



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 21, 2024 – 01:21 AM EDT

PDB ID : 1R5J
Title : Crystal Structure of a Phosphotransacetylase from *Streptococcus pyogenes*
Authors : Xu, Q.S.; Shin, D.H.; Pufan, R.; Yokota, H.; Kim, R.; Kim, S.H.; Berkeley Structural Genomics Center (BSGC)
Deposited on : 2003-10-10
Resolution : 2.70 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

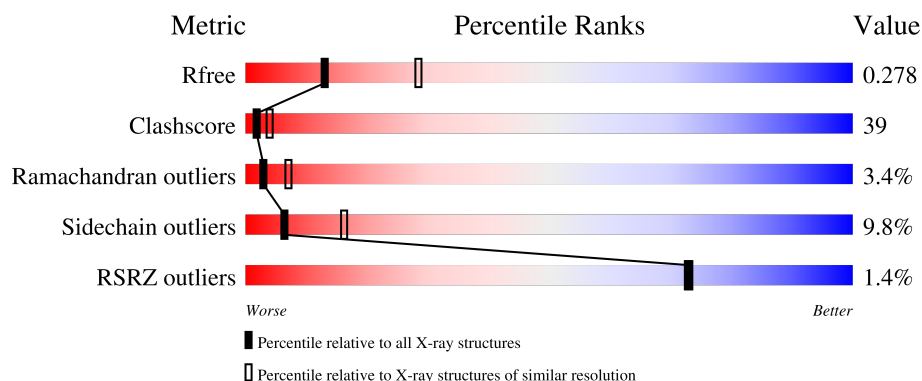
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	164625	3333 (2.70-2.70)
Clashscore	180529	3684 (2.70-2.70)
Ramachandran outliers	177936	3633 (2.70-2.70)
Sidechain outliers	177891	3633 (2.70-2.70)
RSRZ outliers	164620	3333 (2.70-2.70)

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	337	
1	B	337	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5034 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 putative phosphotransacetylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	Se	0	0	0
			2502	1582	424	485	1	10			
1	B	329	Total	C	N	O	S	Se	0	0	0
			2502	1582	424	485	1	10			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	cloning artifact	UNP Q99ZQ5
A	-4	GLY	-	cloning artifact	UNP Q99ZQ5
A	-3	GLY	-	cloning artifact	UNP Q99ZQ5
A	-2	GLY	-	cloning artifact	UNP Q99ZQ5
A	-1	GLY	-	cloning artifact	UNP Q99ZQ5
A	0	GLY	-	cloning artifact	UNP Q99ZQ5
A	1	MSE	MET	modified residue	UNP Q99ZQ5
A	19	MSE	MET	modified residue	UNP Q99ZQ5
A	83	MSE	MET	modified residue	UNP Q99ZQ5
A	103	MSE	MET	modified residue	UNP Q99ZQ5
A	113	MSE	MET	modified residue	UNP Q99ZQ5
A	117	MSE	MET	modified residue	UNP Q99ZQ5
A	123	MSE	MET	modified residue	UNP Q99ZQ5
A	157	MSE	MET	modified residue	UNP Q99ZQ5
A	204	MSE	MET	modified residue	UNP Q99ZQ5
A	288	MSE	MET	modified residue	UNP Q99ZQ5
B	-5	GLY	-	cloning artifact	UNP Q99ZQ5
B	-4	GLY	-	cloning artifact	UNP Q99ZQ5
B	-3	GLY	-	cloning artifact	UNP Q99ZQ5
B	-2	GLY	-	cloning artifact	UNP Q99ZQ5
B	-1	GLY	-	cloning artifact	UNP Q99ZQ5
B	0	GLY	-	cloning artifact	UNP Q99ZQ5
B	1	MSE	MET	modified residue	UNP Q99ZQ5
B	19	MSE	MET	modified residue	UNP Q99ZQ5
B	83	MSE	MET	modified residue	UNP Q99ZQ5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	103	MSE	MET	modified residue	UNP Q99ZQ5
B	113	MSE	MET	modified residue	UNP Q99ZQ5
B	117	MSE	MET	modified residue	UNP Q99ZQ5
B	123	MSE	MET	modified residue	UNP Q99ZQ5
B	157	MSE	MET	modified residue	UNP Q99ZQ5
B	204	MSE	MET	modified residue	UNP Q99ZQ5
B	288	MSE	MET	modified residue	UNP Q99ZQ5

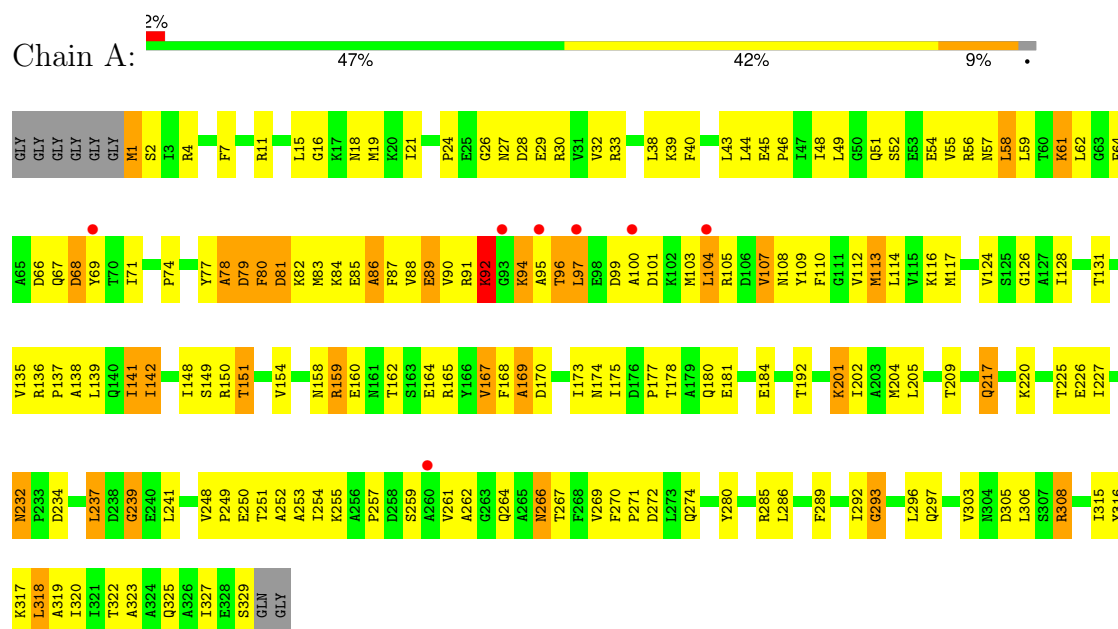
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	12	Total O 12 12	0	0
2	B	18	Total O 18 18	0	0

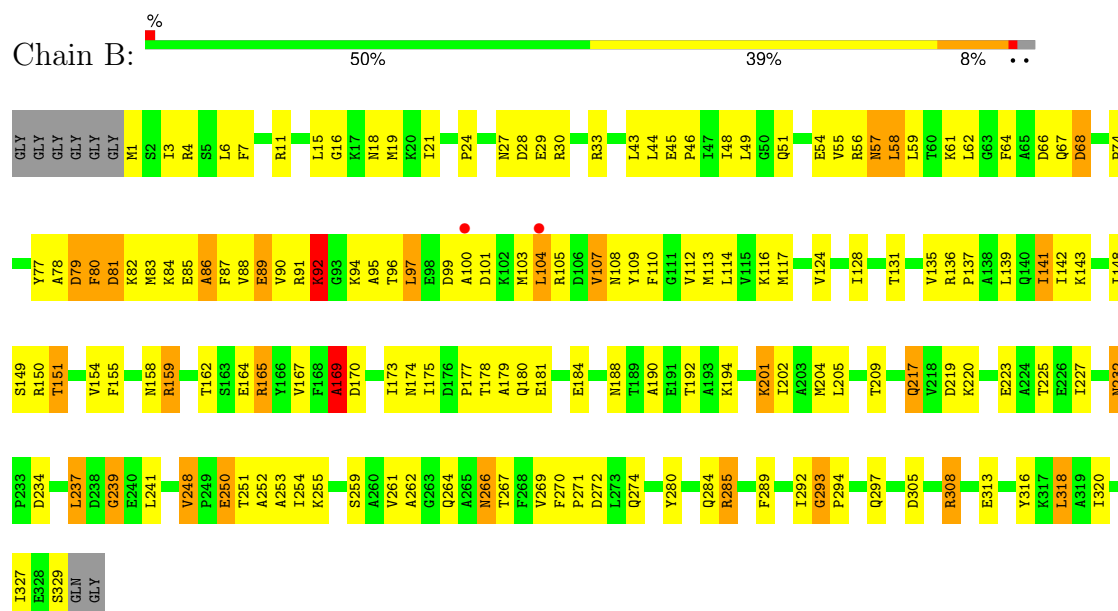
3 Residue-property plots

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: putative phosphotransacetylase



- Molecule 1: putative phosphotransacetylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 3 2	Depositor
Cell constants a, b, c, α , β , γ	173.72Å 173.72Å 173.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.93 – 2.70 19.93 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.93-2.70) 99.6 (19.93-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.71Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.229 , 0.281 0.227 , 0.278	Depositor DCC
R_{free} test set	2537 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å ²)	67.7	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 54.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5034	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.28% 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	0.62	5/2529 (0.2%)	0.78	4/3400 (0.1%)
1	B	0.63	3/2529 (0.1%)	0.79	3/3400 (0.1%)
All	All	0.63	8/5058 (0.2%)	0.78	7/6800 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	92	LYS	CE-NZ	7.80	1.68	1.49
1	A	96	THR	CB-CG2	6.92	1.75	1.52
1	A	94	LYS	CB-CG	6.36	1.69	1.52
1	B	92	LYS	CE-NZ	6.35	1.65	1.49
1	B	94	LYS	CD-CE	5.98	1.66	1.51
1	B	94	LYS	CE-NZ	5.52	1.62	1.49
1	A	92	LYS	CB-CG	5.42	1.67	1.52
1	A	94	LYS	CD-CE	5.35	1.64	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	ALA	N-CA-C	5.92	126.98	111.00
1	A	239	GLY	N-CA-C	5.75	127.49	113.10
1	B	169	ALA	N-CA-C	5.65	126.25	111.00
1	A	201	LYS	N-CA-C	-5.36	96.54	111.00
1	B	201	LYS	N-CA-C	-5.33	96.62	111.00
1	A	142	ILE	N-CA-C	-5.28	96.74	111.00
1	B	239	GLY	N-CA-C	5.16	126.00	113.10

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	2502	0	2542	216	0
1	B	2502	0	2542	180	0
2	A	12	0	0	1	0
2	B	18	0	0	1	0
All	All	5034	0	5084	393	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 39.

All (393) 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:96:THR:CB	1:A:96:THR:CG2	1.75	1.63
1:A:92:LYS:CE	1:A:92:LYS:NZ	1.68	1.56
1:A:169:ALA:CB	1:A:269:VAL:HA	1.71	1.19
1:B:169:ALA:CB	1:B:269:VAL:HA	1.74	1.18
1:A:92:LYS:HD2	1:A:92:LYS:N	1.58	1.15
1:B:92:LYS:HD2	1:B:92:LYS:N	1.66	1.10
1:A:169:ALA:HB3	1:A:270:PHE:H	1.20	1.07
1:B:82:LYS:HE3	1:B:117:MSE:HE1	1.37	1.06
1:A:83:MSE:HE3	1:A:104:LEU:HB3	1.37	1.03
1:A:178:THR:HG22	1:A:180:GLN:H	1.21	1.02
1:B:169:ALA:HB2	1:B:269:VAL:CA	1.91	1.00
1:B:178:THR:HG22	1:B:180:GLN:H	1.27	1.00
1:A:169:ALA:HB2	1:A:269:VAL:HA	0.98	0.97
1:A:82:LYS:HE3	1:A:117:MSE:HE1	1.45	0.96
1:A:169:ALA:HB2	1:A:269:VAL:CA	1.95	0.95
1:B:169:ALA:HB3	1:B:270:PHE:H	1.30	0.95
1:A:83:MSE:CE	1:A:104:LEU:HB3	1.95	0.95
1:B:169:ALA:HB2	1:B:269:VAL:HA	0.94	0.94
1:A:28:ASP:HB2	1:A:128:ILE:HD11	1.51	0.92
1:A:92:LYS:HD2	1:A:92:LYS:H	1.31	0.92
1:B:92:LYS:HD2	1:B:92:LYS:H	1.35	0.91
1:B:88:VAL:HB	1:B:95:ALA:HB1	1.54	0.88
1:B:178:THR:HB	1:B:181:GLU:HG3	1.56	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:MSE:CE	1:B:104:LEU:HB3	2.04	0.87
1:B:83:MSE:HE3	1:B:104:LEU:HB3	1.55	0.87
1:B:169:ALA:CB	1:B:270:PHE:H	1.87	0.86
1:A:225:THR:HG21	1:A:239:GLY:HA3	1.58	0.85
1:B:202:ILE:HB	1:B:237:LEU:HD22	1.59	0.84
1:A:92:LYS:N	1:A:92:LYS:CD	2.42	0.83
1:B:84:LYS:HD3	1:B:101:ASP:HB2	1.61	0.83
1:B:21:ILE:HD13	1:B:44:LEU:HD22	1.61	0.83
1:B:112:VAL:HG21	1:B:137:PRO:HB2	1.60	0.83
1:A:169:ALA:CB	1:A:270:PHE:H	1.91	0.82
1:A:84:LYS:HD3	1:A:101:ASP:HB2	1.61	0.82
1:A:202:ILE:HB	1:A:237:LEU:HD22	1.62	0.82
1:B:92:LYS:N	1:B:92:LYS:CD	2.42	0.82
1:B:175:ILE:HD12	1:B:175:ILE:H	1.46	0.81
1:A:136:ARG:HB3	1:A:137:PRO:HD3	1.61	0.81
1:A:112:VAL:HG21	1:A:137:PRO:HB2	1.60	0.81
1:B:178:THR:HG22	1:B:180:GLN:N	1.97	0.79
1:A:110:PHE:HE1	1:A:114:LEU:HD11	1.46	0.79
1:A:178:THR:HG22	1:A:180:GLN:N	1.97	0.79
1:B:308:ARG:H	1:B:308:ARG:HD3	1.47	0.78
1:A:88:VAL:HB	1:A:95:ALA:HB1	1.65	0.78
1:B:232:ASN:C	1:B:232:ASN:HD22	1.88	0.77
1:B:28:ASP:HB2	1:B:128:ILE:HD11	1.66	0.76
1:A:33:ARG:HG3	1:A:64:PHE:HZ	1.51	0.76
1:A:88:VAL:CA	1:A:95:ALA:HB1	2.15	0.75
1:B:96:THR:HB	1:B:99:ASP:OD2	1.87	0.75
1:A:232:ASN:C	1:A:232:ASN:HD22	1.88	0.75
1:A:266:ASN:HD22	1:A:267:THR:N	1.86	0.74
1:A:175:ILE:HD12	1:A:175:ILE:H	1.51	0.74
1:B:169:ALA:HB3	1:B:270:PHE:N	2.04	0.73
1:A:169:ALA:HB3	1:A:270:PHE:N	1.99	0.73
1:A:96:THR:HB	1:A:99:ASP:OD2	1.89	0.73
1:A:201:LYS:H	1:A:266:ASN:HD21	1.37	0.72
1:A:83:MSE:HE3	1:A:104:LEU:HD13	1.70	0.72
1:A:167:VAL:HG23	1:A:267:THR:HA	1.72	0.72
1:A:33:ARG:HG3	1:A:64:PHE:CZ	2.24	0.71
1:A:92:LYS:O	1:A:95:ALA:HB3	1.90	0.70
1:B:92:LYS:O	1:B:95:ALA:HB3	1.91	0.70
1:B:110:PHE:HE1	1:B:114:LEU:HD11	1.57	0.70
1:A:96:THR:CG2	1:A:96:THR:CA	2.68	0.69
1:A:110:PHE:CE1	1:A:114:LEU:HD11	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:ARG:H	1:A:308:ARG:HD3	1.56	0.69
1:A:88:VAL:CB	1:A:95:ALA:HB1	2.23	0.68
1:B:88:VAL:CB	1:B:95:ALA:HB1	2.24	0.68
1:A:178:THR:HB	1:A:181:GLU:HG3	1.73	0.68
1:B:83:MSE:HE3	1:B:104:LEU:HD13	1.75	0.68
1:B:162:THR:HG21	1:B:164:GLU:OE1	1.93	0.68
1:B:33:ARG:HG3	1:B:64:PHE:CZ	2.28	0.68
1:B:225:THR:HG21	1:B:239:GLY:HA3	1.77	0.67
1:B:175:ILE:HD12	1:B:175:ILE:N	2.09	0.67
1:B:136:ARG:HB3	1:B:137:PRO:HD3	1.77	0.67
1:A:169:ALA:CB	1:A:269:VAL:CA	2.62	0.67
1:A:79:ASP:HB3	1:A:82:LYS:HE2	1.77	0.67
1:B:201:LYS:H	1:B:266:ASN:HD21	1.41	0.67
1:B:308:ARG:H	1:B:308:ARG:CD	2.05	0.67
1:A:28:ASP:OD2	1:A:30:ARG:HB2	1.94	0.67
1:A:160:GLU:HG2	2:A:332:HOH:O	1.94	0.67
1:B:24:PRO:HA	1:B:49:LEU:HD12	1.77	0.66
1:B:109:TYR:CE1	1:B:137:PRO:HG3	2.30	0.66
1:A:29:GLU:HB2	1:A:62:LEU:HD11	1.78	0.66
1:A:162:THR:HG21	1:A:164:GLU:OE1	1.96	0.66
1:B:33:ARG:HG3	1:B:64:PHE:HZ	1.60	0.66
1:A:253:ALA:O	1:A:257:PRO:HG3	1.95	0.66
1:B:28:ASP:OD2	1:B:30:ARG:HB2	1.95	0.66
1:A:79:ASP:CB	1:A:82:LYS:HE2	2.26	0.65
1:B:167:VAL:HG23	1:B:267:THR:HA	1.78	0.65
1:A:225:THR:CG2	1:A:239:GLY:HA3	2.27	0.65
1:B:96:THR:HB	1:B:99:ASP:CG	2.17	0.65
1:A:79:ASP:C	1:A:81:ASP:H	2.00	0.65
1:B:167:VAL:CG2	1:B:267:THR:HG23	2.27	0.64
1:A:150:ARG:HG3	1:A:150:ARG:HH11	1.62	0.64
1:B:79:ASP:C	1:B:81:ASP:H	2.01	0.64
1:B:162:THR:HG22	1:B:162:THR:O	1.98	0.64
1:A:108:ASN:O	1:A:112:VAL:HG23	1.98	0.63
1:A:96:THR:HB	1:A:99:ASP:CG	2.19	0.63
1:A:54:GLU:O	1:A:57:ASN:HB2	1.99	0.62
1:A:175:ILE:HD12	1:A:175:ILE:N	2.14	0.62
1:B:151:THR:HG22	1:B:297:GLN:O	1.99	0.62
1:A:113:MSE:HE3	1:A:116:LYS:HB3	1.80	0.62
1:B:167:VAL:HG23	1:B:267:THR:HG23	1.82	0.62
1:A:308:ARG:H	1:A:308:ARG:CD	2.12	0.62
1:B:108:ASN:O	1:B:112:VAL:HG23	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:MSE:HE3	1:A:104:LEU:CB	2.24	0.62
1:B:154:VAL:HG11	1:B:192:THR:HG21	1.81	0.62
1:B:15:LEU:HD12	1:B:16:GLY:N	2.15	0.61
1:B:51:GLN:HB2	1:B:54:GLU:HG3	1.83	0.61
1:B:29:GLU:HB2	1:B:62:LEU:HD11	1.82	0.61
1:B:141:ILE:N	1:B:141:ILE:HD12	2.15	0.61
1:B:169:ALA:CB	1:B:269:VAL:CA	2.63	0.61
1:A:49:LEU:HD12	1:A:107:VAL:HG12	1.83	0.60
1:B:79:ASP:HB3	1:B:82:LYS:HE2	1.83	0.60
1:A:92:LYS:N	1:A:95:ALA:HB2	2.16	0.60
1:A:56:ARG:HG2	1:A:66:ASP:OD1	2.02	0.60
1:A:169:ALA:CB	1:A:270:PHE:N	2.63	0.60
1:A:128:ILE:HG22	1:A:128:ILE:O	2.02	0.59
1:B:56:ARG:HG2	1:B:66:ASP:OD1	2.02	0.59
1:A:159:ARG:HD3	1:A:162:THR:OG1	2.02	0.59
1:A:205:LEU:O	1:A:271:PRO:HG3	2.03	0.59
1:A:167:VAL:CG2	1:A:267:THR:HG23	2.33	0.59
1:B:92:LYS:N	1:B:95:ALA:HB2	2.18	0.59
1:B:148:ILE:HG12	1:B:188:ASN:ND2	2.17	0.59
1:B:54:GLU:O	1:B:57:ASN:HB2	2.03	0.59
1:B:113:MSE:HE3	1:B:116:LYS:HB3	1.85	0.59
1:B:80:PHE:O	1:B:84:LYS:HB2	2.03	0.59
1:B:175:ILE:H	1:B:175:ILE:CD1	2.15	0.59
1:A:91:ARG:C	1:A:92:LYS:HD2	2.21	0.59
1:A:162:THR:O	1:A:162:THR:HG22	2.02	0.59
1:B:103:MSE:C	1:B:105:ARG:H	2.06	0.59
1:B:28:ASP:HB2	1:B:128:ILE:CD1	2.33	0.58
1:A:59:LEU:O	1:A:64:PHE:HB2	2.03	0.58
1:B:148:ILE:N	1:B:148:ILE:HD12	2.18	0.58
1:A:167:VAL:HG23	1:A:267:THR:HG23	1.84	0.58
1:B:88:VAL:C	1:B:90:VAL:H	2.07	0.58
1:A:159:ARG:HB2	1:A:289:PHE:CD2	2.39	0.58
1:B:59:LEU:O	1:B:64:PHE:HB2	2.04	0.58
1:A:109:TYR:CE1	1:A:137:PRO:HG3	2.38	0.57
1:A:141:ILE:HD12	1:A:141:ILE:N	2.19	0.57
1:A:272:ASP:OD2	1:A:274:GLN:HB2	2.04	0.57
1:A:88:VAL:HG12	1:A:96:THR:H	1.68	0.57
1:A:88:VAL:O	1:A:95:ALA:CB	2.52	0.57
1:B:79:ASP:CB	1:B:82:LYS:HE2	2.34	0.57
1:B:110:PHE:CE1	1:B:114:LEU:HD11	2.39	0.57
1:B:169:ALA:CB	1:B:270:PHE:N	2.60	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:ARG:HB2	1:B:289:PHE:CD2	2.40	0.57
1:B:80:PHE:CE1	1:B:105:ARG:NH2	2.73	0.57
1:B:162:THR:CG2	1:B:164:GLU:OE1	2.52	0.57
1:B:248:VAL:HG22	1:B:251:THR:HG23	1.86	0.57
1:A:322:THR:O	1:A:325:GLN:HB2	2.04	0.57
1:A:88:VAL:HG12	1:A:96:THR:N	2.20	0.57
1:A:48:ILE:HG21	1:A:55:VAL:HG11	1.86	0.57
1:B:51:GLN:HB2	1:B:54:GLU:CG	2.35	0.57
1:A:80:PHE:CE1	1:A:105:ARG:NH2	2.72	0.57
1:B:204:MSE:SE	1:B:237:LEU:HD11	2.55	0.57
1:A:112:VAL:HG21	1:A:137:PRO:CB	2.33	0.56
1:A:175:ILE:H	1:A:175:ILE:CD1	2.17	0.56
1:A:225:THR:HG21	1:A:239:GLY:CA	2.31	0.56
1:A:148:ILE:HD12	1:A:148:ILE:N	2.21	0.56
1:A:26:GLY:O	1:A:55:VAL:HG22	2.06	0.56
1:A:151:THR:HG22	1:A:297:GLN:O	2.06	0.56
1:B:225:THR:CG2	1:B:239:GLY:HA3	2.36	0.56
1:B:266:ASN:HD22	1:B:267:THR:N	2.04	0.55
1:A:135:VAL:O	1:A:139:LEU:HG	2.06	0.55
1:B:201:LYS:N	1:B:266:ASN:HD21	2.03	0.55
1:A:28:ASP:HB2	1:A:128:ILE:CD1	2.31	0.55
1:A:19:MSE:SE	1:A:327:ILE:HD13	2.57	0.55
1:B:201:LYS:HD3	1:B:264:GLN:HE21	1.71	0.55
1:B:209:THR:HG23	1:B:241:LEU:HD23	1.87	0.55
1:B:248:VAL:HG23	1:B:250:GLU:HG2	1.88	0.55
1:A:88:VAL:C	1:A:90:VAL:H	2.10	0.55
1:A:27:ASN:O	1:A:58:LEU:CD1	2.55	0.54
1:A:80:PHE:HE1	1:A:105:ARG:NH2	2.06	0.54
1:A:82:LYS:HE3	1:A:117:MSE:CE	2.29	0.54
1:A:103:MSE:C	1:A:105:ARG:H	2.10	0.54
1:B:87:PHE:HB2	1:B:104:LEU:HD11	1.90	0.54
1:B:88:VAL:CA	1:B:95:ALA:HB1	2.37	0.54
1:B:158:ASN:OD1	1:B:165:ARG:HD3	2.08	0.54
1:B:232:ASN:C	1:B:232:ASN:ND2	2.59	0.54
1:A:292:ILE:O	1:A:293:GLY:O	2.26	0.54
1:B:150:ARG:HG3	1:B:150:ARG:HH11	1.73	0.54
1:A:209:THR:HG23	1:A:241:LEU:HD23	1.88	0.54
1:B:202:ILE:O	1:B:261:VAL:HG11	2.08	0.54
1:B:308:ARG:HD3	1:B:308:ARG:N	2.21	0.54
1:A:80:PHE:O	1:A:84:LYS:HB2	2.08	0.53
1:B:131:THR:HG23	1:B:305:ASP:OD2	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:LEU:CD1	1:A:107:VAL:HG12	2.38	0.53
1:A:51:GLN:HB2	1:A:54:GLU:HG3	1.88	0.53
1:B:173:ILE:HG22	1:B:174:ASN:N	2.24	0.53
1:A:162:THR:CG2	1:A:164:GLU:OE1	2.57	0.53
1:A:83:MSE:HG2	1:A:104:LEU:HD13	1.91	0.53
1:B:292:ILE:O	1:B:293:GLY:O	2.27	0.53
1:A:79:ASP:C	1:A:81:ASP:N	2.61	0.53
1:B:96:THR:O	1:B:99:ASP:N	2.42	0.53
1:A:51:GLN:HB2	1:A:54:GLU:CG	2.39	0.52
1:B:79:ASP:C	1:B:81:ASP:N	2.62	0.52
1:B:56:ARG:HD3	1:B:66:ASP:OD2	2.08	0.52
1:B:248:VAL:CG2	1:B:251:THR:HG23	2.39	0.52
1:A:83:MSE:HE3	1:A:104:LEU:CD1	2.38	0.52
1:A:204:MSE:SE	1:A:237:LEU:HD11	2.59	0.52
1:A:148:ILE:HG22	1:A:149:SER:N	2.25	0.52
1:A:201:LYS:N	1:A:266:ASN:HD21	2.06	0.52
1:A:136:ARG:HB3	1:A:137:PRO:CD	2.38	0.52
1:A:202:ILE:HD13	1:A:237:LEU:HD22	1.92	0.52
1:A:306:LEU:HD23	1:A:318:LEU:HD12	1.91	0.52
1:B:91:ARG:C	1:B:92:LYS:HD2	2.30	0.52
1:B:174:ASN:HB3	1:B:177:PRO:HG3	1.92	0.52
1:A:150:ARG:HG3	1:A:150:ARG:NH1	2.25	0.52
1:A:24:PRO:HA	1:A:49:LEU:HD12	1.92	0.51
1:B:91:ARG:HG2	1:B:91:ARG:HH11	1.74	0.51
1:A:154:VAL:HG11	1:A:192:THR:HG21	1.91	0.51
1:B:51:GLN:CD	1:B:54:GLU:HG3	2.31	0.51
1:B:205:LEU:O	1:B:271:PRO:HG3	2.11	0.51
1:B:219:ASP:O	1:B:223:GLU:HB2	2.09	0.51
1:A:88:VAL:HG11	1:A:97:LEU:N	2.25	0.51
1:B:178:THR:HB	1:B:181:GLU:CG	2.36	0.51
1:A:88:VAL:HA	1:A:95:ALA:HB1	1.90	0.51
1:A:217:GLN:CD	1:A:217:GLN:H	2.13	0.51
1:B:48:ILE:HG21	1:B:55:VAL:HG11	1.92	0.51
1:B:225:THR:OG1	1:B:239:GLY:HA3	2.10	0.51
1:B:86:ALA:O	1:B:90:VAL:HG23	2.11	0.50
1:A:92:LYS:NZ	1:A:92:LYS:CD	2.65	0.50
1:A:80:PHE:CE2	1:A:83:MSE:HE2	2.46	0.50
1:A:7:PHE:CE2	1:A:317:LYS:HG2	2.47	0.50
1:A:79:ASP:O	1:A:81:ASP:N	2.45	0.50
1:B:84:LYS:HD2	1:B:97:LEU:CD2	2.41	0.50
1:B:83:MSE:HG2	1:B:104:LEU:HD13	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:TYR:O	1:A:320:ILE:HG13	2.11	0.49
1:B:88:VAL:HG12	1:B:96:THR:H	1.77	0.49
1:A:254:ILE:HG22	1:A:255:LYS:N	2.26	0.49
1:B:80:PHE:HE1	1:B:105:ARG:NH2	2.09	0.49
1:B:259:SER:HB3	1:B:262:ALA:HB3	1.94	0.49
1:A:48:ILE:HD12	1:A:48:ILE:N	2.28	0.49
1:A:167:VAL:O	1:A:168:PHE:HD2	1.96	0.49
1:A:232:ASN:C	1:A:232:ASN:ND2	2.59	0.49
1:A:87:PHE:HB2	1:A:104:LEU:HD11	1.94	0.48
1:A:201:LYS:HD3	1:A:264:GLN:HE21	1.78	0.48
1:A:67:GLN:O	1:A:68:ASP:HB2	2.13	0.48
1:A:86:ALA:O	1:A:90:VAL:HG23	2.14	0.48
1:B:79:ASP:O	1:B:81:ASP:N	2.46	0.48
1:B:266:ASN:HD22	1:B:266:ASN:H	1.60	0.48
1:A:88:VAL:O	1:A:95:ALA:HB1	2.13	0.48
1:B:272:ASP:OD2	1:B:274:GLN:HB2	2.14	0.48
1:B:128:ILE:HG22	1:B:128:ILE:O	2.13	0.48
1:B:177:PRO:O	1:B:220:LYS:HD2	2.13	0.48
1:A:128:ILE:O	1:A:128:ILE:CG2	2.62	0.48
1:A:232:ASN:ND2	1:A:234:ASP:H	2.11	0.48
1:A:248:VAL:CG2	1:A:251:THR:HG23	2.43	0.48
1:A:38:LEU:HD23	1:A:43:LEU:HD12	1.95	0.48
1:A:61:LYS:HE3	1:A:61:LYS:HA	1.95	0.48
1:A:110:PHE:CE1	1:A:114:LEU:HD21	2.49	0.48
1:A:266:ASN:HD22	1:A:266:ASN:H	1.61	0.47
1:A:113:MSE:HE3	1:A:113:MSE:HA	1.95	0.47
1:B:77:TYR:HB3	1:B:80:PHE:HB2	1.95	0.47
1:B:141:ILE:N	1:B:141:ILE:CD1	2.77	0.47
1:A:158:ASN:OD1	1:A:165:ARG:HD3	2.14	0.47
1:B:128:ILE:O	1:B:128:ILE:CG2	2.62	0.47
1:B:148:ILE:HG22	1:B:149:SER:N	2.29	0.47
1:B:174:ASN:CB	1:B:177:PRO:HG3	2.44	0.47
1:A:51:GLN:CD	1:A:54:GLU:HG3	2.35	0.47
1:A:79:ASP:HB3	1:A:82:LYS:HG2	1.97	0.47
1:A:88:VAL:C	1:A:95:ALA:HB1	2.35	0.47
1:A:173:ILE:HG22	1:A:174:ASN:N	2.29	0.47
1:A:266:ASN:HD22	1:A:267:THR:H	1.62	0.47
1:B:150:ARG:HD3	1:B:174:ASN:OD1	2.14	0.47
1:B:159:ARG:HD3	1:B:162:THR:OG1	2.14	0.47
1:A:7:PHE:O	1:A:11:ARG:HB2	2.14	0.47
1:B:77:TYR:C	1:B:79:ASP:H	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:LYS:HG3	1:A:83:MSE:N	2.30	0.47
1:B:135:VAL:O	1:B:139:LEU:HG	2.15	0.47
1:B:167:VAL:HG21	1:B:267:THR:HG23	1.96	0.47
1:B:184:GLU:O	1:B:188:ASN:HB2	2.15	0.47
1:A:87:PHE:HD2	1:A:100:ALA:HB2	1.81	0.46
1:A:91:ARG:C	1:A:95:ALA:HB2	2.36	0.46
1:A:52:SER:HA	1:A:71:ILE:HG21	1.97	0.46
1:A:39:LYS:HD3	1:A:69:TYR:CE2	2.51	0.46
1:B:136:ARG:HH11	1:B:136:ARG:HG2	1.81	0.46
1:B:217:GLN:CD	1:B:217:GLN:H	2.18	0.46
1:A:177:PRO:O	1:A:220:LYS:HD2	2.16	0.46
1:A:136:ARG:HG2	1:A:136:ARG:HH11	1.80	0.46
1:A:21:ILE:HD13	1:A:44:LEU:HD22	1.97	0.46
1:B:27:ASN:O	1:B:58:LEU:CD1	2.63	0.46
1:A:87:PHE:HE1	1:A:141:ILE:HD11	1.81	0.45
1:A:88:VAL:O	1:A:95:ALA:HB2	2.17	0.45
1:B:202:ILE:CB	1:B:237:LEU:HD22	2.37	0.45
1:A:46:PRO:HB2	1:A:48:ILE:HD11	1.98	0.45
1:A:202:ILE:HD13	1:A:237:LEU:CD2	2.45	0.45
1:B:232:ASN:ND2	1:B:234:ASP:H	2.14	0.45
1:A:266:ASN:HD22	1:A:266:ASN:C	2.19	0.45
1:B:49:LEU:CD1	1:B:107:VAL:HG12	2.47	0.45
1:B:103:MSE:C	1:B:105:ARG:N	2.70	0.45
1:A:180:GLN:NE2	1:A:227:ILE:HG21	2.32	0.45
1:A:56:ARG:HD3	1:A:66:ASP:OD2	2.17	0.44
1:A:96:THR:CB	1:A:99:ASP:OD2	2.63	0.44
1:A:308:ARG:HD3	1:A:308:ARG:N	2.29	0.44
1:B:150:ARG:HG3	1:B:150:ARG:NH1	2.31	0.44
1:B:88:VAL:O	1:B:90:VAL:N	2.50	0.44
1:B:92:LYS:O	1:B:95:ALA:CB	2.64	0.44
1:A:83:MSE:CE	1:A:104:LEU:HD13	2.44	0.44
1:A:85:GLU:C	1:A:87:PHE:N	2.70	0.44
1:B:88:VAL:C	1:B:90:VAL:N	2.69	0.44
1:B:113:MSE:HE3	1:B:113:MSE:HA	1.99	0.44
1:A:32:VAL:HG21	1:A:58:LEU:HD13	2.00	0.44
1:A:85:GLU:O	1:A:87:PHE:N	2.50	0.44
1:A:131:THR:HG23	1:A:305:ASP:OD2	2.16	0.44
1:A:327:ILE:C	1:A:329:SER:H	2.21	0.44
1:A:274:GLN:NE2	1:B:271:PRO:HB3	2.33	0.44
1:B:49:LEU:HD12	1:B:107:VAL:HG12	1.98	0.44
1:B:85:GLU:C	1:B:87:PHE:N	2.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:VAL:O	1:B:95:ALA:CB	2.66	0.44
1:A:126:GLY:HA3	1:A:305:ASP:OD1	2.17	0.44
1:A:306:LEU:HD21	1:A:315:ILE:HA	2.00	0.44
1:A:178:THR:HB	1:A:181:GLU:CG	2.46	0.43
1:B:155:PHE:CZ	1:B:294:PRO:HG3	2.54	0.43
1:B:48:ILE:N	1:B:48:ILE:HD12	2.32	0.43
1:A:148:ILE:CG2	1:A:149:SER:N	2.81	0.43
1:B:11:ARG:HD3	1:B:43:LEU:HD11	2.00	0.43
1:B:21:ILE:HG21	1:B:124:VAL:HG23	2.00	0.43
1:A:80:PHE:HE2	1:A:83:MSE:HE2	1.84	0.43
1:A:318:LEU:HD23	1:A:318:LEU:HA	1.90	0.43
1:B:148:ILE:HG22	1:B:150:ARG:H	1.83	0.43
1:B:16:GLY:C	1:B:18:ASN:H	2.22	0.43
1:A:79:ASP:HB3	1:A:82:LYS:CE	2.48	0.43
1:A:232:ASN:HD22	1:A:234:ASP:H	1.66	0.43
1:B:223:GLU:O	1:B:227:ILE:HG13	2.18	0.43
1:A:27:ASN:O	1:A:58:LEU:HD12	2.19	0.43
1:B:88:VAL:HG12	1:B:96:THR:N	2.32	0.43
1:A:88:VAL:HA	1:A:95:ALA:CB	2.49	0.43
1:A:180:GLN:O	1:A:184:GLU:HG3	2.19	0.43
1:A:225:THR:HG21	1:A:239:GLY:N	2.34	0.43
1:B:327:ILE:C	1:B:329:SER:H	2.21	0.43
1:A:77:TYR:CD2	1:A:78:ALA:N	2.87	0.42
1:B:82:LYS:HG3	1:B:83:MSE:N	2.33	0.42
1:A:21:ILE:N	1:A:21:ILE:HD12	2.34	0.42
1:A:202:ILE:CB	1:A:237:LEU:HD22	2.41	0.42
1:B:6:LEU:HD12	1:B:165:ARG:CZ	2.48	0.42
1:B:266:ASN:HD22	1:B:266:ASN:N	2.16	0.42
1:A:252:ALA:O	1:A:253:ALA:C	2.56	0.42
1:B:190:ALA:O	1:B:194:LYS:HG3	2.19	0.42
1:A:110:PHE:HE1	1:A:114:LEU:CD1	2.26	0.42
1:B:316:TYR:O	1:B:320:ILE:HG13	2.20	0.42
1:A:138:ALA:O	1:A:142:ILE:O	2.37	0.42
1:A:205:LEU:HA	1:A:241:LEU:O	2.20	0.42
1:B:3:ILE:HD12	2:B:348:HOH:O	2.20	0.42
1:B:33:ARG:HH11	1:B:33:ARG:HB2	1.84	0.42
1:B:103:MSE:O	1:B:105:ARG:N	2.53	0.42
1:B:178:THR:O	1:B:179:ALA:C	2.54	0.42
1:B:318:LEU:HD23	1:B:318:LEU:HA	1.89	0.42
1:A:77:TYR:C	1:A:79:ASP:H	2.23	0.42
1:A:271:PRO:HB3	1:B:274:GLN:NE2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:ARG:HG2	1:A:136:ARG:NH1	2.34	0.42
1:A:148:ILE:HG22	1:A:150:ARG:H	1.85	0.42
1:A:296:LEU:HD12	1:A:303:VAL:O	2.18	0.42
1:B:19:MSE:HA	1:B:19:MSE:HE2	2.01	0.42
1:A:141:ILE:N	1:A:141:ILE:CD1	2.83	0.42
1:A:202:ILE:O	1:A:261:VAL:HG11	2.20	0.42
1:B:284:GLN:HG2	1:B:285:ARG:HD2	2.01	0.42
1:A:92:LYS:O	1:A:95:ALA:N	2.53	0.41
1:A:78:ALA:O	1:A:79:ASP:CG	2.58	0.41
1:B:173:ILE:N	1:B:173:ILE:HD12	2.35	0.41
1:A:1:MSE:HE2	1:A:165:ARG:NH2	2.35	0.41
1:B:33:ARG:NH1	1:B:33:ARG:CB	2.84	0.41
1:A:124:VAL:HG21	1:A:319:ALA:HA	2.00	0.41
1:B:84:LYS:HD2	1:B:97:LEU:HD21	2.02	0.41
1:B:252:ALA:O	1:B:253:ALA:C	2.59	0.41
1:A:15:LEU:HD12	1:A:16:GLY:N	2.35	0.41
1:A:202:ILE:HD12	1:A:202:ILE:H	1.85	0.41
1:A:248:VAL:HA	1:A:249:PRO:HD3	1.90	0.41
1:B:88:VAL:HG11	1:B:97:LEU:N	2.35	0.41
1:A:56:ARG:HG2	1:A:66:ASP:CG	2.41	0.41
1:A:259:SER:HB3	1:A:262:ALA:HB3	2.03	0.41
1:A:286:LEU:O	1:B:159:ARG:NH2	2.53	0.41
1:B:87:PHE:HD2	1:B:100:ALA:HB2	1.86	0.41
1:B:92:LYS:O	1:B:95:ALA:N	2.53	0.41
1:A:40:PHE:HE1	1:A:69:TYR:HH	1.67	0.41
1:A:91:ARG:HH11	1:A:91:ARG:HG2	1.84	0.41
1:A:45:GLU:HA	1:A:46:PRO:HD3	1.87	0.41
1:A:109:TYR:CZ	1:A:137:PRO:HG3	2.56	0.41
1:B:67:GLN:O	1:B:68:ASP:HB2	2.21	0.41
1:B:85:GLU:O	1:B:87:PHE:N	2.53	0.41
1:A:52:SER:OG	1:A:56:ARG:NH2	2.44	0.41
1:A:88:VAL:C	1:A:90:VAL:N	2.73	0.41
1:A:261:VAL:O	1:A:262:ALA:C	2.59	0.40
1:B:142:ILE:HG22	1:B:143:LYS:N	2.36	0.40
1:B:45:GLU:HA	1:B:46:PRO:HD3	1.91	0.40
1:B:254:ILE:HG22	1:B:255:LYS:N	2.36	0.40
1:A:323:ALA:O	1:A:327:ILE:HG12	2.20	0.40
1:A:74:PRO:HA	1:A:110:PHE:CE2	2.57	0.40
1:A:83:MSE:HG3	1:A:117:MSE:SE	2.72	0.40
1:B:46:PRO:HB2	1:B:48:ILE:HD11	2.03	0.40
1:A:83:MSE:HG2	1:A:104:LEU:CD1	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:PHE:O	1:B:11:ARG:HB2	2.21	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	327/337 (97%)	284 (87%)	32 (10%)	11 (3%)	3	7
1	B	327/337 (97%)	284 (87%)	32 (10%)	11 (3%)	3	7
All	All	654/674 (97%)	568 (87%)	64 (10%)	22 (3%)	3	7

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	92	LYS
1	A	293	GLY
1	B	293	GLY
1	B	78	ALA
1	B	92	LYS
1	B	170	ASP
1	A	18	ASN
1	A	78	ALA
1	A	79	ASP
1	A	80	PHE
1	A	89	GLU
1	A	170	ASP
1	B	79	ASP
1	B	80	PHE
1	B	89	GLU
1	B	104	LEU
1	A	141	ILE

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Mol	Chain	Res	Type
1	B	86	ALA
1	A	86	ALA
1	A	104	LEU
1	B	169	ALA
1	B	141	ILE

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	265/256 (104%)	239 (90%)	26 (10%)	6	16
1	B	265/256 (104%)	239 (90%)	26 (10%)	6	16
All	All	530/512 (104%)	478 (90%)	52 (10%)	6	16

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	2	SER
1	A	4	ARG
1	A	58	LEU
1	A	61	LYS
1	A	68	ASP
1	A	81	ASP
1	A	89	GLU
1	A	92	LYS
1	A	94	LYS
1	A	97	LEU
1	A	107	VAL
1	A	113	MSE
1	A	151	THR
1	A	159	ARG
1	A	167	VAL
1	A	217	GLN
1	A	226	GLU

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Mol	Chain	Res	Type
1	A	232	ASN
1	A	237	LEU
1	A	250	GLU
1	A	266	ASN
1	A	280	TYR
1	A	285	ARG
1	A	308	ARG
1	A	318	LEU
1	B	1	MSE
1	B	4	ARG
1	B	57	ASN
1	B	58	LEU
1	B	61	LYS
1	B	68	ASP
1	B	74	PRO
1	B	81	ASP
1	B	89	GLU
1	B	92	LYS
1	B	97	LEU
1	B	107	VAL
1	B	151	THR
1	B	159	ARG
1	B	165	ARG
1	B	217	GLN
1	B	232	ASN
1	B	237	LEU
1	B	248	VAL
1	B	250	GLU
1	B	266	ASN
1	B	280	TYR
1	B	285	ARG
1	B	308	ARG
1	B	313	GLU
1	B	318	LEU

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	57	ASN
1	A	129	HIS
1	A	180	GLN
1	A	188	ASN

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Mol	Chain	Res	Type
1	A	232	ASN
1	A	264	GLN
1	A	266	ASN
1	B	57	ASN
1	B	129	HIS
1	B	180	GLN
1	B	188	ASN
1	B	232	ASN
1	B	264	GLN
1	B	266	ASN
1	B	274	GLN

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 oligosaccharides 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, 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	319/337 (94%)	-0.01	7 (2%) 62 61	42, 74, 121, 138	0
1	B	319/337 (94%)	-0.35	2 (0%) 85 85	40, 61, 106, 131	0
All	All	638/674 (94%)	-0.18	9 (1%) 73 73	40, 67, 116, 138	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	95	ALA	5.7
1	A	93	GLY	5.6
1	A	69	TYR	3.2
1	A	100	ALA	2.9
1	B	104	LEU	2.5
1	B	100	ALA	2.3
1	A	260	ALA	2.2
1	A	104	LEU	2.2
1	A	97	LEU	2.1

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.