



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2021 – 06:52 PM EDT

PDB ID : 1P4L
Title : Crystal structure of NK receptor Ly49C mutant with its MHC class I ligand H-2Kb
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Deposited on : 2003-04-23
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

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
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)	:	2.23.2

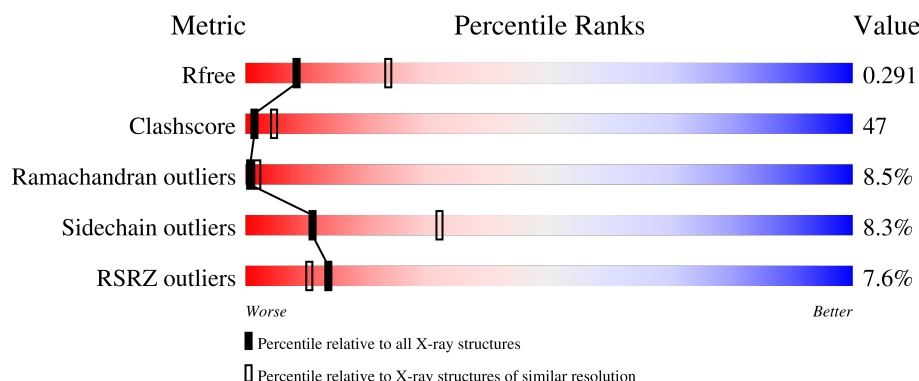
1 Overall quality at a glance

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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (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 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	274	
2	B	99	
3	P	8	
4	D	122	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4134 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 MHC CLASS I H-2KB HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2228	1406	392	421	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			818	523	138	150	7			

- Molecule 3 is a protein called Ovalbumin peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	8	Total	C	N	O	0	0	0
			68	45	10	13			

- Molecule 4 is a protein called LY49-C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	122	Total	C	N	O	S	0	0	0
			1020	664	170	175	11			

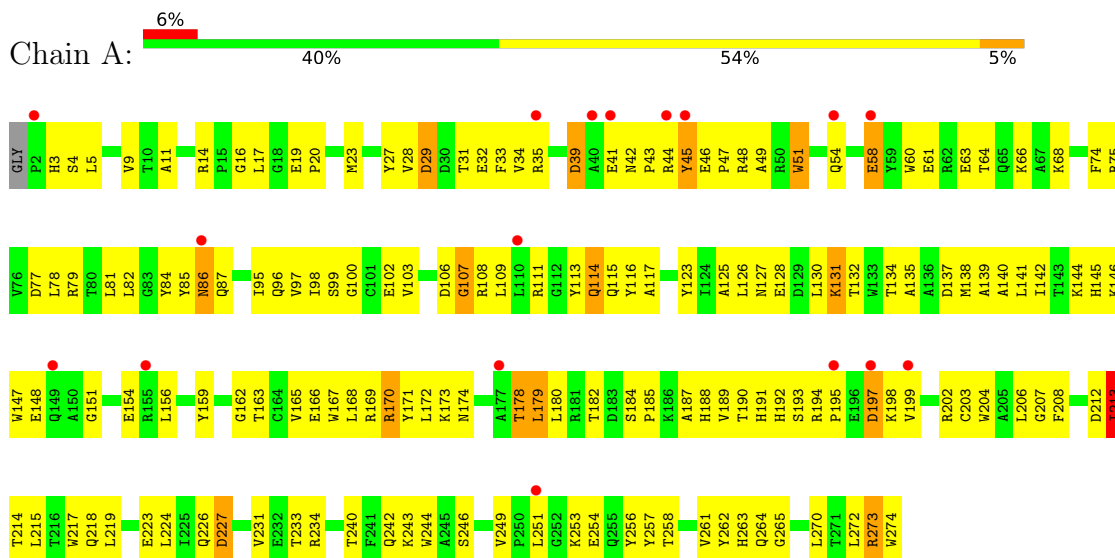
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	171	GLY	SER	engineered mutation	UNP Q64329
D	193	GLY	GLU	engineered mutation	UNP Q64329
D	223	LYS	ARG	engineered mutation	UNP Q64329

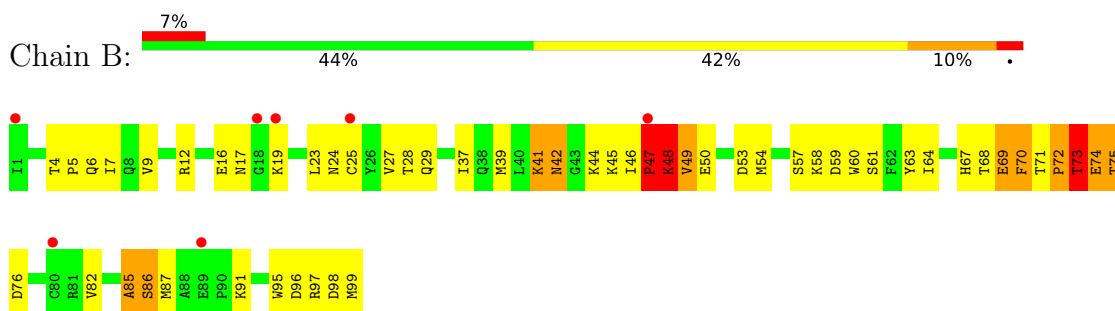
3 Residue-property plots

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.

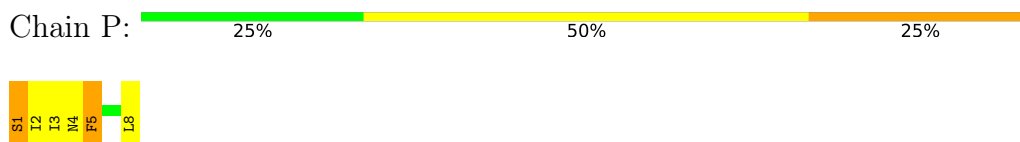
• Molecule 1: MHC CLASS I H-2KB HEAVY CHAIN



• Molecule 2: Beta-2-microglobulin

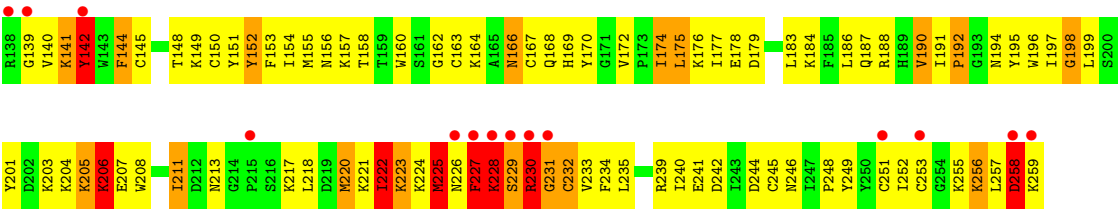


• Molecule 3: Ovalbumin peptide



• Molecule 4: LY49-C





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	152.01Å 152.01Å 64.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 2.90 19.96 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.6 (19.96-2.90) 95.5 (19.96-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.10 (at 2.88Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.269 , 0.295 0.267 , 0.291	Depositor DCC
R_{free} test set	820 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	69.0	Xtriage
Anisotropy	0.306	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 75.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4134	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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	0.37	0/2289	0.64	1/3107 (0.0%)
2	B	0.48	0/844	0.71	1/1144 (0.1%)
3	P	1.08	0/68	1.07	1/88 (1.1%)
4	D	0.74	2/1049 (0.2%)	0.94	4/1406 (0.3%)
All	All	0.52	2/4250 (0.0%)	0.74	7/5745 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	225	MET	SD-CE	5.66	2.09	1.77
4	D	142	TYR	CE1-CZ	5.24	1.45	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	231	GLY	N-CA-C	8.26	133.74	113.10
4	D	232	CYS	N-CA-C	7.38	130.92	111.00
4	D	225	MET	N-CA-C	6.64	128.94	111.00
4	D	230	ARG	N-CA-C	5.93	127.00	111.00
2	B	74	GLU	N-CA-C	-5.30	96.70	111.00
3	P	5	PHE	N-CA-C	5.13	124.84	111.00
1	A	213	ILE	N-CA-C	-5.12	97.19	111.00

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	2228	0	2118	175	0
2	B	818	0	797	84	0
3	P	68	0	74	15	0
4	D	1020	0	1013	125	0
All	All	4134	0	4002	381	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 47.

All (381) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:225:MET:SD	4:D:225:MET:CE	2.09	1.38
2:B:72:PRO:O	2:B:73:THR:HG22	1.41	1.20
2:B:72:PRO:O	2:B:73:THR:CG2	2.05	1.05
4:D:154:ILE:HB	4:D:251:CYS:HB2	1.49	0.92
2:B:46:ILE:HG23	2:B:47:PRO:HD2	1.49	0.92
4:D:256:LYS:H	4:D:256:LYS:HD2	1.35	0.90
1:A:81:LEU:HD23	1:A:84:TYR:HD2	1.37	0.89
2:B:73:THR:HG23	2:B:75:THR:HG23	1.54	0.89
2:B:41:LYS:N	2:B:41:LYS:HE3	1.88	0.89
2:B:27:VAL:HG21	2:B:37:ILE:HD12	1.54	0.88
2:B:41:LYS:HE3	2:B:41:LYS:H	1.39	0.88
2:B:41:LYS:HZ2	2:B:42:ASN:HB2	1.40	0.87
1:A:4:SER:HB3	1:A:102:GLU:HG2	1.54	0.87
2:B:96:ASP:HB3	2:B:99:MET:HB2	1.57	0.84
2:B:73:THR:OG1	2:B:74:GLU:N	2.05	0.84
4:D:174:ILE:HD12	4:D:198:GLY:N	1.92	0.84
1:A:131:LYS:HE2	1:A:131:LYS:HA	1.59	0.84
2:B:73:THR:CG2	2:B:75:THR:HG23	2.08	0.84
4:D:174:ILE:HG22	4:D:211:ILE:HG13	1.60	0.83
2:B:41:LYS:NZ	2:B:42:ASN:HB2	1.93	0.83
1:A:213:ILE:HD11	1:A:261:VAL:HG13	1.61	0.82
4:D:201:TYR:HB3	4:D:230:ARG:O	1.79	0.82
1:A:84:TYR:CE1	1:A:142:ILE:HG21	2.16	0.81
1:A:51:TRP:HB2	1:A:174:ASN:O	1.81	0.81
1:A:193:SER:O	1:A:194:ARG:HG3	1.81	0.81
4:D:230:ARG:NH1	4:D:245:CYS:HB2	1.97	0.80
3:P:3:ILE:HG12	3:P:4:ASN:N	1.95	0.80
1:A:131:LYS:HE3	1:A:154:GLU:HA	1.62	0.80
1:A:207:GLY:HA2	1:A:240:THR:HB	1.61	0.79
4:D:177:ILE:HD13	4:D:183:LEU:HD13	1.62	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ALA:HB1	1:A:140:ALA:HB3	1.63	0.78
2:B:46:ILE:O	2:B:48:LYS:N	2.17	0.78
4:D:225:MET:SD	4:D:225:MET:O	2.43	0.77
1:A:226:GLN:OE1	1:A:226:GLN:HA	1.83	0.77
1:A:114:GLN:HB2	1:A:156:LEU:HD21	1.67	0.76
4:D:148:THR:HG22	4:D:148:THR:O	1.84	0.76
4:D:194:ASN:ND2	4:D:234:PHE:CE1	2.53	0.76
1:A:206:LEU:HD23	1:A:242:GLN:HB2	1.68	0.75
2:B:23:LEU:HD13	2:B:70:PHE:CZ	2.22	0.75
1:A:4:SER:CB	1:A:102:GLU:HG2	2.15	0.75
1:A:81:LEU:HD23	1:A:84:TYR:CD2	2.21	0.74
4:D:167:CYS:HB3	4:D:172:VAL:O	1.87	0.74
1:A:231:VAL:O	1:A:243:LYS:HE3	1.87	0.74
2:B:6:GLN:HG3	2:B:29:GLN:OE1	1.88	0.74
2:B:73:THR:O	2:B:97:ARG:NH2	2.21	0.74
2:B:42:ASN:OD1	2:B:76:ASP:HB3	1.87	0.73
1:A:128:GLU:O	1:A:130:LEU:HG	1.89	0.72
2:B:17:ASN:HB3	2:B:73:THR:HB	1.71	0.72
4:D:175:LEU:HD12	4:D:176:LYS:H	1.53	0.72
4:D:225:MET:O	4:D:225:MET:CG	2.38	0.71
1:A:213:ILE:HD13	1:A:214:THR:N	2.05	0.71
1:A:64:THR:O	1:A:68:LYS:HG3	1.91	0.71
4:D:158:THR:HB	4:D:162:GLY:HA3	1.72	0.71
4:D:191:ILE:HG23	4:D:192:PRO:HD2	1.74	0.70
4:D:175:LEU:HD22	4:D:252:ILE:HG21	1.73	0.69
2:B:27:VAL:HG21	2:B:37:ILE:CD1	2.22	0.69
2:B:41:LYS:HE2	2:B:44:LYS:O	1.92	0.69
1:A:128:GLU:CD	1:A:128:GLU:H	1.95	0.69
4:D:187:GLN:HG2	4:D:235:LEU:HD11	1.74	0.69
4:D:225:MET:CE	4:D:230:ARG:H	2.07	0.68
1:A:244:TRP:HZ3	1:A:246:SER:HG	1.41	0.68
4:D:141:LYS:NZ	4:D:141:LYS:HB2	2.08	0.68
2:B:42:ASN:HD21	2:B:76:ASP:CG	1.97	0.68
4:D:149:LYS:HE3	4:D:256:LYS:HG3	1.74	0.68
2:B:39:MET:HG3	2:B:49:VAL:HG21	1.74	0.67
4:D:174:ILE:HD13	4:D:175:LEU:N	2.09	0.67
2:B:29:GLN:HA	2:B:61:SER:HB2	1.76	0.67
4:D:225:MET:HE1	4:D:230:ARG:H	1.60	0.67
4:D:140:VAL:HG12	4:D:155:MET:HB2	1.76	0.67
1:A:68:LYS:NZ	1:A:68:LYS:HB3	2.10	0.67
1:A:144:LYS:O	1:A:148:GLU:HG3	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:41:LYS:HE2	2:B:44:LYS:HB2	1.77	0.66
2:B:16:GLU:O	2:B:19:LYS:HG2	1.96	0.66
4:D:174:ILE:HD12	4:D:198:GLY:H	1.60	0.66
1:A:9:VAL:HB	1:A:97:VAL:HB	1.77	0.66
4:D:208:TRP:HZ2	4:D:225:MET:HB3	1.60	0.66
1:A:199:VAL:HG23	1:A:249:VAL:HG22	1.78	0.65
4:D:164:LYS:O	4:D:168:GLN:HG2	1.97	0.65
1:A:273:ARG:NH1	1:A:273:ARG:HG2	2.11	0.65
1:A:35:ARG:CZ	2:B:53:ASP:HB3	2.27	0.65
1:A:84:TYR:HE1	1:A:142:ILE:HG21	1.60	0.65
1:A:137:ASP:OD1	1:A:139:ALA:HB3	1.97	0.64
1:A:85:TYR:O	1:A:86:ASN:HB3	1.95	0.64
1:A:199:VAL:HG22	1:A:249:VAL:O	1.97	0.64
4:D:256:LYS:H	4:D:256:LYS:CD	2.08	0.64
1:A:273:ARG:HG2	1:A:273:ARG:HH11	1.63	0.64
1:A:77:ASP:HB3	3:P:8:LEU:HD12	1.78	0.63
1:A:23:MET:HE2	1:A:35:ARG:HE	1.62	0.63
2:B:45:LYS:NZ	2:B:48:LYS:HG3	2.14	0.63
2:B:97:ARG:HG3	2:B:98:ASP:N	2.14	0.63
1:A:111:ARG:CZ	4:D:222:ILE:HG22	2.28	0.63
4:D:224:LYS:HB3	4:D:242:ASP:OD2	1.99	0.63
1:A:111:ARG:NH1	4:D:222:ILE:HG22	2.14	0.62
4:D:160:TRP:CZ3	4:D:164:LYS:HD2	2.33	0.62
1:A:79:ARG:NH1	1:A:82:LEU:HD12	2.15	0.62
1:A:47:PRO:HB3	1:A:60:TRP:CH2	2.35	0.61
2:B:25:CYS:HB2	2:B:39:MET:CE	2.30	0.61
4:D:225:MET:HE1	4:D:230:ARG:N	2.15	0.61
4:D:140:VAL:CG1	4:D:155:MET:HB2	2.30	0.61
3:P:3:ILE:CG1	3:P:4:ASN:N	2.62	0.61
1:A:63:GLU:OE1	3:P:1:SER:HA	2.01	0.60
1:A:197:ASP:OD2	1:A:251:LEU:HB3	2.01	0.60
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.35	0.60
4:D:157:LYS:HA	4:D:249:TYR:O	2.01	0.60
1:A:68:LYS:HB3	1:A:68:LYS:HZ2	1.66	0.60
4:D:178:GLU:O	4:D:217:LYS:HD2	2.01	0.60
4:D:187:GLN:CG	4:D:235:LEU:HD11	2.31	0.60
1:A:5:LEU:HB2	1:A:168:LEU:HD13	1.82	0.60
1:A:138:MET:HA	1:A:141:LEU:HD22	1.84	0.60
2:B:72:PRO:O	2:B:73:THR:HG23	2.00	0.60
4:D:225:MET:CE	4:D:225:MET:HB2	2.31	0.60
2:B:46:ILE:HG23	2:B:47:PRO:CD	2.29	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:GLY:O	1:A:17:LEU:HD23	2.01	0.60
1:A:135:ALA:HB1	1:A:140:ALA:CB	2.32	0.59
1:A:42:ASN:HD21	1:A:44:ARG:HG3	1.65	0.59
1:A:47:PRO:HB3	1:A:60:TRP:HH2	1.66	0.59
1:A:207:GLY:HA2	1:A:240:THR:CB	2.33	0.59
2:B:73:THR:HG23	2:B:75:THR:CG2	2.30	0.59
4:D:190:VAL:HG13	4:D:195:TYR:CE1	2.37	0.59
1:A:19:GLU:OE2	1:A:75:ARG:HD3	2.03	0.59
1:A:162:GLY:O	1:A:166:GLU:HG3	2.02	0.59
1:A:170:ARG:HG2	1:A:174:ASN:HD21	1.67	0.58
1:A:187:ALA:HA	1:A:204:TRP:O	2.04	0.58
4:D:148:THR:O	4:D:148:THR:CG2	2.49	0.58
4:D:172:VAL:HG13	4:D:255:LYS:HB2	1.86	0.58
1:A:178:THR:C	1:A:180:LEU:H	2.07	0.58
4:D:167:CYS:SG	4:D:174:ILE:HA	2.43	0.58
4:D:175:LEU:HD22	4:D:252:ILE:CG2	2.33	0.57
2:B:41:LYS:CE	2:B:44:LYS:HB2	2.34	0.57
4:D:153:PHE:HB3	4:D:155:MET:HE1	1.85	0.57
4:D:230:ARG:HD3	4:D:231:GLY:N	2.20	0.57
1:A:170:ARG:HG2	1:A:174:ASN:ND2	2.19	0.57
2:B:41:LYS:HE2	2:B:44:LYS:CA	2.35	0.57
4:D:208:TRP:CZ2	4:D:225:MET:HB3	2.40	0.57
1:A:218:GLN:HG2	1:A:223:GLU:HA	1.86	0.57
1:A:106:ASP:O	1:A:108:ARG:N	2.37	0.57
1:A:126:LEU:HD22	1:A:156:LEU:HD23	1.87	0.57
2:B:69:GLU:N	2:B:69:GLU:OE1	2.38	0.57
4:D:225:MET:SD	4:D:230:ARG:N	2.77	0.57
4:D:141:LYS:NZ	4:D:141:LYS:CB	2.68	0.56
4:D:196:TRP:HH2	4:D:244:ASP:O	1.87	0.56
1:A:35:ARG:NH2	2:B:53:ASP:HB3	2.21	0.56
1:A:44:ARG:O	1:A:46:GLU:N	2.39	0.56
1:A:99:SER:OG	1:A:114:GLN:HG3	2.06	0.56
4:D:149:LYS:HG2	4:D:256:LYS:HA	1.86	0.56
1:A:4:SER:HB3	1:A:102:GLU:CG	2.31	0.56
1:A:98:ILE:O	1:A:114:GLN:HA	2.05	0.56
1:A:106:ASP:O	1:A:108:ARG:HG3	2.06	0.56
2:B:47:PRO:O	2:B:49:VAL:N	2.38	0.56
1:A:3:HIS:HD1	1:A:29:ASP:CG	2.09	0.55
1:A:144:LYS:NZ	1:A:148:GLU:OE2	2.40	0.55
2:B:47:PRO:C	2:B:49:VAL:H	2.10	0.55
1:A:66:LYS:HD3	3:P:2:ILE:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:ASP:OD2	4:D:229:SER:HB2	2.07	0.54
1:A:42:ASN:HD21	1:A:44:ARG:CG	2.21	0.54
1:A:194:ARG:HB3	1:A:195:PRO:HD2	1.90	0.54
1:A:219:LEU:HD12	1:A:256:TYR:O	2.07	0.54
1:A:258:THR:HG22	1:A:273:ARG:CD	2.37	0.54
2:B:42:ASN:CG	2:B:76:ASP:HB3	2.28	0.54
4:D:225:MET:O	4:D:225:MET:HG3	2.07	0.54
4:D:230:ARG:NH1	4:D:245:CYS:CB	2.70	0.54
4:D:230:ARG:HH12	4:D:245:CYS:HB2	1.72	0.54
1:A:189:VAL:HA	1:A:202:ARG:O	2.07	0.54
1:A:258:THR:CG2	1:A:273:ARG:HE	2.21	0.53
1:A:274:TRP:OXT	1:A:274:TRP:CG	2.61	0.53
4:D:257:LEU:O	4:D:259:LYS:N	2.42	0.53
1:A:11:ALA:HB3	1:A:95:ILE:HB	1.91	0.53
2:B:17:ASN:CB	2:B:73:THR:HB	2.38	0.53
2:B:17:ASN:CG	2:B:73:THR:HA	2.28	0.53
1:A:23:MET:CE	1:A:35:ARG:HH21	2.21	0.53
1:A:226:GLN:O	1:A:227:ASP:HB2	2.09	0.53
2:B:41:LYS:H	2:B:41:LYS:CE	2.13	0.53
1:A:28:VAL:O	1:A:29:ASP:HB2	2.08	0.52
1:A:106:ASP:C	1:A:108:ARG:H	2.12	0.52
1:A:159:TYR:CE1	3:P:3:ILE:HB	2.44	0.52
1:A:202:ARG:HD2	1:A:204:TRP:NE1	2.24	0.52
1:A:202:ARG:NH1	1:A:244:TRP:CZ3	2.77	0.52
2:B:41:LYS:HE3	2:B:41:LYS:CA	2.38	0.52
4:D:220:MET:O	4:D:220:MET:HG2	2.07	0.52
4:D:145:CYS:HA	4:D:150:CYS:HA	1.91	0.52
4:D:195:TYR:CD2	4:D:252:ILE:HD11	2.45	0.52
1:A:33:PHE:O	1:A:34:VAL:HG13	2.10	0.52
4:D:154:ILE:HD12	4:D:154:ILE:N	2.25	0.52
1:A:194:ARG:HG2	1:A:194:ARG:HH11	1.75	0.52
4:D:205:LYS:HB2	4:D:205:LYS:NZ	2.25	0.52
1:A:58:GLU:CD	1:A:58:GLU:H	2.13	0.51
2:B:69:GLU:H	2:B:69:GLU:CD	2.14	0.51
1:A:215:LEU:CD2	1:A:261:VAL:HG22	2.40	0.51
2:B:85:ALA:O	2:B:86:SER:C	2.49	0.51
4:D:228:LYS:O	4:D:229:SER:HB3	2.10	0.51
2:B:28:THR:HB	2:B:63:TYR:CE2	2.46	0.51
2:B:41:LYS:HE2	2:B:44:LYS:CB	2.41	0.51
1:A:258:THR:HG23	1:A:273:ARG:HE	1.76	0.51
1:A:261:VAL:HB	1:A:270:LEU:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:141:LYS:CB	4:D:141:LYS:HZ3	2.23	0.51
1:A:213:ILE:HG12	1:A:263:HIS:HB2	1.93	0.51
4:D:144:PHE:C	4:D:144:PHE:CD2	2.84	0.51
2:B:60:TRP:CZ2	4:D:239:ARG:NH1	2.79	0.50
1:A:32:GLU:O	1:A:48:ARG:HB2	2.11	0.50
1:A:5:LEU:O	1:A:100:GLY:HA3	2.11	0.50
2:B:46:ILE:CG2	2:B:47:PRO:HD2	2.33	0.50
4:D:240:ILE:HG23	4:D:240:ILE:O	2.10	0.50
1:A:20:PRO:HD2	1:A:75:ARG:HG2	1.93	0.50
1:A:234:ARG:HE	1:A:242:GLN:HG3	1.77	0.50
4:D:152:TYR:HB2	4:D:253:CYS:HB2	1.93	0.50
4:D:190:VAL:CG1	4:D:195:TYR:CE1	2.95	0.50
1:A:44:ARG:HH22	1:A:61:GLU:HA	1.77	0.49
1:A:188:HIS:N	1:A:272:LEU:HD21	2.27	0.49
4:D:204:LYS:O	4:D:206:LYS:N	2.45	0.49
4:D:222:ILE:O	4:D:223:LYS:O	2.30	0.49
2:B:41:LYS:CE	2:B:44:LYS:H	2.24	0.49
1:A:3:HIS:ND1	1:A:29:ASP:OD2	2.45	0.49
1:A:258:THR:HG22	1:A:273:ARG:HD2	1.94	0.49
2:B:41:LYS:HE2	2:B:44:LYS:C	2.32	0.49
2:B:59:ASP:O	2:B:60:TRP:HB2	2.13	0.49
2:B:23:LEU:HB2	2:B:70:PHE:CD1	2.48	0.49
4:D:178:GLU:HA	4:D:178:GLU:OE1	2.11	0.49
1:A:197:ASP:OD2	1:A:197:ASP:C	2.51	0.49
1:A:253:LYS:HE3	1:A:256:TYR:CD2	2.48	0.48
1:A:170:ARG:O	1:A:173:LYS:N	2.45	0.48
1:A:202:ARG:HD2	1:A:204:TRP:HE1	1.78	0.48
4:D:141:LYS:HD2	4:D:170:TYR:OH	2.13	0.48
4:D:158:THR:O	4:D:248:PRO:HA	2.14	0.48
4:D:196:TRP:CE2	4:D:249:TYR:HD1	2.32	0.48
4:D:196:TRP:CD1	4:D:249:TYR:HB2	2.47	0.48
1:A:102:GLU:O	1:A:109:LEU:HD12	2.13	0.48
4:D:257:LEU:O	4:D:259:LYS:HG2	2.14	0.48
1:A:14:ARG:HH22	1:A:39:ASP:CG	2.16	0.48
4:D:206:LYS:O	4:D:207:GLU:CB	2.62	0.48
1:A:66:LYS:HD3	3:P:2:ILE:HG23	1.95	0.48
1:A:273:ARG:HH11	1:A:273:ARG:CG	2.26	0.48
4:D:221:LYS:C	4:D:223:LYS:H	2.17	0.48
1:A:81:LEU:HD12	1:A:95:ILE:HD11	1.95	0.47
1:A:85:TYR:O	1:A:86:ASN:CB	2.60	0.47
1:A:125:ALA:O	1:A:134:THR:HG22	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:2:ILE:O	3:P:2:ILE:HG23	2.13	0.47
1:A:199:VAL:HG23	1:A:249:VAL:CG2	2.43	0.47
2:B:28:THR:OG1	2:B:29:GLN:HG3	2.14	0.47
1:A:41:GLU:CD	1:A:41:GLU:H	2.18	0.47
1:A:49:ALA:HB1	1:A:51:TRP:NE1	2.30	0.47
1:A:167:TRP:HB3	1:A:171:TYR:CE2	2.50	0.47
2:B:25:CYS:HB2	2:B:39:MET:HE3	1.95	0.47
4:D:163:CYS:O	4:D:164:LYS:C	2.53	0.47
4:D:175:LEU:CD1	4:D:176:LYS:H	2.26	0.47
4:D:177:ILE:CD1	4:D:183:LEU:HD13	2.38	0.47
4:D:226:ASN:O	4:D:227:PHE:CB	2.62	0.47
4:D:227:PHE:CD1	4:D:227:PHE:C	2.88	0.47
1:A:60:TRP:HA	1:A:60:TRP:CE3	2.50	0.47
2:B:72:PRO:C	2:B:73:THR:CG2	2.73	0.47
2:B:73:THR:HG21	2:B:75:THR:HG23	1.93	0.47
3:P:3:ILE:HG12	3:P:4:ASN:H	1.77	0.47
4:D:151:TYR:HB2	4:D:153:PHE:HE1	1.80	0.47
4:D:153:PHE:HB3	4:D:155:MET:CE	2.45	0.46
4:D:205:LYS:O	4:D:206:LYS:C	2.53	0.46
1:A:35:ARG:HD3	1:A:48:ARG:NH2	2.30	0.46
4:D:204:LYS:C	4:D:206:LYS:H	2.18	0.46
1:A:34:VAL:HB	1:A:45:TYR:CE1	2.51	0.46
1:A:97:VAL:HG11	3:P:5:PHE:CE1	2.50	0.46
2:B:9:VAL:O	2:B:95:TRP:HD1	1.97	0.46
2:B:45:LYS:HZ1	2:B:48:LYS:HG3	1.81	0.46
1:A:138:MET:O	1:A:141:LEU:HB2	2.15	0.46
2:B:68:THR:HG23	2:B:68:THR:O	2.15	0.46
1:A:82:LEU:HD23	1:A:87:GLN:HB2	1.98	0.46
1:A:81:LEU:HD21	1:A:123:TYR:CZ	2.51	0.46
2:B:41:LYS:NZ	2:B:44:LYS:H	2.13	0.46
2:B:45:LYS:HZ2	2:B:48:LYS:HG3	1.80	0.46
1:A:78:LEU:HD13	1:A:95:ILE:HD13	1.98	0.46
2:B:4:THR:HG23	2:B:5:PRO:HD2	1.98	0.46
1:A:182:THR:HG21	1:A:265:GLY:HA2	1.98	0.46
1:A:168:LEU:O	1:A:172:LEU:HD12	2.16	0.45
4:D:175:LEU:HG	4:D:176:LYS:N	2.31	0.45
1:A:127:ASN:ND2	1:A:132:THR:O	2.49	0.45
3:P:2:ILE:HD11	3:P:5:PHE:CE2	2.51	0.45
1:A:42:ASN:ND2	1:A:44:ARG:CG	2.79	0.45
4:D:220:MET:O	4:D:220:MET:CG	2.64	0.45
1:A:224:LEU:C	1:A:226:GLN:N	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:9:VAL:HA	2:B:24:ASN:O	2.15	0.45
4:D:211:ILE:HD13	4:D:211:ILE:O	2.16	0.45
1:A:96:GLN:O	1:A:116:TYR:HA	2.15	0.45
1:A:106:ASP:C	1:A:108:ARG:N	2.70	0.45
4:D:190:VAL:HG13	4:D:195:TYR:CZ	2.52	0.45
4:D:141:LYS:HB2	4:D:141:LYS:HZ3	1.80	0.45
4:D:197:ILE:O	4:D:199:LEU:N	2.48	0.45
1:A:191:HIS:O	1:A:192:HIS:ND1	2.50	0.45
2:B:6:GLN:O	2:B:27:VAL:HA	2.16	0.45
2:B:97:ARG:HG3	2:B:98:ASP:H	1.82	0.45
1:A:5:LEU:O	1:A:100:GLY:CA	2.65	0.45
4:D:191:ILE:CG2	4:D:192:PRO:HD2	2.45	0.44
1:A:82:LEU:HD22	1:A:87:GLN:O	2.16	0.44
4:D:160:TRP:HZ3	4:D:164:LYS:HD2	1.81	0.44
1:A:165:VAL:O	1:A:169:ARG:HG3	2.17	0.44
1:A:167:TRP:CD1	3:P:1:SER:HB3	2.52	0.44
4:D:244:ASP:OD2	4:D:245:CYS:N	2.51	0.44
1:A:213:ILE:HD13	1:A:214:THR:H	1.82	0.44
1:A:170:ARG:NH1	1:A:170:ARG:HG3	2.32	0.44
2:B:57:SER:O	2:B:58:LYS:C	2.56	0.44
1:A:44:ARG:O	1:A:46:GLU:HG2	2.18	0.44
2:B:41:LYS:HE2	2:B:44:LYS:N	2.32	0.44
4:D:150:CYS:SG	4:D:257:LEU:HD11	2.58	0.44
2:B:42:ASN:ND2	2:B:76:ASP:CG	2.68	0.44
4:D:166:ASN:O	4:D:169:HIS:N	2.50	0.44
4:D:175:LEU:CG	4:D:176:LYS:N	2.81	0.44
1:A:97:VAL:HA	1:A:115:GLN:O	2.17	0.44
2:B:60:TRP:NE1	4:D:239:ARG:NH2	2.66	0.44
4:D:150:CYS:SG	4:D:257:LEU:HD21	2.58	0.44
1:A:42:ASN:ND2	1:A:44:ARG:HG2	2.33	0.43
1:A:262:TYR:CD2	4:D:203:LYS:HE2	2.53	0.43
4:D:175:LEU:HD21	4:D:186:LEU:HD11	2.00	0.43
1:A:103:VAL:HB	1:A:107:GLY:HA2	1.99	0.43
2:B:85:ALA:O	2:B:87:MET:N	2.51	0.43
1:A:27:TYR:HA	1:A:31:THR:O	2.18	0.43
1:A:194:ARG:CB	1:A:195:PRO:HD2	2.47	0.43
1:A:224:LEU:C	1:A:226:GLN:H	2.22	0.43
4:D:175:LEU:HD12	4:D:176:LYS:N	2.27	0.43
2:B:41:LYS:HE3	2:B:42:ASN:N	2.34	0.43
2:B:59:ASP:O	2:B:60:TRP:CB	2.66	0.43
2:B:96:ASP:CB	2:B:99:MET:HB2	2.37	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:184:LYS:O	4:D:188:ARG:HG3	2.19	0.43
1:A:162:GLY:O	1:A:163:THR:C	2.56	0.43
1:A:134:THR:HG23	1:A:134:THR:O	2.19	0.43
1:A:170:ARG:HG3	1:A:170:ARG:HH11	1.83	0.43
1:A:178:THR:HG22	1:A:179:LEU:HD23	2.00	0.43
1:A:79:ARG:HH11	1:A:82:LEU:HD12	1.84	0.43
4:D:142:TYR:HE1	4:D:144:PHE:HB2	1.83	0.43
4:D:148:THR:HG22	4:D:258:ASP:OD2	2.18	0.43
1:A:35:ARG:NH1	2:B:53:ASP:HB3	2.34	0.43
4:D:164:LYS:HE3	4:D:211:ILE:HD11	2.00	0.43
1:A:96:GLN:HB2	1:A:117:ALA:HB3	2.01	0.42
1:A:103:VAL:HB	1:A:107:GLY:C	2.39	0.42
1:A:44:ARG:HH12	1:A:61:GLU:CD	2.21	0.42
1:A:159:TYR:CD1	3:P:3:ILE:HB	2.54	0.42
1:A:206:LEU:HD23	1:A:242:GLN:CB	2.41	0.42
4:D:258:ASP:OD1	4:D:258:ASP:N	2.51	0.42
1:A:44:ARG:HD2	1:A:64:THR:HG21	2.01	0.42
1:A:60:TRP:HA	1:A:60:TRP:HE3	1.83	0.42
2:B:96:ASP:O	2:B:99:MET:HB2	2.20	0.42
4:D:177:ILE:CG2	4:D:218:LEU:HD22	2.49	0.42
4:D:201:TYR:CD2	4:D:230:ARG:O	2.72	0.42
4:D:141:LYS:HE2	4:D:170:TYR:CZ	2.54	0.42
1:A:189:VAL:HG12	1:A:190:THR:N	2.35	0.42
4:D:233:VAL:HG12	4:D:234:PHE:N	2.34	0.42
1:A:74:PHE:O	1:A:77:ASP:N	2.53	0.42
1:A:233:THR:OG1	1:A:243:LYS:HD2	2.20	0.42
2:B:7:ILE:CD1	2:B:82:VAL:HG21	2.49	0.42
3:P:3:ILE:CG1	3:P:4:ASN:H	2.29	0.42
4:D:205:LYS:HB2	4:D:205:LYS:HZ3	1.83	0.42
1:A:46:GLU:HG2	1:A:46:GLU:H	1.48	0.42
1:A:114:GLN:CB	1:A:156:LEU:HD21	2.42	0.42
1:A:97:VAL:HG22	1:A:116:TYR:CD2	2.54	0.41
1:A:82:LEU:CD2	1:A:87:GLN:HB2	2.51	0.41
2:B:71:THR:HA	2:B:72:PRO:HD2	1.87	0.41
4:D:174:ILE:HD13	4:D:174:ILE:C	2.41	0.41
2:B:41:LYS:HE2	2:B:44:LYS:H	1.86	0.41
4:D:141:LYS:HB2	4:D:141:LYS:HZ2	1.80	0.41
1:A:63:GLU:OE1	3:P:1:SER:CA	2.68	0.41
1:A:215:LEU:HD21	1:A:261:VAL:HG22	2.03	0.41
4:D:141:LYS:CE	4:D:170:TYR:CZ	3.03	0.41
4:D:175:LEU:HD21	4:D:186:LEU:CD1	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:257:LEU:C	4:D:259:LYS:H	2.23	0.41
2:B:87:MET:SD	2:B:91:LYS:HE3	2.60	0.41
1:A:144:LYS:O	1:A:145:HIS:C	2.59	0.41
1:A:198:LYS:C	1:A:199:VAL:HG13	2.41	0.41
1:A:151:GLY:O	1:A:154:GLU:HB3	2.21	0.41
2:B:69:GLU:N	2:B:69:GLU:CD	2.74	0.41
4:D:233:VAL:CG1	4:D:234:PHE:N	2.83	0.41
1:A:184:SER:HA	1:A:185:PRO:HD3	1.85	0.41
4:D:174:ILE:HD11	4:D:197:ILE:HA	2.03	0.41
2:B:37:ILE:CD1	2:B:82:VAL:HG22	2.50	0.41
2:B:50:GLU:HB2	2:B:67:HIS:CE1	2.56	0.40
4:D:151:TYR:HB2	4:D:153:PHE:CE1	2.56	0.40
1:A:170:ARG:O	1:A:171:TYR:C	2.58	0.40
1:A:185:PRO:HG3	1:A:208:PHE:CB	2.51	0.40
1:A:249:VAL:HB	1:A:257:TYR:CE1	2.55	0.40
1:A:258:THR:CG2	1:A:273:ARG:NE	2.84	0.40
4:D:190:VAL:CG1	4:D:191:ILE:N	2.84	0.40
4:D:205:LYS:HZ3	4:D:205:LYS:CB	2.34	0.40
1:A:74:PHE:O	1:A:75:ARG:C	2.57	0.40
1:A:226:GLN:O	1:A:227:ASP:CB	2.69	0.40
2:B:29:GLN:HA	2:B:61:SER:CB	2.47	0.40
4:D:196:TRP:CE2	4:D:249:TYR:CD1	3.09	0.40
4:D:205:LYS:NZ	4:D:205:LYS:CB	2.84	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	271/274 (99%)	226 (83%)	30 (11%)	15 (6%)	2	5
2	B	97/99 (98%)	73 (75%)	13 (13%)	11 (11%)	0	1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	P	6/8 (75%)	4 (67%)	2 (33%)	0	100	100
4	D	120/122 (98%)	86 (72%)	18 (15%)	16 (13%)	0	0
All	All	494/503 (98%)	389 (79%)	63 (13%)	42 (8%)	1	2

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	45	TYR
2	B	42	ASN
2	B	47	PRO
2	B	48	LYS
2	B	49	VAL
2	B	73	THR
4	D	205	LYS
4	D	223	LYS
4	D	225	MET
4	D	227	PHE
4	D	228	LYS
4	D	232	CYS
1	A	107	GLY
1	A	227	ASP
2	B	75	THR
2	B	86	SER
4	D	139	GLY
4	D	175	LEU
4	D	198	GLY
4	D	229	SER
4	D	230	ARG
4	D	258	ASP
1	A	51	TRP
1	A	54	GLN
1	A	86	ASN
1	A	146	LYS
1	A	179	LEU
1	A	197	ASP
2	B	12	ARG
2	B	72	PRO
4	D	206	LYS
1	A	29	ASP
1	A	43	PRO
1	A	147	TRP

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Mol	Chain	Res	Type
1	A	170	ARG
2	B	85	ALA
1	A	178	THR
1	A	264	GLN
4	D	222	ILE
4	D	190	VAL
4	D	192	PRO
2	B	64	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	232/232 (100%)	224 (97%)	8 (3%)	37	71
2	B	93/93 (100%)	86 (92%)	7 (8%)	13	37
3	P	8/8 (100%)	7 (88%)	1 (12%)	4	14
4	D	111/111 (100%)	90 (81%)	21 (19%)	1	4
All	All	444/444 (100%)	407 (92%)	37 (8%)	11	32

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ASP
1	A	58	GLU
1	A	113	TYR
1	A	114	GLN
1	A	131	LYS
1	A	213	ILE
1	A	254	GLU
1	A	273	ARG
2	B	41	LYS
2	B	47	PRO
2	B	48	LYS
2	B	54	MET
2	B	69	GLU

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Mol	Chain	Res	Type
2	B	70	PHE
2	B	73	THR
3	P	1	SER
4	D	141	LYS
4	D	142	TYR
4	D	144	PHE
4	D	152	TYR
4	D	156	ASN
4	D	166	ASN
4	D	174	ILE
4	D	179	ASP
4	D	206	LYS
4	D	211	ILE
4	D	213	ASN
4	D	220	MET
4	D	222	ILE
4	D	225	MET
4	D	227	PHE
4	D	228	LYS
4	D	230	ARG
4	D	241	GLU
4	D	246	ASN
4	D	256	LYS
4	D	258	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	54	GLN
1	A	72	GLN
1	A	114	GLN
1	A	115	GLN
1	A	174	ASN
1	A	191	HIS
1	A	242	GLN
2	B	42	ASN
2	B	67	HIS
4	D	166	ASN
4	D	168	GLN
4	D	187	GLN
4	D	213	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	273/274 (99%)	0.30	17 (6%) 20 16	22, 74, 124, 190	0
2	B	99/99 (100%)	0.35	7 (7%) 16 12	37, 72, 114, 137	0
3	P	8/8 (100%)	-0.13	0 100 100	56, 59, 81, 81	0
4	D	122/122 (100%)	0.52	14 (11%) 4 3	28, 62, 133, 181	0
All	All	502/503 (99%)	0.36	38 (7%) 13 10	22, 71, 125, 190	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	227	PHE	11.9
4	D	229	SER	5.9
4	D	226	ASN	4.1
4	D	230	ARG	4.0
4	D	259	LYS	3.9
2	B	18	GLY	3.5
4	D	258	ASP	3.4
1	A	44	ARG	3.3
1	A	195	PRO	3.2
1	A	177	ALA	3.0
1	A	251	LEU	3.0
4	D	231	GLY	2.9
1	A	110	LEU	2.8
4	D	251	CYS	2.6
2	B	19	LYS	2.6
1	A	2	PRO	2.6
4	D	215	PRO	2.6
4	D	228	LYS	2.5
4	D	253	CYS	2.5
4	D	138	ARG	2.5
4	D	142	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	89	GLU	2.4
2	B	47	PRO	2.3
1	A	54	GLN	2.3
2	B	80	CYS	2.3
1	A	86	ASN	2.3
1	A	35	ARG	2.2
1	A	58	GLU	2.2
1	A	197	ASP	2.2
1	A	41	GLU	2.2
1	A	149	GLN	2.2
1	A	199	VAL	2.2
2	B	1	ILE	2.1
2	B	25	CYS	2.1
4	D	139	GLY	2.1
1	A	155	ARG	2.1
1	A	45	TYR	2.0
1	A	40	ALA	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.