



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 09:30 am GMT

PDB ID : 2QC7
Title : Crystal structure of the protein-disulfide isomerase related chaperone ERp29
Authors : Barak, N.N.; Sevvana, M.; Neumann, P.; Malesevic, M.; Naumann, K.; Fischer, G.; Sheldrick, G.M.; Stubbs, M.T.; Ferrari, D.M.
Deposited on : 2007-06-19
Resolution : 2.90 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

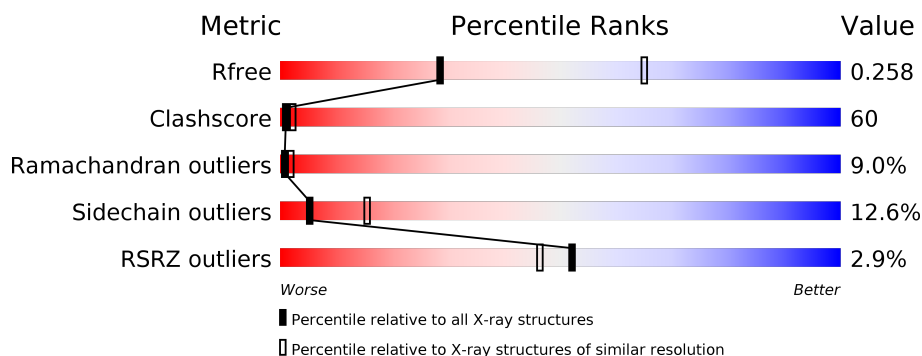
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

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

Mol	Chain	Length	Quality of chain
1	A	240	<div> <div>2%</div> <div> <div></div> <div>28%</div> <div>52%</div> <div>14%</div> <div>7%</div> </div> </div>
1	B	240	<div> <div>3%</div> <div> <div></div> <div>26%</div> <div>50%</div> <div>16%</div> <div>7%</div> </div> </div>

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 3526 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 Endoplasmic reticulum protein ERp29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	224	Total	C	N	O	S	0	0	0
			1763	1131	291	335	6			
1	B	224	Total	C	N	O	S	0	0	0
			1763	1131	291	335	6			

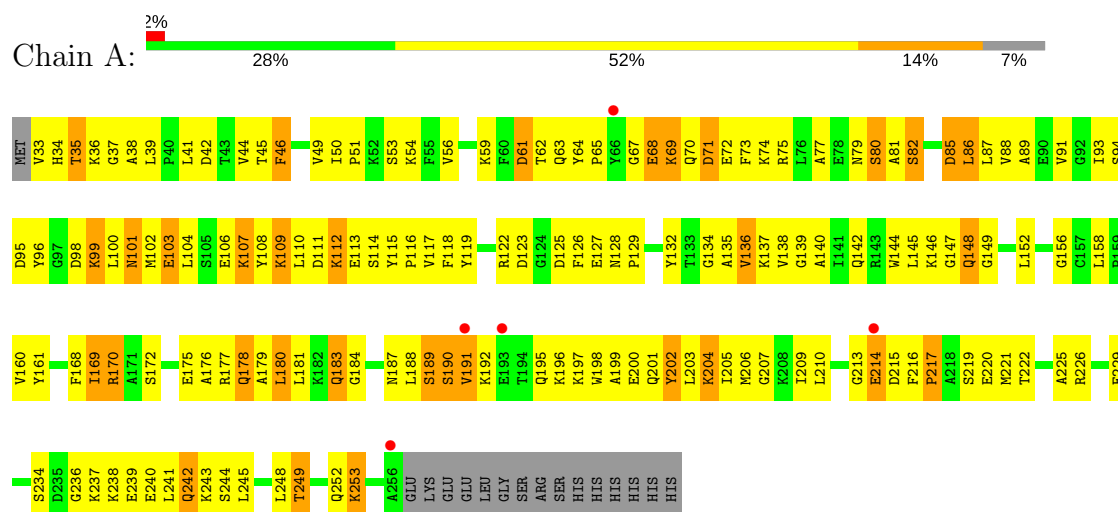
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	32	MET	-	INITIATING METHIONINE	UNP P30040
A	33	VAL	-	EXPRESSION TAG	UNP P30040
A	262	GLY	-	EXPRESSION TAG	UNP P30040
A	263	SER	-	EXPRESSION TAG	UNP P30040
A	264	ARG	-	EXPRESSION TAG	UNP P30040
A	265	SER	-	EXPRESSION TAG	UNP P30040
A	266	HIS	-	EXPRESSION TAG	UNP P30040
A	267	HIS	-	EXPRESSION TAG	UNP P30040
A	268	HIS	-	EXPRESSION TAG	UNP P30040
A	269	HIS	-	EXPRESSION TAG	UNP P30040
A	270	HIS	-	EXPRESSION TAG	UNP P30040
A	271	HIS	-	EXPRESSION TAG	UNP P30040
B	32	MET	-	INITIATING METHIONINE	UNP P30040
B	33	VAL	-	EXPRESSION TAG	UNP P30040
B	262	GLY	-	EXPRESSION TAG	UNP P30040
B	263	SER	-	EXPRESSION TAG	UNP P30040
B	264	ARG	-	EXPRESSION TAG	UNP P30040
B	265	SER	-	EXPRESSION TAG	UNP P30040
B	266	HIS	-	EXPRESSION TAG	UNP P30040
B	267	HIS	-	EXPRESSION TAG	UNP P30040
B	268	HIS	-	EXPRESSION TAG	UNP P30040
B	269	HIS	-	EXPRESSION TAG	UNP P30040
B	270	HIS	-	EXPRESSION TAG	UNP P30040
B	271	HIS	-	EXPRESSION TAG	UNP P30040

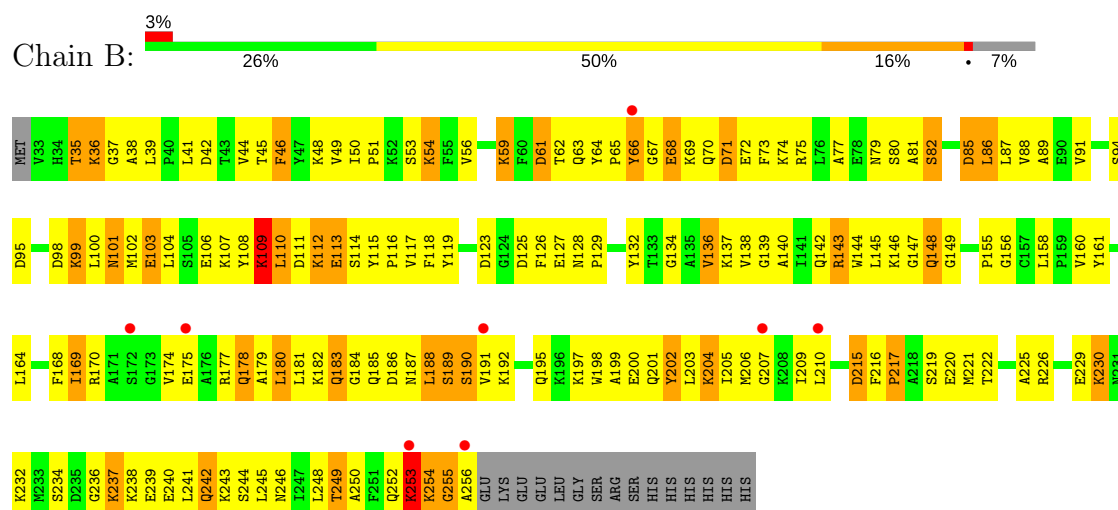
3 Residue-property plots

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

• Molecule 1: Endoplasmic reticulum protein ERp29



• Molecule 1: Endoplasmic reticulum protein ERp29



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	58.04Å 68.03Å 70.09Å 90.00° 107.46° 90.00°	Depositor
Resolution (Å)	28.98 – 2.90 28.82 – 2.90	Depositor EDS
% Data completeness (in resolution range)	94.7 (28.98-2.90) 94.7 (28.82-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.265 , 0.279 0.260 , 0.258	Depositor DCC
R_{free} test set	545 reflections (5.18%)	DCC
Wilson B-factor (Å ²)	77.2	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 70.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3526	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.22% 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 [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	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.75	0/1797	0.79	1/2417 (0.0%)
1	B	0.68	0/1797	0.76	2/2417 (0.1%)
All	All	0.72	0/3594	0.78	3/4834 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	86	LEU	CA-CB-CG	-5.61	102.40	115.30
1	B	123	ASP	CA-CB-CG	5.15	124.72	113.40
1	A	86	LEU	CA-CB-CG	-5.02	103.76	115.30

There are no chirality outliers.

There are no planarity outliers.

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	A	1763	0	1765	198	0
1	B	1763	0	1765	226	0
All	All	3526	0	3530	421	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 60.

All (421) 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:183:GLN:HG3	1:A:184:GLY:H	1.14	1.12
1:A:192:LYS:HB2	1:A:195:GLN:HG3	1.35	1.07
1:B:71:ASP:HA	1:B:74:LYS:HE2	1.41	1.03
1:A:71:ASP:HA	1:A:74:LYS:HD2	1.38	1.00
1:B:143:ARG:HG3	1:B:143:ARG:HH11	1.27	0.99
1:B:207:GLY:HA2	1:B:210:LEU:HD12	1.45	0.98
1:B:68:GLU:HG2	1:B:69:LYS:N	1.81	0.96
1:A:132:TYR:HB2	1:A:144:TRP:CZ3	2.02	0.94
1:A:207:GLY:HA2	1:A:210:LEU:HD12	1.48	0.94
1:B:183:GLN:HG3	1:B:184:GLY:H	1.31	0.93
1:A:50:ILE:HB	1:A:51:PRO:HD3	1.50	0.92
1:B:132:TYR:HB2	1:B:144:TRP:CZ3	2.03	0.92
1:B:174:VAL:HA	1:B:177:ARG:HH21	1.35	0.92
1:B:50:ILE:HB	1:B:51:PRO:HD3	1.52	0.91
1:A:33:VAL:HG12	1:A:34:HIS:N	1.85	0.91
1:B:143:ARG:CG	1:B:143:ARG:HH11	1.84	0.90
1:A:253:LYS:H	1:A:253:LYS:HD2	1.31	0.90
1:A:33:VAL:HG12	1:A:34:HIS:H	1.35	0.90
1:B:185:GLN:O	1:B:188:LEU:HD11	1.73	0.89
1:B:234:SER:OG	1:B:237:LYS:HD3	1.73	0.88
1:A:95:ASP:HB2	1:A:112:LYS:HE2	1.57	0.86
1:B:175:GLU:O	1:B:178:GLN:HB3	1.78	0.83
1:A:75:ARG:HD2	1:A:138:VAL:HG21	1.60	0.83
1:B:75:ARG:HD2	1:B:138:VAL:HG21	1.60	0.83
1:A:39:LEU:O	1:A:89:ALA:HB1	1.80	0.82
1:B:138:VAL:O	1:B:142:GLN:HG3	1.78	0.82
1:A:71:ASP:HA	1:A:74:LYS:CD	2.11	0.81
1:B:139:GLY:C	1:B:143:ARG:HH12	1.81	0.81
1:B:174:VAL:HA	1:B:177:ARG:NH2	1.95	0.81
1:B:109:LYS:HD2	1:B:109:LYS:N	1.94	0.81
1:B:109:LYS:HD2	1:B:109:LYS:H	1.46	0.80
1:A:108:TYR:O	1:A:110:LEU:HG	1.82	0.80
1:B:175:GLU:HB3	1:B:178:GLN:NE2	1.97	0.79
1:B:39:LEU:O	1:B:89:ALA:HB1	1.82	0.79
1:A:192:LYS:CB	1:A:195:GLN:HG3	2.11	0.79
1:B:139:GLY:O	1:B:143:ARG:NH1	2.18	0.77
1:B:143:ARG:HG3	1:B:143:ARG:NH1	1.97	0.77
1:A:175:GLU:O	1:A:178:GLN:HB3	1.84	0.76
1:A:183:GLN:HG3	1:A:184:GLY:N	1.97	0.76
1:A:138:VAL:O	1:A:142:GLN:HG3	1.85	0.76
1:A:241:LEU:O	1:A:244:SER:N	2.19	0.76
1:B:175:GLU:HB3	1:B:178:GLN:HE22	1.52	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:LYS:HG3	1:A:123:ASP:OD1	1.87	0.73
1:B:241:LEU:O	1:B:244:SER:N	2.21	0.73
1:A:132:TYR:CZ	1:A:134:GLY:HA3	2.24	0.73
1:A:175:GLU:HB3	1:A:178:GLN:NE2	2.04	0.71
1:B:252:GLN:O	1:B:255:GLY:HA3	1.90	0.71
1:B:175:GLU:OE1	1:B:178:GLN:NE2	2.23	0.71
1:B:108:TYR:O	1:B:110:LEU:HD23	1.91	0.70
1:B:160:VAL:HG13	1:B:161:TYR:HD1	1.56	0.70
1:B:132:TYR:CZ	1:B:134:GLY:HA3	2.27	0.70
1:A:117:VAL:HG12	1:A:119:TYR:HE1	1.57	0.70
1:B:117:VAL:HG12	1:B:119:TYR:HE1	1.55	0.69
1:B:186:ASP:O	1:B:189:SER:N	2.23	0.69
1:A:253:LYS:CD	1:A:253:LYS:H	2.05	0.69
1:B:205:ILE:O	1:B:209:ILE:HG13	1.93	0.69
1:B:175:GLU:CA	1:B:178:GLN:HE21	2.05	0.69
1:B:183:GLN:CG	1:B:184:GLY:H	2.05	0.68
1:A:175:GLU:CA	1:A:178:GLN:HE21	2.06	0.68
1:A:236:GLY:O	1:A:240:GLU:HG3	1.93	0.68
1:B:110:LEU:HD11	1:B:117:VAL:HG11	1.74	0.68
1:A:205:ILE:O	1:A:209:ILE:HG13	1.94	0.68
1:A:221:MET:CE	1:A:249:THR:HG22	2.23	0.68
1:B:59:LYS:HD3	1:B:61:ASP:HB2	1.75	0.68
1:A:44:VAL:O	1:A:45:THR:C	2.31	0.68
1:A:221:MET:HE3	1:A:249:THR:HG22	1.76	0.67
1:B:44:VAL:O	1:B:45:THR:C	2.30	0.67
1:A:160:VAL:HG13	1:A:161:TYR:HD1	1.59	0.67
1:B:100:LEU:C	1:B:102:MET:H	1.98	0.67
1:B:35:THR:CG2	1:B:74:LYS:HA	2.25	0.67
1:A:100:LEU:C	1:A:102:MET:H	1.98	0.67
1:A:175:GLU:HB3	1:A:178:GLN:HE22	1.61	0.66
1:B:160:VAL:HG13	1:B:161:TYR:CD1	2.31	0.66
1:B:169:ILE:HG22	1:B:170:ARG:N	2.11	0.66
1:B:221:MET:CE	1:B:249:THR:HG22	2.26	0.66
1:B:197:LYS:HE2	1:B:198:TRP:NE1	2.12	0.66
1:B:86:LEU:HG	1:B:87:LEU:N	2.11	0.66
1:B:95:ASP:OD1	1:B:102:MET:HE3	1.96	0.65
1:A:111:ASP:OD2	1:A:113:GLU:HB3	1.96	0.65
1:B:185:GLN:O	1:B:188:LEU:CD1	2.42	0.65
1:B:236:GLY:O	1:B:240:GLU:HG2	1.95	0.65
1:A:197:LYS:HE2	1:A:198:TRP:NE1	2.12	0.65
1:B:71:ASP:HA	1:B:74:LYS:CE	2.20	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:GLN:HG3	1:B:184:GLY:N	2.09	0.65
1:B:35:THR:HB	1:B:38:ALA:HB3	1.78	0.65
1:B:195:GLN:HG2	1:B:198:TRP:CZ3	2.32	0.64
1:A:86:LEU:HG	1:A:87:LEU:N	2.13	0.64
1:B:56:VAL:HA	1:B:87:LEU:O	1.98	0.64
1:A:253:LYS:HD2	1:A:253:LYS:N	2.10	0.64
1:A:183:GLN:CG	1:A:184:GLY:H	1.97	0.64
1:B:230:LYS:HE2	1:B:230:LYS:HA	1.79	0.63
1:B:80:SER:O	1:B:82:SER:N	2.30	0.63
1:A:160:VAL:HG13	1:A:161:TYR:CD1	2.32	0.63
1:B:109:LYS:O	1:B:110:LEU:CD2	2.46	0.63
1:B:111:ASP:HB2	1:B:113:GLU:HG2	1.79	0.62
1:B:221:MET:HE3	1:B:249:THR:HG22	1.80	0.62
1:B:59:LYS:HD3	1:B:61:ASP:CB	2.29	0.62
1:A:50:ILE:CB	1:A:51:PRO:HD3	2.28	0.62
1:A:50:ILE:HB	1:A:51:PRO:CD	2.27	0.61
1:A:241:LEU:O	1:A:242:GLN:C	2.38	0.61
1:B:109:LYS:CD	1:B:109:LYS:H	2.13	0.61
1:B:206:MET:HA	1:B:209:ILE:HD12	1.82	0.61
1:B:50:ILE:HB	1:B:51:PRO:CD	2.28	0.61
1:B:109:LYS:O	1:B:110:LEU:HD22	1.99	0.61
1:B:175:GLU:CB	1:B:178:GLN:NE2	2.62	0.60
1:B:188:LEU:HD12	1:B:189:SER:N	2.16	0.60
1:B:230:LYS:HG3	1:B:232:LYS:HG3	1.83	0.60
1:B:86:LEU:HD12	1:B:87:LEU:H	1.66	0.60
1:A:200:GLU:HA	1:A:203:LEU:HD12	1.83	0.60
1:A:175:GLU:CB	1:A:178:GLN:NE2	2.64	0.60
1:A:77:ALA:O	1:A:80:SER:OG	2.20	0.60
1:B:200:GLU:HA	1:B:203:LEU:HD12	1.84	0.60
1:A:241:LEU:O	1:A:243:LYS:N	2.35	0.59
1:A:86:LEU:HD12	1:A:87:LEU:H	1.66	0.59
1:A:67:GLY:O	1:A:70:GLN:N	2.36	0.59
1:B:241:LEU:O	1:B:242:GLN:C	2.39	0.59
1:B:143:ARG:CG	1:B:143:ARG:NH1	2.56	0.59
1:B:230:LYS:HA	1:B:230:LYS:CE	2.32	0.58
1:A:56:VAL:HA	1:A:87:LEU:O	2.02	0.58
1:A:96:TYR:HE2	1:B:112:LYS:O	1.86	0.58
1:A:80:SER:O	1:A:82:SER:N	2.37	0.58
1:B:195:GLN:HG2	1:B:198:TRP:HZ3	1.66	0.58
1:A:219:SER:OG	1:A:220:GLU:N	2.35	0.58
1:A:39:LEU:HD22	1:A:41:LEU:HD21	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:ARG:CB	1:B:143:ARG:HH11	2.16	0.58
1:B:144:TRP:O	1:B:148:GLN:HB2	2.03	0.58
1:A:213:GLY:C	1:A:215:ASP:H	2.07	0.58
1:A:175:GLU:CA	1:A:178:GLN:NE2	2.67	0.58
1:B:164:LEU:HD21	1:B:182:LYS:HD2	1.86	0.58
1:B:77:ALA:O	1:B:80:SER:OG	2.22	0.58
1:A:237:LYS:HA	1:A:240:GLU:OE1	2.04	0.57
1:B:80:SER:C	1:B:82:SER:H	2.06	0.57
1:B:175:GLU:HA	1:B:178:GLN:HE21	1.70	0.57
1:B:175:GLU:CA	1:B:178:GLN:NE2	2.67	0.57
1:A:206:MET:HA	1:A:209:ILE:HD12	1.85	0.57
1:A:201:GLN:O	1:A:205:ILE:HG13	2.05	0.57
1:A:199:ALA:O	1:A:202:TYR:HB2	2.04	0.57
1:A:209:ILE:HG23	1:A:214:GLU:HA	1.85	0.57
1:A:144:TRP:O	1:A:148:GLN:HB2	2.05	0.56
1:A:200:GLU:HB3	1:A:204:LYS:HD3	1.86	0.56
1:B:241:LEU:O	1:B:243:LYS:N	2.39	0.56
1:A:45:THR:O	1:A:49:VAL:HG23	2.05	0.56
1:A:108:TYR:O	1:A:110:LEU:CG	2.54	0.56
1:A:109:LYS:O	1:A:110:LEU:HD23	2.05	0.56
1:A:209:ILE:CG2	1:A:214:GLU:HA	2.36	0.56
1:A:67:GLY:O	1:A:68:GLU:C	2.44	0.56
1:B:219:SER:OG	1:B:220:GLU:N	2.39	0.56
1:B:53:SER:HB2	1:B:56:VAL:CG2	2.35	0.56
1:B:182:LYS:HB3	1:B:186:ASP:OD2	2.06	0.56
1:B:216:PHE:HB3	1:B:217:PRO:HD3	1.88	0.56
1:B:185:GLN:O	1:B:188:LEU:CG	2.54	0.56
1:B:200:GLU:HB3	1:B:204:LYS:HD3	1.88	0.56
1:A:170:ARG:HB3	1:A:170:ARG:HH11	1.71	0.55
1:A:226:ARG:HH11	1:A:226:ARG:HG2	1.71	0.55
1:B:117:VAL:CG1	1:B:119:TYR:HE1	2.18	0.55
1:B:116:PRO:HB2	1:B:118:PHE:HE2	1.72	0.55
1:A:148:GLN:HG3	1:A:148:GLN:O	2.07	0.55
1:A:102:MET:O	1:A:104:LEU:N	2.40	0.55
1:A:117:VAL:CG1	1:A:119:TYR:HE1	2.20	0.55
1:B:35:THR:HB	1:B:38:ALA:CB	2.37	0.55
1:B:117:VAL:HG12	1:B:119:TYR:CE1	2.39	0.55
1:B:201:GLN:O	1:B:205:ILE:HG13	2.07	0.55
1:A:221:MET:SD	1:A:252:GLN:OE1	2.65	0.54
1:B:70:GLN:O	1:B:72:GLU:N	2.40	0.54
1:A:95:ASP:CB	1:A:112:LYS:HE2	2.34	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:PHE:CE1	1:A:206:MET:HE3	2.43	0.54
1:B:188:LEU:HD12	1:B:189:SER:HB2	1.88	0.54
1:A:114:SER:O	1:A:115:TYR:C	2.46	0.54
1:A:85:ASP:N	1:A:85:ASP:OD1	2.36	0.54
1:B:98:ASP:O	1:B:99:LYS:C	2.46	0.54
1:A:197:LYS:HE2	1:A:198:TRP:HE1	1.72	0.54
1:A:80:SER:C	1:A:82:SER:H	2.11	0.54
1:B:230:LYS:CA	1:B:230:LYS:CE	2.86	0.54
1:A:191:VAL:CG1	1:A:195:GLN:OE1	2.56	0.54
1:A:217:PRO:HB2	1:A:253:LYS:HZ2	1.71	0.54
1:B:175:GLU:HA	1:B:178:GLN:HB3	1.90	0.54
1:B:179:ALA:O	1:B:181:LEU:N	2.41	0.54
1:B:197:LYS:HE2	1:B:198:TRP:HE1	1.72	0.54
1:B:192:LYS:HB3	1:B:195:GLN:H	1.73	0.54
1:A:179:ALA:O	1:A:181:LEU:N	2.42	0.53
1:B:175:GLU:HA	1:B:178:GLN:NE2	2.23	0.53
1:B:221:MET:CG	1:B:248:LEU:HB3	2.38	0.53
1:B:39:LEU:HD22	1:B:41:LEU:HD21	1.90	0.53
1:B:68:GLU:HG2	1:B:69:LYS:H	1.71	0.53
1:A:146:LYS:HG2	1:A:152:LEU:HD12	1.91	0.53
1:A:39:LEU:H	1:A:89:ALA:HA	1.73	0.53
1:A:112:LYS:HB2	1:A:112:LYS:NZ	2.23	0.53
1:B:116:PRO:HB2	1:B:118:PHE:CE2	2.44	0.53
1:B:69:LYS:HB3	1:B:136:VAL:HG21	1.91	0.53
1:A:116:PRO:HB2	1:A:118:PHE:CE2	2.43	0.53
1:A:116:PRO:HB2	1:A:118:PHE:HE2	1.74	0.53
1:B:102:MET:O	1:B:104:LEU:N	2.41	0.53
1:B:180:LEU:HD13	1:B:206:MET:HE2	1.91	0.53
1:A:100:LEU:C	1:A:102:MET:N	2.62	0.53
1:A:175:GLU:HA	1:A:178:GLN:HE21	1.73	0.52
1:A:221:MET:CG	1:A:248:LEU:HB3	2.39	0.52
1:B:199:ALA:O	1:B:202:TYR:HB2	2.09	0.52
1:B:253:LYS:C	1:B:255:GLY:N	2.63	0.52
1:B:229:GLU:OE1	1:B:229:GLU:HA	2.10	0.52
1:A:112:LYS:HG3	1:A:115:TYR:CE2	2.45	0.52
1:B:180:LEU:O	1:B:183:GLN:HG2	2.09	0.52
1:A:175:GLU:HA	1:A:178:GLN:NE2	2.24	0.52
1:B:80:SER:C	1:B:82:SER:N	2.62	0.52
1:A:70:GLN:O	1:A:72:GLU:N	2.44	0.51
1:B:175:GLU:C	1:B:178:GLN:HB3	2.30	0.51
1:B:41:LEU:HD22	1:B:49:VAL:HG21	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:226:ARG:HH11	1:B:226:ARG:HG2	1.75	0.51
1:B:148:GLN:O	1:B:148:GLN:HG3	2.10	0.51
1:A:125:ASP:OD2	1:A:128:ASN:HB2	2.09	0.51
1:A:132:TYR:HD2	1:A:144:TRP:CD2	2.28	0.51
1:B:70:GLN:O	1:B:73:PHE:N	2.43	0.51
1:A:213:GLY:C	1:A:215:ASP:N	2.63	0.51
1:A:33:VAL:CG1	1:A:34:HIS:N	2.56	0.51
1:B:45:THR:O	1:B:49:VAL:HG23	2.11	0.51
1:B:114:SER:O	1:B:115:TYR:C	2.49	0.51
1:A:175:GLU:HA	1:A:178:GLN:HB3	1.93	0.51
1:A:145:LEU:O	1:A:146:LYS:C	2.49	0.51
1:B:158:LEU:O	1:B:161:TYR:N	2.44	0.51
1:B:86:LEU:CG	1:B:87:LEU:N	2.74	0.51
1:A:158:LEU:O	1:A:161:TYR:N	2.44	0.50
1:B:100:LEU:C	1:B:102:MET:N	2.63	0.50
1:B:249:THR:HA	1:B:252:GLN:HG3	1.92	0.50
1:B:39:LEU:H	1:B:89:ALA:HA	1.77	0.50
1:A:71:ASP:CA	1:A:74:LYS:HD2	2.27	0.50
1:A:100:LEU:O	1:A:100:LEU:HD23	2.11	0.50
1:B:95:ASP:HA	1:B:102:MET:HE3	1.94	0.50
1:A:35:THR:HB	1:A:38:ALA:HB3	1.94	0.49
1:B:109:LYS:O	1:B:110:LEU:HD23	2.12	0.49
1:A:190:SER:OG	1:A:196:LYS:HG3	2.12	0.49
1:B:54:LYS:NZ	1:B:85:ASP:HB3	2.26	0.49
1:A:117:VAL:HG12	1:A:119:TYR:CE1	2.41	0.49
1:A:86:LEU:CG	1:A:87:LEU:N	2.74	0.49
1:B:132:TYR:HD2	1:B:144:TRP:CD2	2.30	0.49
1:A:96:TYR:CE2	1:B:112:LYS:O	2.64	0.49
1:B:127:GLU:C	1:B:129:PRO:HD3	2.32	0.49
1:B:145:LEU:O	1:B:146:LYS:C	2.51	0.49
1:B:216:PHE:N	1:B:217:PRO:CD	2.75	0.49
1:B:170:ARG:HE	1:B:254:LYS:HG3	1.78	0.49
1:B:221:MET:HG2	1:B:248:LEU:HB3	1.95	0.49
1:B:255:GLY:O	1:B:256:ALA:HB3	2.13	0.49
1:A:192:LYS:HB3	1:A:195:GLN:H	1.78	0.49
1:A:98:ASP:O	1:A:99:LYS:C	2.50	0.49
1:B:125:ASP:OD2	1:B:128:ASN:HB2	2.12	0.49
1:A:112:LYS:HG3	1:A:115:TYR:CD2	2.48	0.49
1:B:59:LYS:HG3	1:B:118:PHE:CZ	2.48	0.49
1:B:98:ASP:O	1:B:99:LYS:O	2.31	0.49
1:A:147:GLY:C	1:A:149:GLY:H	2.15	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:LEU:O	1:B:249:THR:HG23	2.13	0.48
1:B:147:GLY:C	1:B:149:GLY:H	2.16	0.48
1:A:245:LEU:O	1:A:249:THR:HG23	2.13	0.48
1:A:70:GLN:O	1:A:73:PHE:N	2.47	0.48
1:B:185:GLN:O	1:B:188:LEU:HD21	2.13	0.48
1:B:71:ASP:CA	1:B:74:LYS:HE2	2.28	0.48
1:B:110:LEU:HD21	1:B:119:TYR:OH	2.13	0.48
1:B:66:TYR:CG	1:B:67:GLY:N	2.81	0.48
1:A:127:GLU:C	1:A:129:PRO:HD3	2.34	0.48
1:A:175:GLU:C	1:A:178:GLN:HB3	2.34	0.48
1:A:187:ASN:C	1:A:189:SER:N	2.65	0.48
1:A:91:VAL:HG13	1:A:101:ASN:OD1	2.13	0.48
1:A:222:THR:HG23	1:A:226:ARG:HH12	1.78	0.48
1:B:177:ARG:O	1:B:178:GLN:C	2.52	0.48
1:A:132:TYR:CD2	1:A:144:TRP:CE3	3.02	0.48
1:A:175:GLU:C	1:A:178:GLN:HE21	2.17	0.48
1:B:222:THR:HG23	1:B:226:ARG:HH12	1.79	0.48
1:B:61:ASP:CG	1:B:62:THR:N	2.66	0.48
1:B:132:TYR:HD2	1:B:144:TRP:CE3	2.31	0.47
1:A:221:MET:HG2	1:A:248:LEU:HB3	1.96	0.47
1:A:132:TYR:HD2	1:A:144:TRP:CE3	2.31	0.47
1:B:100:LEU:O	1:B:102:MET:N	2.47	0.47
1:A:168:PHE:CE1	1:A:176:ALA:HB1	2.50	0.47
1:B:118:PHE:C	1:B:119:TYR:HD1	2.18	0.47
1:B:132:TYR:CD2	1:B:144:TRP:CE3	3.03	0.47
1:B:216:PHE:N	1:B:217:PRO:HD2	2.29	0.47
1:B:234:SER:HG	1:B:237:LYS:HD3	1.78	0.47
1:A:80:SER:C	1:A:82:SER:N	2.66	0.47
1:A:172:SER:HB3	1:A:175:GLU:OE1	2.15	0.47
1:A:187:ASN:C	1:A:189:SER:H	2.17	0.47
1:B:221:MET:HE1	1:B:249:THR:HG22	1.97	0.47
1:A:100:LEU:O	1:A:102:MET:N	2.48	0.46
1:A:135:ALA:O	1:A:137:LYS:N	2.44	0.46
1:A:170:ARG:HB3	1:A:170:ARG:NH1	2.30	0.46
1:A:50:ILE:N	1:A:51:PRO:CD	2.78	0.46
1:A:221:MET:HE1	1:A:249:THR:HG22	1.96	0.46
1:B:202:TYR:O	1:B:205:ILE:HB	2.16	0.46
1:B:175:GLU:C	1:B:178:GLN:HE21	2.18	0.46
1:B:254:LYS:O	1:B:255:GLY:C	2.53	0.46
1:B:175:GLU:OE1	1:B:178:GLN:CD	2.53	0.46
1:B:168:PHE:CE1	1:B:206:MET:HE3	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:LYS:HE2	1:A:139:GLY:HA3	1.97	0.46
1:A:216:PHE:N	1:A:217:PRO:CD	2.79	0.46
1:A:61:ASP:CG	1:A:62:THR:N	2.69	0.46
1:A:62:THR:O	1:A:65:PRO:HD3	2.15	0.46
1:B:54:LYS:HZ1	1:B:85:ASP:HB3	1.81	0.46
1:A:202:TYR:O	1:A:205:ILE:HB	2.15	0.46
1:B:46:PHE:CD2	1:B:46:PHE:C	2.89	0.46
1:A:180:LEU:HD13	1:A:206:MET:HE2	1.96	0.46
1:B:79:ASN:HD22	1:B:142:GLN:HE22	1.64	0.46
1:B:206:MET:HE3	1:B:206:MET:HB3	1.79	0.46
1:A:46:PHE:CD2	1:A:46:PHE:C	2.89	0.46
1:A:68:GLU:O	1:A:69:LYS:C	2.53	0.46
1:A:132:TYR:HB2	1:A:144:TRP:CE3	2.50	0.45
1:A:35:THR:O	1:A:38:ALA:N	2.41	0.45
1:B:255:GLY:O	1:B:256:ALA:CB	2.65	0.45
1:A:146:LYS:CG	1:A:152:LEU:HD12	2.46	0.45
1:A:79:ASN:HD22	1:A:142:GLN:HE22	1.64	0.45
1:A:106:GLU:O	1:A:107:LYS:C	2.54	0.45
1:A:191:VAL:HG13	1:A:195:GLN:OE1	2.17	0.45
1:A:42:ASP:H	1:A:45:THR:HB	1.81	0.45
1:B:108:TYR:O	1:B:109:LYS:C	2.53	0.45
1:A:216:PHE:HB3	1:A:217:PRO:HD3	1.98	0.45
1:A:180:LEU:HB2	1:A:206:MET:CE	2.47	0.45
1:B:41:LEU:CD2	1:B:49:VAL:HG21	2.45	0.45
1:B:253:LYS:O	1:B:255:GLY:N	2.50	0.45
1:B:180:LEU:HB2	1:B:206:MET:CE	2.46	0.45
1:A:177:ARG:O	1:A:181:LEU:HG	2.17	0.45
1:A:33:VAL:CG1	1:A:34:HIS:H	2.03	0.45
1:A:192:LYS:HB2	1:A:195:GLN:CG	2.25	0.45
1:B:222:THR:O	1:B:225:ALA:HB3	2.16	0.44
1:B:254:LYS:HD3	1:B:254:LYS:HA	1.71	0.44
1:A:41:LEU:HD22	1:A:49:VAL:HG21	1.98	0.44
1:B:156:GLY:HA2	1:B:198:TRP:CE3	2.52	0.44
1:B:42:ASP:O	1:B:44:VAL:N	2.50	0.44
1:A:93:ILE:HD11	1:A:102:MET:SD	2.58	0.44
1:A:169:ILE:HA	1:A:169:ILE:HD13	1.56	0.44
1:A:177:ARG:O	1:A:178:GLN:C	2.56	0.44
1:B:143:ARG:HB2	1:B:143:ARG:HH11	1.82	0.44
1:B:241:LEU:C	1:B:243:LYS:N	2.71	0.44
1:A:144:TRP:O	1:A:148:GLN:CB	2.66	0.44
1:B:144:TRP:O	1:B:148:GLN:CB	2.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:LYS:O	1:B:239:GLU:C	2.56	0.44
1:B:70:GLN:C	1:B:72:GLU:N	2.69	0.44
1:A:137:LYS:O	1:A:140:ALA:HB3	2.18	0.44
1:B:155:PRO:HG2	1:B:246:ASN:ND2	2.33	0.44
1:A:42:ASP:HA	1:A:101:ASN:HD21	1.83	0.44
1:A:118:PHE:N	1:A:118:PHE:CD2	2.85	0.44
1:B:95:ASP:CA	1:B:102:MET:HE3	2.48	0.44
1:A:119:TYR:HB3	1:A:129:PRO:HB2	2.00	0.44
1:A:136:VAL:O	1:A:136:VAL:HG23	2.18	0.43
1:B:39:LEU:C	1:B:39:LEU:HD23	2.39	0.43
1:A:100:LEU:C	1:A:100:LEU:HD23	2.38	0.43
1:A:241:LEU:C	1:A:243:LYS:N	2.70	0.43
1:B:62:THR:O	1:B:65:PRO:HD3	2.18	0.43
1:B:119:TYR:HB3	1:B:129:PRO:HB2	2.01	0.43
1:A:238:LYS:O	1:A:239:GLU:C	2.56	0.43
1:A:71:ASP:O	1:A:74:LYS:HB2	2.18	0.43
1:B:66:TYR:CD1	1:B:67:GLY:N	2.87	0.43
1:A:63:GLN:HG3	1:A:64:TYR:CD2	2.53	0.43
1:B:118:PHE:N	1:B:118:PHE:CD2	2.86	0.43
1:B:183:GLN:C	1:B:187:ASN:HD22	2.22	0.43
1:B:42:ASP:H	1:B:45:THR:HB	1.82	0.43
1:A:59:LYS:NZ	1:A:70:GLN:OE1	2.51	0.43
1:B:146:LYS:HB2	1:B:146:LYS:HE3	1.74	0.43
1:A:226:ARG:HG2	1:A:226:ARG:NH1	2.33	0.43
1:A:100:LEU:HD23	1:A:101:ASN:HB2	2.00	0.43
1:A:156:GLY:HA2	1:A:198:TRP:CE3	2.54	0.43
1:A:197:LYS:HE2	1:A:198:TRP:CD1	2.54	0.43
1:A:61:ASP:OD2	1:A:65:PRO:HB3	2.19	0.43
1:B:53:SER:HB2	1:B:56:VAL:HG23	2.01	0.43
1:A:69:LYS:HG2	1:A:69:LYS:H	1.48	0.42
1:A:203:LEU:O	1:A:206:MET:HB2	2.19	0.42
1:B:50:ILE:N	1:B:51:PRO:CD	2.81	0.42
1:B:136:VAL:HG23	1:B:136:VAL:O	2.19	0.42
1:A:170:ARG:CB	1:A:170:ARG:HH11	2.31	0.42
1:B:63:GLN:HG3	1:B:64:TYR:CD2	2.54	0.42
1:B:86:LEU:HD12	1:B:87:LEU:N	2.33	0.42
1:A:68:GLU:O	1:A:71:ASP:N	2.52	0.42
1:A:70:GLN:C	1:A:72:GLU:N	2.72	0.42
1:B:206:MET:O	1:B:209:ILE:N	2.52	0.42
1:B:175:GLU:CA	1:B:178:GLN:HB3	2.49	0.42
1:B:177:ARG:O	1:B:180:LEU:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:GLY:HA3	1:B:77:ALA:HB1	2.01	0.42
1:B:91:VAL:HG13	1:B:101:ASN:OD1	2.20	0.42
1:B:250:ALA:O	1:B:253:LYS:HB2	2.20	0.42
1:A:204:LYS:HD2	1:A:204:LYS:N	2.35	0.42
1:B:108:TYR:OH	1:B:126:PHE:O	2.25	0.42
1:A:132:TYR:CZ	1:A:134:GLY:CA	3.01	0.42
1:A:198:TRP:O	1:A:201:GLN:HB3	2.20	0.42
1:A:37:GLY:HA3	1:A:77:ALA:HB1	2.02	0.42
1:B:137:LYS:O	1:B:140:ALA:HB3	2.19	0.42
1:B:197:LYS:HE2	1:B:198:TRP:CD1	2.55	0.42
1:A:102:MET:C	1:A:104:LEU:N	2.71	0.41
1:B:142:GLN:O	1:B:143:ARG:C	2.59	0.41
1:B:106:GLU:O	1:B:107:LYS:C	2.57	0.41
1:B:170:ARG:NE	1:B:254:LYS:HG3	2.35	0.41
1:A:101:ASN:O	1:A:102:MET:C	2.57	0.41
1:A:86:LEU:HD12	1:A:87:LEU:N	2.34	0.41
1:A:98:ASP:O	1:A:99:LYS:O	2.38	0.41
1:B:132:TYR:HB2	1:B:144:TRP:CE3	2.51	0.41
1:B:204:LYS:HD2	1:B:204:LYS:N	2.35	0.41
1:A:187:ASN:O	1:A:189:SER:N	2.53	0.41
1:A:39:LEU:C	1:A:39:LEU:HD23	2.41	0.41
1:A:102:MET:O	1:A:103:GLU:C	2.57	0.41
1:B:61:ASP:OD2	1:B:65:PRO:HB3	2.21	0.41
1:B:70:GLN:C	1:B:72:GLU:H	2.23	0.41
1:B:86:LEU:CD1	1:B:87:LEU:H	2.32	0.41
1:A:118:PHE:C	1:A:119:TYR:HD1	2.23	0.41
1:A:225:ALA:O	1:A:229:GLU:HG2	2.21	0.41
1:B:50:ILE:CB	1:B:51:PRO:HD3	2.29	0.41
1:B:59:LYS:O	1:B:59:LYS:HD2	2.20	0.41
1:A:234:SER:H	1:A:237:LYS:HD2	1.86	0.41
1:A:125:ASP:CG	1:A:128:ASN:HB2	2.41	0.41
1:B:169:ILE:HG23	1:B:169:ILE:HD12	1.86	0.41
1:A:206:MET:O	1:A:209:ILE:N	2.54	0.41
1:A:236:GLY:O	1:A:240:GLU:CG	2.63	0.41
1:B:198:TRP:O	1:B:201:GLN:HB3	2.20	0.41
1:B:203:LEU:O	1:B:204:LYS:C	2.59	0.41
1:B:226:ARG:NH1	1:B:226:ARG:HG2	2.35	0.41
1:B:102:MET:O	1:B:103:GLU:C	2.59	0.41
1:B:42:ASP:C	1:B:44:VAL:N	2.73	0.41
1:A:41:LEU:CD2	1:A:49:VAL:HG21	2.51	0.41
1:A:61:ASP:OD1	1:A:62:THR:O	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:ASN:O	1:B:102:MET:C	2.59	0.41
1:B:95:ASP:OD1	1:B:102:MET:CE	2.68	0.41
1:A:101:ASN:HD22	1:A:101:ASN:HA	1.62	0.40
1:A:96:TYR:CD2	1:B:113:GLU:HA	2.56	0.40
1:B:44:VAL:O	1:B:46:PHE:N	2.54	0.40
1:B:179:ALA:O	1:B:180:LEU:C	2.59	0.40
1:B:177:ARG:O	1:B:181:LEU:HG	2.21	0.40
1:B:190:SER:O	1:B:190:SER:OG	2.36	0.40
1:B:35:THR:O	1:B:36:LYS:C	2.58	0.40
1:B:102:MET:C	1:B:104:LEU:N	2.75	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	222/240 (92%)	155 (70%)	48 (22%)	19 (9%)	1	2
1	B	222/240 (92%)	157 (71%)	44 (20%)	21 (10%)	1	1
All	All	444/480 (92%)	312 (70%)	92 (21%)	40 (9%)	1	2

All (40) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	36	LYS
1	A	81	ALA
1	A	183	GLN
1	A	191	VAL
1	B	81	ALA
1	B	183	GLN
1	B	191	VAL
1	B	253	LYS

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Mol	Chain	Res	Type
1	A	71	ASP
1	A	99	LYS
1	A	103	GLU
1	A	109	LYS
1	A	136	VAL
1	A	180	LEU
1	A	188	LEU
1	A	190	SER
1	A	242	GLN
1	B	71	ASP
1	B	99	LYS
1	B	103	GLU
1	B	109	LYS
1	B	136	VAL
1	B	180	LEU
1	B	190	SER
1	B	202	TYR
1	A	46	PHE
1	A	148	GLN
1	A	202	TYR
1	A	253	LYS
1	B	46	PHE
1	B	148	GLN
1	B	215	ASP
1	B	242	GLN
1	B	254	LYS
1	B	255	GLY
1	A	217	PRO
1	B	101	ASN
1	B	217	PRO
1	A	101	ASN
1	B	36	LYS

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	186/205 (91%)	165 (89%)	21 (11%)	7	21
1	B	186/205 (91%)	160 (86%)	26 (14%)	4	12
All	All	372/410 (91%)	325 (87%)	47 (13%)	5	15

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

Mol	Chain	Res	Type
1	A	35	THR
1	A	53	SER
1	A	61	ASP
1	A	68	GLU
1	A	69	LYS
1	A	80	SER
1	A	82	SER
1	A	85	ASP
1	A	88	VAL
1	A	94	SER
1	A	107	LYS
1	A	112	LYS
1	A	122	ARG
1	A	126	PHE
1	A	169	ILE
1	A	170	ARG
1	A	178	GLN
1	A	189	SER
1	A	204	LYS
1	A	214	GLU
1	A	249	THR
1	B	35	THR
1	B	48	LYS
1	B	54	LYS
1	B	59	LYS
1	B	61	ASP
1	B	66	TYR
1	B	68	GLU
1	B	82	SER
1	B	85	ASP
1	B	88	VAL
1	B	94	SER
1	B	109	LYS
1	B	110	LEU
1	B	112	LYS

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Mol	Chain	Res	Type
1	B	113	GLU
1	B	143	ARG
1	B	169	ILE
1	B	178	GLN
1	B	188	LEU
1	B	189	SER
1	B	204	LYS
1	B	215	ASP
1	B	230	LYS
1	B	237	LYS
1	B	249	THR
1	B	253	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	ASN
1	A	142	GLN
1	A	178	GLN
1	A	187	ASN
1	A	212	GLN
1	B	101	ASN
1	B	142	GLN
1	B	178	GLN
1	B	187	ASN
1	B	212	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 carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	224/240 (93%)	-0.08	5 (2%) 62 59	48, 54, 59, 76	0
1	B	224/240 (93%)	-0.07	8 (3%) 43 37	48, 54, 59, 70	0
All	All	448/480 (93%)	-0.08	13 (2%) 52 46	48, 54, 59, 76	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	191	VAL	5.1
1	B	191	VAL	5.0
1	A	256	ALA	3.5
1	B	175	GLU	3.2
1	A	66	TYR	2.9
1	B	210	LEU	2.7
1	B	256	ALA	2.7
1	B	66	TYR	2.3
1	A	193	GLU	2.2
1	B	207	GLY	2.2
1	A	214	GLU	2.2
1	B	172	SER	2.1
1	B	253	LYS	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 carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.