS:HE3	1.57	0.86
1:B:30:PHE:HD1	1:B:91:TYR:CD2	1.94	0.86
1:B:97:SER:HB3	1:B:99:LEU:N	1.89	0.85
1:A:30:PHE:HD1	1:A:91:TYR:CD2	1.95	0.84

	A L O	Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
1:A:96:TYR:HA	1:A:97:SER:OG	1.77	0.84
1:A:201:SER:CB	1:A:202:LEU:HA	2.08	0.84
1:A:97:SER:HB3	1:A:99:LEU:N	1.92	0.83
1:A:80:LEU:HB3	1:A:113:VAL:HG11	1.59	0.83
1:B:96:TYR:HA	1:B:97:SER:OG	1.80	0.82
1:A:292:LEU:HD12	1:A:292:LEU:H	1.45	0.81
1:A:201:SER:OG	1:A:202:LEU:HD13	1.79	0.81
1:A:218:LYS:HG2	1:A:251:ILE:HD13	1.61	0.80
1:B:292:LEU:HD12	1:B:292:LEU:H	1.46	0.79
1:A:233:ASN:HD22	1:A:235:TYR:H	1.28	0.79
1:B:170:LYS:NZ	1:B:175:LYS:NZ	2.31	0.78
1:A:311:ILE:O	1:A:312:ILE:HD12	1.84	0.78
1:A:310:LYS:HE3	1:A:328:ARG:NH2	1.99	0.78
1:B:310:LYS:HE3	1:B:328:ARG:NH2	1.99	0.77
1:B:233:ASN:HD22	1:B:235:TYR:H	1.29	0.77
1:A:236:ASP:OD1	1:A:286:ARG:NH2	2.16	0.77
1:A:176:VAL:HB	1:A:189:SER:HB2	1.67	0.76
1:A:33:TYR:OH	1:A:228:LEU:HA	1.85	0.76
1:A:70:ASP:HA	1:A:71:ASP:C	2.06	0.76
1:A:176:VAL:HA	1:A:189:SER:HA	1.67	0.76
1:B:170:LYS:HZ3	1:B:175:LYS:CE	1.98	0.75
1:B:79:TYR:HE1	1:B:89:THR:HB	1.51	0.75
1:B:261:ARG:O	1:B:262:VAL:HG22	1.85	0.75
1:B:353:VAL:HG12	1:B:356:VAL:CG1	2.18	0.74
1:B:70:ASP:HA	1:B:71:ASP:C	2.07	0.74
1:B:176:VAL:HA	1:B:189:SER:HA	1.70	0.74
1:B:33:TYR:OH	1:B:228:LEU:HA	1.88	0.73
1:B:41:LYS:O	1:B:63:THR:N	2.21	0.73
1:B:80:LEU:HB3	1:B:113:VAL:HG11	1.69	0.73
1:A:347:THR:OG1	1:A:350:VAL:HG12	1.89	0.73
1:A:79:TYR:HE1	1:A:89:THR:HB	1.52	0.73
1:B:178:THR:HG22	1:B:178:THR:O	1.88	0.72
1:A:79:TYR:CE1	1:A:89:THR:HB	2.25	0.72
1:B:353:VAL:HG12	1:B:356:VAL:HG11	1.71	0.71
1:A:178:THR:HG22	1:A:178:THR:O	1.91	0.71
1:A:74:PHE:HB2	1:A:342:TYR:HB3	1.72	0.71
1:A:221:ILE:HG21	1:A:271:PHE:HE1	1.54	0.71
1:A:170:LYS:NZ	1:A:175:LYS:NZ	2.39	0.71
1:B:217:PHE:HB3	1:B:251:ILE:HD11	1.72	0.70
1:A:164:GLU:HB2	1:A:181:GLU:HA	1.71	0.70
1:A:262:VAL:HB	1:A:263:PRO:CD	2.21	0.70

	lo do pagom	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:179:PRO:HD2	1:B:186:VAL:HG12	1.73	0.70
1:B:350:VAL:HG12	1:B:351:GLY:H	1.56	0.70
1:B:79:TYR:CE1	1:B:89:THR:HB	2.26	0.70
1:B:167:GLU:C	1:B:168:ASN:HD22	1.95	0.70
1:A:167:GLU:C	1:A:168:ASN:HD22	1.95	0.70
1:A:170:LYS:HZ3	1:A:175:LYS:CE	2.05	0.69
1:B:176:VAL:HB	1:B:189:SER:HB2	1.74	0.69
1:B:41:LYS:H	1:B:63:THR:N	1.90	0.69
1:B:236:ASP:OD1	1:B:286:ARG:NH2	2.23	0.69
1:B:170:LYS:HZ3	1:B:175:LYS:NZ	1.90	0.69
1:B:311:ILE:O	1:B:312:ILE:HD12	1.91	0.69
1:A:179:PRO:HD2	1:A:186:VAL:HG12	1.75	0.69
1:A:168:ASN:HD22	1:A:168:ASN:N	1.91	0.69
1:B:73:ILE:HG21	1:B:342:TYR:CE2	2.28	0.69
1:B:170:LYS:HZ3	1:B:175:LYS:HE2	1.57	0.69
1:A:71:ASP:C	1:A:73:ILE:H	1.94	0.69
1:A:201:SER:OG	1:A:202:LEU:HD22	1.93	0.69
1:B:168:ASN:HD22	1:B:168:ASN:N	1.91	0.68
1:B:177:VAL:HG23	1:B:179:PRO:HD3	1.75	0.68
1:B:74:PHE:HB2	1:B:342:TYR:HB3	1.76	0.68
1:A:217:PHE:HB3	1:A:251:ILE:HD11	1.75	0.68
1:B:71:ASP:C	1:B:73:ILE:H	1.95	0.68
1:B:254:TYR:CE2	1:B:258:ILE:HG21	2.28	0.68
1:A:258:ILE:HG12	1:A:258:ILE:O	1.92	0.68
1:A:73:ILE:HG21	1:A:342:TYR:CE2	2.28	0.67
1:B:109:ALA:HB1	1:B:213:VAL:HG12	1.76	0.67
1:B:41:LYS:N	1:B:63:THR:N	2.42	0.67
1:A:195:ASN:HB3	1:A:198:SER:HB3	1.75	0.67
1:A:239:LEU:O	1:A:281:PRO:HA	1.93	0.67
1:B:239:LEU:O	1:B:281:PRO:HA	1.95	0.67
1:A:170:LYS:HZ3	1:A:175:LYS:HE2	1.60	0.66
1:A:66:ILE:HG12	1:A:235:TYR:CD2	2.30	0.66
1:A:177:VAL:HG23	1:A:179:PRO:HD3	1.77	0.66
1:A:109:ALA:HB1	1:A:213:VAL:HG12	1.77	0.66
1:B:41:LYS:CA	1:B:63:THR:N	2.59	0.65
1:B:258:ILE:HG12	1:B:258:ILE:O	1.94	0.65
1:A:261:ARG:O	1:A:262:VAL:HG13	1.97	0.65
1:A:273:GLN:NE2	1:B:287:PHE:HE2	1.93	0.65
1:A:66:ILE:HG21	1:A:235:TYR:CG	2.31	0.65
1:B:366:LEU:O	1:B:367:ASP:HB2	1.97	0.65
1:A:308:LEU:HB3	1:A:309:PHE:CE2	2.32	0.65

	lo uo pugo	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:221:ILE:HG21	1:B:271:PHE:HE1	1.61	0.65
1:B:242:GLU:HA	1:B:278:ASP:HB2	1.79	0.65
1:A:173:THR:HG23	1:A:249:ILE:HG22	1.78	0.65
1:A:311:ILE:HD12	1:A:311:ILE:H	1.61	0.65
1:B:66:ILE:HG12	1:B:235:TYR:CD2	2.32	0.65
1:B:193:ILE:HG12	1:B:193:ILE:O	1.97	0.65
1:B:104:ARG:HG3	1:B:105:ASP:H	1.61	0.64
1:A:254:TYR:CE2	1:A:258:ILE:HG21	2.31	0.64
1:B:164:GLU:HB2	1:B:181:GLU:HA	1.80	0.64
1:B:114:ARG:HH11	1:B:117:PHE:HD2	1.46	0.63
1:A:29:THR:HG23	1:A:32:SER:HB3	1.81	0.63
1:B:190:LEU:N	1:B:190:LEU:HD23	2.13	0.63
1:A:193:ILE:HG12	1:A:193:ILE:O	1.99	0.62
1:B:247:LEU:HD23	1:B:247:LEU:H	1.63	0.62
1:A:233:ASN:HD22	1:A:235:TYR:N	1.97	0.62
1:A:175:LYS:HZ2	1:A:217:PHE:HZ	1.47	0.62
1:B:66:ILE:HD12	1:B:66:ILE:H	1.64	0.62
1:A:170:LYS:HZ3	1:A:175:LYS:NZ	1.97	0.62
1:A:174:PRO:HG2	1:A:176:VAL:CG1	2.30	0.62
1:A:77:LEU:HD11	1:A:170:LYS:HZ2	1.64	0.62
1:B:29:THR:HG23	1:B:32:SER:HB3	1.81	0.62
1:A:114:ARG:HH11	1:A:117:PHE:HD2	1.47	0.62
1:A:172:VAL:HG13	1:A:173:THR:HA	1.80	0.62
1:A:25:GLN:HE21	1:A:26:GLU:H	1.47	0.61
1:A:30:PHE:CD1	1:A:91:TYR:CD2	2.84	0.61
1:A:292:LEU:HB3	1:B:288:GLN:HE22	1.66	0.61
1:A:310:LYS:HE3	1:A:328:ARG:CZ	2.31	0.61
1:A:74:PHE:C	1:A:74:PHE:CD2	2.72	0.61
1:A:82:ASP:HB3	1:A:84:THR:HG23	1.83	0.61
1:B:127:LYS:HE2	1:B:182:GLU:OE1	2.01	0.61
1:A:190:LEU:HD23	1:A:190:LEU:N	2.15	0.61
1:A:247:LEU:O	1:A:251:ILE:HG13	2.00	0.61
1:B:172:VAL:HG13	1:B:173:THR:HA	1.81	0.61
1:A:74:PHE:HZ	1:A:166:GLU:HB3	1.66	0.60
1:A:292:LEU:HD12	1:A:292:LEU:N	2.16	0.60
1:B:310:LYS:HE3	1:B:328:ARG:CZ	2.30	0.60
1:B:41:LYS:C	1:B:63:THR:N	2.55	0.60
1:B:270:ILE:O	1:B:274:LEU:HG	2.00	0.60
1:B:174:PRO:HG2	1:B:176:VAL:CG1	2.32	0.60
1:B:341:ARG:C	1:B:342:TYR:HD1	2.05	0.60
1:A:270:ILE:O	1:A:274:LEU:HG	2.02	0.60

		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:364:LEU:HD12	1:A:364:LEU:O	2.01	0.60
1:B:311:ILE:HD12	1:B:311:ILE:H	1.67	0.60
1:B:66:ILE:HG21	1:B:235:TYR:CG	2.37	0.59
1:A:353:VAL:HG22	1:A:356:VAL:HG11	1.85	0.59
1:A:221:ILE:HG21	1:A:271:PHE:CE1	2.35	0.59
1:A:295:THR:HG22	1:A:299:ILE:CG2	2.32	0.59
1:B:74:PHE:HZ	1:B:166:GLU:HB3	1.68	0.59
1:B:325:VAL:HG22	1:B:337:LEU:HD23	1.84	0.59
1:B:30:PHE:CD1	1:B:91:TYR:CD2	2.83	0.59
1:B:82:ASP:HB3	1:B:84:THR:HG23	1.85	0.59
1:B:295:THR:HG22	1:B:299:ILE:CG2	2.33	0.59
1:B:308:LEU:HB3	1:B:309:PHE:CE2	2.38	0.59
1:A:201:SER:HB2	1:A:202:LEU:HA	1.83	0.59
1:A:192:PRO:HD2	1:A:210:GLU:OE2	2.01	0.59
1:A:104:ARG:HG3	1:A:105:ASP:H	1.67	0.58
1:A:288:GLN:HE22	1:B:292:LEU:HB3	1.67	0.58
1:B:173:THR:HG23	1:B:249:ILE:HG22	1.83	0.58
1:A:127:LYS:O	1:A:128:ASP:CB	2.51	0.58
1:B:252:ILE:HD11	1:B:327:PHE:CZ	2.38	0.58
1:A:74:PHE:CZ	1:A:166:GLU:HB3	2.38	0.58
1:B:94:PHE:HB2	1:B:96:TYR:CE1	2.39	0.58
1:A:308:LEU:HB3	1:A:309:PHE:CD2	2.39	0.57
1:B:292:LEU:HD12	1:B:292:LEU:N	2.17	0.57
1:B:97:SER:CB	1:B:99:LEU:H	2.06	0.57
1:B:233:ASN:HD22	1:B:235:TYR:N	2.00	0.57
1:B:104:ARG:HG3	1:B:105:ASP:N	2.19	0.57
1:A:287:PHE:HE2	1:B:273:GLN:NE2	2.02	0.57
1:B:191:LEU:HB3	1:B:194:LEU:HD12	1.85	0.57
1:B:171:PRO:HD2	1:B:176:VAL:O	2.04	0.57
1:B:339:ILE:HD12	1:B:339:ILE:H	1.70	0.57
1:B:341:ARG:C	1:B:342:TYR:CD1	2.78	0.57
1:A:180:LYS:HG3	1:A:185:THR:HA	1.87	0.57
1:B:74:PHE:C	1:B:74:PHE:CD2	2.77	0.57
1:A:242:GLU:HA	1:A:278:ASP:HB2	1.87	0.56
1:B:247:LEU:HD23	1:B:247:LEU:N	2.20	0.56
1:A:325:VAL:HG22	1:A:337:LEU:HD23	1.87	0.56
1:B:74:PHE:CZ	1:B:166:GLU:HB3	2.41	0.56
1:A:171:PRO:HD2	1:A:176:VAL:O	2.05	0.56
1:A:258:ILE:O	1:A:262:VAL:HG23	2.05	0.56
1:A:254:TYR:O	1:A:258:ILE:HG22	2.06	0.56
1:A:69:LEU:O	1:A:71:ASP:HB3	2.06	0.56

		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:348:ARG:O	1:A:349:ILE:HB	2.05	0.55
1:B:180:LYS:HG3	1:B:185:THR:HA	1.89	0.55
1:B:295:THR:HG22	1:B:299:ILE:HG22	1.87	0.55
1:B:122:GLU:HG3	1:B:123:ASN:N	2.21	0.55
1:A:68:LEU:O	1:A:70:ASP:N	2.40	0.55
1:A:261:ARG:O	1:A:262:VAL:HG22	2.06	0.55
1:A:262:VAL:O	1:A:263:PRO:C	2.43	0.55
1:B:77:LEU:HD11	1:B:170:LYS:HZ2	1.72	0.55
1:B:170:LYS:NZ	1:B:175:LYS:HZ3	2.05	0.55
1:B:258:ILE:O	1:B:262:VAL:HG23	2.07	0.54
1:A:122:GLU:HG3	1:A:123:ASN:N	2.22	0.54
1:A:191:LEU:HB3	1:A:194:LEU:HD12	1.89	0.54
1:A:339:ILE:HD12	1:A:339:ILE:H	1.72	0.54
1:B:339:ILE:HD12	1:B:339:ILE:N	2.23	0.54
1:A:201:SER:OG	1:A:202:LEU:HA	2.06	0.54
1:A:33:TYR:HB2	1:A:70:ASP:O	2.08	0.54
1:B:66:ILE:HD12	1:B:66:ILE:N	2.21	0.54
1:B:366:LEU:O	1:B:367:ASP:CB	2.54	0.54
1:B:350:VAL:HG12	1:B:351:GLY:N	2.22	0.54
1:B:68:LEU:O	1:B:70:ASP:N	2.41	0.54
1:B:33:TYR:HB2	1:B:70:ASP:O	2.08	0.54
1:B:178:THR:OG1	1:B:187:GLU:HG3	2.07	0.54
1:A:97:SER:CB	1:A:99:LEU:H	2.09	0.53
1:A:94:PHE:HB2	1:A:96:TYR:CE1	2.43	0.53
1:A:341:ARG:C	1:A:342:TYR:CD1	2.82	0.53
1:A:341:ARG:C	1:A:342:TYR:HD1	2.11	0.53
1:A:349:ILE:HG22	1:A:349:ILE:O	2.07	0.53
1:B:170:LYS:NZ	1:B:175:LYS:HZ1	2.06	0.53
1:B:327:PHE:O	1:B:334:GLU:HA	2.08	0.53
1:A:174:PRO:HB2	1:A:176:VAL:HG12	1.90	0.53
1:A:295:THR:HG22	1:A:299:ILE:HG22	1.88	0.53
1:B:247:LEU:O	1:B:251:ILE:HG13	2.08	0.53
1:B:288:GLN:OE1	1:B:288:GLN:HA	2.08	0.53
1:A:70:ASP:HA	1:A:71:ASP:HB3	1.91	0.53
1:A:71:ASP:O	1:A:73:ILE:N	2.42	0.53
1:A:252:ILE:HD11	1:A:327:PHE:CZ	2.43	0.53
1:B:178:THR:O	1:B:178:THR:CG2	2.55	0.53
1:A:126:LEU:O	1:A:127:LYS:HB2	2.09	0.53
1:B:341:ARG:O	1:B:342:TYR:HD1	1.92	0.53
1:B:109:ALA:O	1:B:113:VAL:HG12	2.09	0.52
1:B:71:ASP:O	1:B:73:ILE:N	2.42	0.52

		Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
1:A:247:LEU:HD23	1:A:247:LEU:H	1.73	0.52
1:A:96:TYR:HA	1:A:97:SER:CB	2.35	0.52
1:A:284:LEU:HD12	1:A:284:LEU:O	2.09	0.52
1:B:192:PRO:HD2	1:B:210:GLU:OE2	2.09	0.52
1:B:308:LEU:HB3	1:B:309:PHE:CD2	2.45	0.52
1:A:178:THR:OG1	1:A:187:GLU:HG3	2.10	0.52
1:A:288:GLN:OE1	1:A:288:GLN:HA	2.10	0.52
1:B:41:LYS:O	1:B:42:ASN:HB3	2.09	0.52
1:B:184:LYS:H	1:B:184:LYS:HD2	1.74	0.52
1:A:91:TYR:N	1:A:91:TYR:CD1	2.75	0.52
1:A:339:ILE:HD12	1:A:339:ILE:N	2.24	0.51
1:A:341:ARG:O	1:A:342:TYR:HD1	1.93	0.51
1:A:33:TYR:OH	1:A:228:LEU:HD12	2.11	0.51
1:A:104:ARG:HG3	1:A:105:ASP:N	2.25	0.51
1:B:175:LYS:HZ2	1:B:217:PHE:HZ	1.58	0.51
1:B:353:VAL:HG12	1:B:356:VAL:HG12	1.92	0.51
1:B:353:VAL:O	1:B:356:VAL:HG12	2.10	0.51
1:A:170:LYS:NZ	1:A:175:LYS:HZ1	2.07	0.51
1:B:212:ALA:O	1:B:216:VAL:HG13	2.11	0.51
1:A:71:ASP:C	1:A:73:ILE:N	2.63	0.51
1:B:221:ILE:HG21	1:B:271:PHE:CE1	2.44	0.51
1:B:254:TYR:O	1:B:258:ILE:HG22	2.11	0.51
1:B:114:ARG:NH1	1:B:117:PHE:HD2	2.09	0.51
1:B:217:PHE:CB	1:B:251:ILE:HD11	2.40	0.50
1:B:284:LEU:HD12	1:B:284:LEU:O	2.11	0.50
1:B:289:LEU:HD12	1:B:292:LEU:HD21	1.93	0.50
1:A:247:LEU:HD23	1:A:247:LEU:N	2.27	0.50
1:A:174:PRO:HG2	1:A:176:VAL:HG12	1.94	0.50
1:B:29:THR:HB	1:B:94:PHE:CD1	2.47	0.50
1:B:28:ARG:HH22	1:B:70:ASP:CB	2.25	0.50
1:A:178:THR:O	1:A:178:THR:CG2	2.59	0.49
1:B:35:ILE:HD13	1:B:239:LEU:CD1	2.42	0.49
1:A:80:LEU:O	1:A:113:VAL:HG13	2.12	0.49
1:B:96:TYR:CA	1:B:97:SER:OG	2.58	0.49
1:A:41:LYS:N	1:A:63:THR:N	2.61	0.49
1:A:262:VAL:HB	1:A:263:PRO:HD2	1.93	0.49
1:B:347:THR:O	1:B:348:ARG:C	2.50	0.49
1:A:101:ASN:N	1:A:101:ASN:HD22	2.10	0.49
1:A:361:PHE:O	1:A:364:LEU:HG	2.12	0.49
1:B:174:PRO:HG2	1:B:176:VAL:HG12	1.95	0.49
1:B:91:TYR:N	1:B:91:TYR:CD1	2.79	0.49

		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:96:TYR:HA	1:B:97:SER:CB	2.37	0.49
1:B:169:SER:HA	1:B:340:LYS:O	2.13	0.49
1:B:97:SER:HB3	1:B:99:LEU:CA	2.42	0.49
1:B:282:PHE:O	1:B:285:LEU:HB3	2.13	0.49
1:A:117:PHE:O	1:A:121:THR:HG23	2.13	0.49
1:A:201:SER:OG	1:A:202:LEU:CD1	2.54	0.49
1:A:79:TYR:HE1	1:A:89:THR:CB	2.23	0.49
1:A:28:ARG:HG3	1:A:32:SER:OG	2.13	0.48
1:A:184:LYS:H	1:A:184:LYS:HD2	1.78	0.48
1:A:221:ILE:CG2	1:A:271:PHE:HE1	2.25	0.48
1:B:69:LEU:O	1:B:71:ASP:HB3	2.13	0.48
1:B:79:TYR:HE1	1:B:89:THR:CB	2.24	0.48
1:B:33:TYR:CB	1:B:70:ASP:O	2.61	0.48
1:B:274:LEU:N	1:B:274:LEU:HD23	2.27	0.48
1:A:68:LEU:O	1:A:68:LEU:HG	2.13	0.48
1:B:303:ILE:O	1:B:306:GLU:O	2.30	0.48
1:A:100:PRO:C	1:A:101:ASN:ND2	2.67	0.48
1:A:289:LEU:HD12	1:A:292:LEU:HD21	1.96	0.48
1:A:173:THR:HG23	1:A:249:ILE:CG2	2.43	0.48
1:A:212:ALA:O	1:A:216:VAL:HG13	2.13	0.48
1:B:248:LEU:HD22	1:B:274:LEU:HB2	1.96	0.48
1:A:33:TYR:CB	1:A:70:ASP:O	2.61	0.48
1:A:327:PHE:O	1:A:334:GLU:HA	2.13	0.48
1:B:30:PHE:CD1	1:B:91:TYR:CE2	3.02	0.48
1:B:41:LYS:O	1:B:42:ASN:CB	2.61	0.48
1:A:70:ASP:CA	1:A:71:ASP:C	2.81	0.48
1:A:201:SER:H	1:A:202:LEU:CD2	2.04	0.48
1:A:303:ILE:O	1:A:306:GLU:O	2.32	0.48
1:A:169:SER:HA	1:A:340:LYS:O	2.14	0.48
1:A:28:ARG:HH22	1:A:70:ASP:CB	2.26	0.47
1:A:30:PHE:CD1	1:A:91:TYR:CE2	3.03	0.47
1:A:82:ASP:CB	1:A:84:THR:HG23	2.44	0.47
1:A:101:ASN:N	1:A:101:ASN:ND2	2.62	0.47
1:B:28:ARG:HH22	1:B:70:ASP:CG	2.17	0.47
1:A:29:THR:HB	1:A:94:PHE:CD1	2.49	0.47
1:B:66:ILE:N	1:B:66:ILE:CD1	2.77	0.47
1:B:364:LEU:HD12	1:B:364:LEU:H	1.78	0.47
1:A:309:PHE:HA	1:A:326:LEU:O	2.14	0.47
1:B:68:LEU:O	1:B:68:LEU:HG	2.14	0.47
1:A:29:THR:CG2	1:A:32:SER:HB3	2.44	0.47
1:A:97:SER:HB3	1:A:99:LEU:CA	2.45	0.47

	lo ao pagom	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:109:ALA:O	1:A:113:VAL:HG12	2.15	0.47
1:A:168:ASN:N	1:A:168:ASN:ND2	2.61	0.47
1:B:295:THR:HA	1:B:330:ILE:HD11	1.97	0.47
1:B:342:TYR:CD1	1:B:342:TYR:N	2.78	0.47
1:A:25:GLN:HE21	1:A:26:GLU:N	2.12	0.47
1:B:261:ARG:O	1:B:262:VAL:HG13	2.15	0.47
1:B:309:PHE:HA	1:B:326:LEU:O	2.15	0.47
1:A:66:ILE:N	1:A:66:ILE:HD12	2.30	0.47
1:A:112:TYR:CG	1:A:209:GLY:HA3	2.50	0.47
1:A:125:ASP:OD2	1:A:127:LYS:HG3	2.15	0.47
1:B:28:ARG:HG3	1:B:32:SER:OG	2.15	0.47
1:A:70:ASP:CA	1:A:71:ASP:HB3	2.45	0.46
1:A:190:LEU:O	1:A:191:LEU:C	2.54	0.46
1:B:70:ASP:HA	1:B:71:ASP:HB3	1.96	0.46
1:B:37:GLU:HG2	1:B:64:LYS:HD3	1.96	0.46
1:B:80:LEU:O	1:B:113:VAL:HG13	2.15	0.46
1:A:65:LYS:C	1:A:66:ILE:HG13	2.35	0.46
1:A:69:LEU:C	1:A:71:ASP:HB3	2.35	0.46
1:B:257:LEU:HD23	1:B:257:LEU:HA	1.65	0.46
1:A:114:ARG:NH1	1:A:117:PHE:HD2	2.12	0.46
1:A:190:LEU:N	1:A:190:LEU:CD2	2.77	0.46
1:A:282:PHE:O	1:A:285:LEU:HB3	2.15	0.46
1:B:73:ILE:CG2	1:B:342:TYR:CD2	2.99	0.46
1:A:257:LEU:HD23	1:A:257:LEU:HA	1.67	0.45
1:B:29:THR:CG2	1:B:32:SER:HB3	2.44	0.45
1:A:289:LEU:HD12	1:A:289:LEU:HA	1.63	0.45
1:A:28:ARG:HH22	1:A:70:ASP:CG	2.20	0.45
1:A:248:LEU:HD22	1:A:274:LEU:HB2	1.99	0.45
1:A:70:ASP:HB3	1:A:72:GLY:HA2	1.98	0.45
1:A:295:THR:HA	1:A:330:ILE:HD11	1.99	0.45
1:A:353:VAL:O	1:A:356:VAL:CG1	2.49	0.45
1:B:290:SER:HA	1:B:291:SER:HA	1.59	0.45
1:A:201:SER:N	1:A:202:LEU:CD2	2.70	0.45
1:B:190:LEU:N	1:B:190:LEU:CD2	2.76	0.45
1:A:201:SER:OG	1:A:202:LEU:CD2	2.64	0.45
1:B:100:PRO:C	1:B:101:ASN:ND2	2.70	0.45
1:B:101:ASN:N	1:B:101:ASN:HD22	2.13	0.45
1:B:170:LYS:HZ1	1:B:175:LYS:NZ	2.11	0.45
1:B:82:ASP:CB	1:B:84:THR:HG23	2.46	0.45
1:A:28:ARG:HG2	1:A:29:THR:N	2.31	0.45
1:A:79:TYR:CE1	1:A:89:THR:CB	2.99	0.45

	loue page	Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlap (Å)
1:A:288:GLN:NE2	1:B:292:LEU:HB3	2.32	0.45
1:B:28:ARG:HG2	1:B:29:THR:N	2.31	0.45
1:A:235:TYR:O	1:A:236:ASP:C	2.55	0.45
1:A:35:ILE:HD13	1:A:239:LEU:CD1	2.47	0.44
1:A:73:ILE:CG2	1:A:342:TYR:CD2	3.00	0.44
1:B:71:ASP:C	1:B:73:ILE:N	2.64	0.44
1:B:101:ASN:ND2	1:B:101:ASN:N	2.66	0.44
1:B:214:VAL:HG21	1:B:254:TYR:CG	2.52	0.44
1:A:73:ILE:HG22	1:A:342:TYR:CD2	2.51	0.44
1:A:342:TYR:CD1	1:A:342:TYR:N	2.83	0.44
1:B:229:GLY:HA2	1:B:286:ARG:HD3	1.99	0.44
1:A:33:TYR:CA	1:A:70:ASP:O	2.66	0.44
1:B:70:ASP:HB3	1:B:72:GLY:HA2	1.99	0.44
1:A:295:THR:HG22	1:A:299:ILE:HG21	2.00	0.44
1:A:306:GLU:HB3	1:A:328:ARG:HH12	1.82	0.44
1:B:33:TYR:OH	1:B:228:LEU:HD12	2.17	0.44
1:A:81:ILE:O	1:A:114:ARG:HD3	2.18	0.44
1:A:177:VAL:CG2	1:A:179:PRO:HD3	2.44	0.44
1:A:229:GLY:HA2	1:A:286:ARG:HD3	2.00	0.44
1:B:164:GLU:HB3	1:B:181:GLU:OE2	2.17	0.44
1:B:168:ASN:N	1:B:168:ASN:ND2	2.62	0.44
1:B:289:LEU:HD12	1:B:289:LEU:HA	1.70	0.44
1:B:73:ILE:HG22	1:B:342:TYR:CD2	2.52	0.44
1:B:303:ILE:HA	1:B:303:ILE:HD12	1.71	0.44
1:A:283:ASP:OD1	1:A:283:ASP:N	2.51	0.43
1:B:176:VAL:O	1:B:176:VAL:HG22	2.16	0.43
1:A:356:VAL:O	1:A:356:VAL:HG13	2.18	0.43
1:B:117:PHE:O	1:B:121:THR:HG23	2.17	0.43
1:A:102:LEU:HD12	1:A:219:MET:SD	2.59	0.43
1:B:170:LYS:CE	1:B:175:LYS:HZ3	2.32	0.43
1:B:174:PRO:HB2	1:B:176:VAL:HG12	2.00	0.43
1:A:96:TYR:CA	1:A:97:SER:OG	2.57	0.43
1:B:28:ARG:NH1	1:B:32:SER:O	2.52	0.43
1:B:306:GLU:HB3	1:B:328:ARG:HH12	1.82	0.43
1:B:35:ILE:HD13	1:B:239:LEU:HD11	2.01	0.43
1:B:170:LYS:NZ	1:B:175:LYS:CE	2.74	0.43
1:B:296:PRO:O	1:B:300:ARG:HG3	2.18	0.43
1:A:66:ILE:HD12	1:A:66:ILE:H	1.83	0.43
1:A:273:GLN:HE21	1:B:287:PHE:HE2	1.63	0.43
1:B:41:LYS:O	1:B:42:ASN:ND2	2.52	0.43
1:B:112:TYR:CG	1:B:209:GLY:HA3	2.54	0.43

	1	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:195:ASN:HB3	1:B:196:ARG:H	1.64	0.43
1:B:79:TYR:CE1	1:B:89:THR:CB	3.00	0.42
1:B:224:PHE:CD2	1:B:225:LEU:N	2.87	0.42
1:A:41:LYS:H	1:A:63:THR:N	2.16	0.42
1:A:348:ARG:O	1:A:349:ILE:CB	2.67	0.42
1:A:290:SER:HA	1:A:291:SER:HA	1.57	0.42
1:B:97:SER:OG	1:B:99:LEU:HB2	2.19	0.42
1:B:261:ARG:C	1:B:262:VAL:HG22	2.39	0.42
1:A:28:ARG:NH1	1:A:32:SER:O	2.53	0.42
1:A:33:TYR:HA	1:A:70:ASP:O	2.20	0.42
1:B:70:ASP:CA	1:B:71:ASP:C	2.83	0.42
1:B:228:LEU:HD11	1:B:282:PHE:HB3	2.02	0.42
1:A:292:LEU:HB3	1:B:288:GLN:NE2	2.34	0.42
1:B:70:ASP:CA	1:B:71:ASP:HB3	2.48	0.42
1:B:174:PRO:O	1:B:175:LYS:HB3	2.19	0.42
1:A:164:GLU:HB3	1:A:181:GLU:OE2	2.19	0.42
1:A:273:GLN:NE2	1:B:287:PHE:CE2	2.82	0.42
1:A:280:TYR:HA	1:A:281:PRO:HD3	1.86	0.42
1:A:28:ARG:O	1:A:91:TYR:HA	2.20	0.42
1:A:37:GLU:HG2	1:A:64:LYS:HD3	2.01	0.42
1:A:173:THR:N	1:A:174:PRO:HD2	2.35	0.42
1:B:190:LEU:O	1:B:191:LEU:C	2.58	0.42
1:A:303:ILE:HD12	1:A:303:ILE:HA	1.72	0.42
1:B:177:VAL:CG2	1:B:179:PRO:HD3	2.49	0.42
1:A:173:THR:O	1:A:175:LYS:N	2.53	0.41
1:A:232:PRO:HA	1:A:286:ARG:NH1	2.35	0.41
1:A:66:ILE:HG21	1:A:235:TYR:CD2	2.56	0.41
1:A:224:PHE:CE2	1:A:225:LEU:HD23	2.55	0.41
1:B:173:THR:N	1:B:174:PRO:HD2	2.35	0.41
1:B:224:PHE:CE2	1:B:225:LEU:HD23	2.55	0.41
1:A:176:VAL:O	1:A:176:VAL:HG22	2.19	0.41
1:B:289:LEU:CD1	1:B:292:LEU:HD21	2.50	0.41
1:B:364:LEU:HD12	1:B:364:LEU:N	2.34	0.41
1:A:170:LYS:NZ	1:A:175:LYS:CE	2.80	0.41
1:A:284:LEU:O	1:A:288:GLN:HB2	2.20	0.41
1:B:69:LEU:C	1:B:71:ASP:HB3	2.41	0.41
1:B:295:THR:HG22	1:B:299:ILE:HG21	2.03	0.41
1:A:169:SER:C	1:A:170:LYS:HG2	2.39	0.41
1:A:232:PRO:HA	1:A:286:ARG:HH12	1.86	0.41
1:A:262:VAL:O	1:A:263:PRO:O	2.39	0.41
1:B:28:ARG:O	1:B:91:TYR:HA	2.21	0.41

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:LEU:O	1:A:68:LEU:CG	2.67	0.40
1:A:73:ILE:CG2	1:A:342:TYR:CE2	3.02	0.40
1:B:240:ASN:ND2	1:B:242:GLU:HG3	2.37	0.40
1:A:199:GLU:O	1:A:200:GLU:C	2.59	0.40
1:A:214:VAL:HG21	1:A:254:TYR:CG	2.56	0.40
1:A:366:LEU:O	1:A:367:ASP:HB2	2.21	0.40
1:B:38:LEU:HB2	1:B:67:PRO:HD3	2.03	0.40
1:A:217:PHE:CB	1:A:251:ILE:HD11	2.49	0.40
1:A:228:LEU:O	1:A:286:ARG:NH1	2.54	0.40
1:B:110:SER:HA	1:B:113:VAL:HG12	2.03	0.40
1:B:113:VAL:O	1:B:116:ASN:HB3	2.20	0.40
1:B:179:PRO:HD2	1:B:186:VAL:CG1	2.48	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:SER:OG	$1:B:215:ASN:OD1[8_555]$	2.01	0.19
1:B:105:ASP:OD1	$1:B:208:GLU:OE2[8_555]$	2.16	0.04

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	Perc	entiles
1	А	271/368~(74%)	243~(90%)	23~(8%)	5(2%)	8	37
1	В	258/368~(70%)	238~(92%)	18 (7%)	2(1%)	19	57
All	All	529/736~(72%)	481 (91%)	41 (8%)	7 (1%)	12	45

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	262	VAL
1	А	349	ILE
1	В	96	TYR
1	А	69	LEU
1	А	96	TYR
1	А	182	GLU
1	В	69	LEU

5.3.2 Protein sidechains (i)

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	Pe	erce	entiles
1	А	258/342~(75%)	204 (79%)	54 (21%)		1	5
1	В	251/342 (73%)	201 (80%)	50 (20%)		1	7
All	All	509/684~(74%)	405 (80%)	104 (20%)		1	6

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	36	GLU
1	А	39	THR
1	А	63	THR
1	А	66	ILE
1	А	74	PHE
1	А	75	GLU
1	А	85	ASN
1	А	90	CYS
1	А	101	ASN
1	А	104	ARG
1	А	126	LEU
1	А	167	GLU
1	А	168	ASN
1	А	172	VAL
1	А	173	THR
1	А	176	VAL
1	А	178	THR

Mol	Chain	Res	Type
1	А	180	LYS
1	А	183	LYS
1	А	184	LYS
1	А	186	VAL
1	А	190	LEU
1	А	193	ILE
1	А	198	SER
1	А	201	SER
1	А	202	LEU
1	А	205	GLU
1	А	206	ILE
1	А	207	LEU
1	А	221	ILE
1	А	224	PHE
1	А	233	ASN
1	А	234	SER
1	А	237	ARG
1	А	247	LEU
1	А	257	LEU
1	А	258	ILE
1	А	262	VAL
1	А	265	LYS
1	А	280	TYR
1	А	290	SER
1	А	292	LEU
1	А	294	SER
1	А	295	THR
1	А	298	LEU
1	А	299	ILE
1	А	303	ILE
1	A	311	ILE
1	А	325	VAL
1	A	335	SER
1	A	338	ASN
1	А	346	LYS
1	A	347	THR
1	Α	348	ARG
1	В	39	THR
1	В	42	ASN
1	В	63	THR
1	В	66	ILE
1	В	75	GLU

Mol	Chain	Res	Type
1	В	85	ASN
1	В	90	CYS
1	В	101	ASN
1	В	104	ARG
1	В	128	ASP
1	В	167	GLU
1	В	168	ASN
1	В	172	VAL
1	В	173	THR
1	В	176	VAL
1	В	178	THR
1	В	180	LYS
1	В	183	LYS
1	В	184	LYS
1	В	186	VAL
1	В	190	LEU
1	В	193	ILE
1	В	202	LEU
1	В	205	GLU
1	В	206	ILE
1	В	221	ILE
1	В	224	PHE
1	В	233	ASN
1	В	234	SER
1	В	237	ARG
1	В	247	LEU
1	В	258	ILE
1	В	260	THR
1	В	265	LYS
1	В	280	TYR
1	В	290	SER
1	В	292	LEU
1	В	294	SER
1	В	295	THR
1	В	298	LEU
1	В	299	ILE
1	В	303	ILE
1	В	311	ILE
1	В	325	VAL
1	В	335	SER
1	В	338	ASN
1	В	346	LYS

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Mol	Chain	Res	Type
1	В	353	VAL
1	В	355	CYS
1	В	364	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	25	GLN
1	А	101	ASN
1	А	168	ASN
1	А	233	ASN
1	В	101	ASN
1	В	168	ASN
1	В	233	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides 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 (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, 95^{th} 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	А	277/368~(75%)	0.13	13 (4%) 31 11	65, 113, 177, 227	0
1	В	270/368~(73%)	0.21	17 (6%) 20 6	61, 109, 167, 223	0
All	All	547/736~(74%)	0.17	30 (5%) 25 9	61, 111, 174, 227	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	349	ILE	5.0
1	В	345	PHE	5.0
1	В	123	ASN	4.6
1	В	265	LYS	4.6
1	В	202	LEU	4.5
1	В	128	ASP	4.3
1	В	63	THR	4.2
1	В	348	ARG	3.5
1	А	38	LEU	3.4
1	В	347	THR	3.1
1	А	125	ASP	3.0
1	В	125	ASP	3.0
1	В	313	THR	2.8
1	А	345	PHE	2.8
1	А	88	LYS	2.8
1	В	124	LEU	2.7
1	А	264	ASN	2.7
1	А	311	ILE	2.7
1	А	40	ILE	2.5
1	А	63	THR	2.4
1	А	348	ARG	2.4
1	В	301	GLU	2.3
1	A	367	ASP	2.3
1	В	196	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	В	184	LYS	2.2
1	А	324	THR	2.1
1	А	184	LYS	2.1
1	В	307	GLY	2.1
1	А	349	ILE	2.0
1	В	312	ILE	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 monosaccharides 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.

