



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:42 AM UTC

PDB ID : 1DPZ / pdb_00001dpz
Title : STRUCTURE OF MODIFIED 3-ISOPROPYLMALATE DEHYDROGENASE AT THE C-TERMINUS, HD711
Authors : Nurachman, Z.; Akanuma, S.; Sato, T.; Oshima, T.; Tanaka, N.
Deposited on : 1999-12-29
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

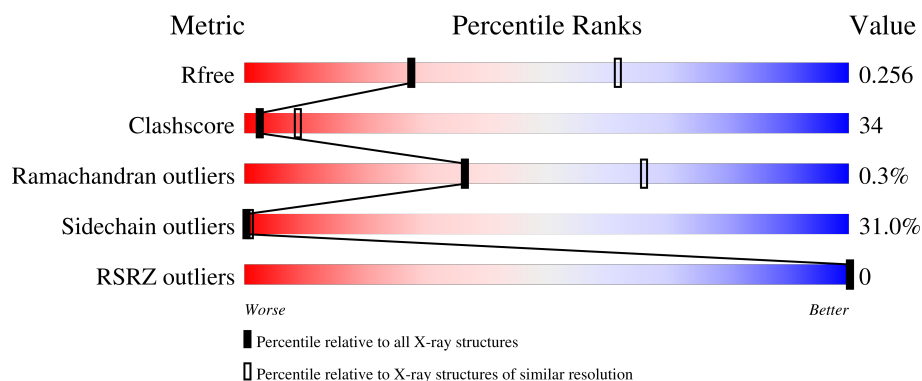
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	349	
1	B	349	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5248 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 3-ISOPROPYLMALATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	0	0
			2592	1654	447	485	6			
1	B	344	Total	C	N	O	S	0	0	0
			2587	1651	446	484	6			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	85	ARG	SER	SEE REMARK 999	UNP Q5SIY4
A	172	LEU	ALA	SEE REMARK 999	UNP Q5SIY4
A	341	GLU	LEU	SEE REMARK 999	UNP Q5SIY4
A	342	ALA	ARG	SEE REMARK 999	UNP Q5SIY4
A	343	PHE	HIS	SEE REMARK 999	UNP Q5SIY4
A	344	THR	LEU	SEE REMARK 999	UNP Q5SIY4
A	345	ALA	ALA	SEE REMARK 999	UNP Q5SIY4
A	346	THR	-	SEE REMARK 999	UNP Q5SIY4
A	347	VAL	-	SEE REMARK 999	UNP Q5SIY4
A	348	GLY	-	SEE REMARK 999	UNP Q5SIY4
A	349	ILE	-	SEE REMARK 999	UNP Q5SIY4
B	85	ARG	SER	SEE REMARK 999	UNP Q5SIY4
B	172	LEU	ALA	SEE REMARK 999	UNP Q5SIY4
B	341	GLU	LEU	SEE REMARK 999	UNP Q5SIY4
B	342	ALA	ARG	SEE REMARK 999	UNP Q5SIY4
B	343	PHE	HIS	SEE REMARK 999	UNP Q5SIY4
B	344	THR	LEU	SEE REMARK 999	UNP Q5SIY4
B	345	ALA	ALA	SEE REMARK 999	UNP Q5SIY4
B	346	THR	-	SEE REMARK 999	UNP Q5SIY4
B	347	VAL	-	SEE REMARK 999	UNP Q5SIY4
B	348	GLY	-	SEE REMARK 999	UNP Q5SIY4
B	349	ILE	-	SEE REMARK 999	UNP Q5SIY4

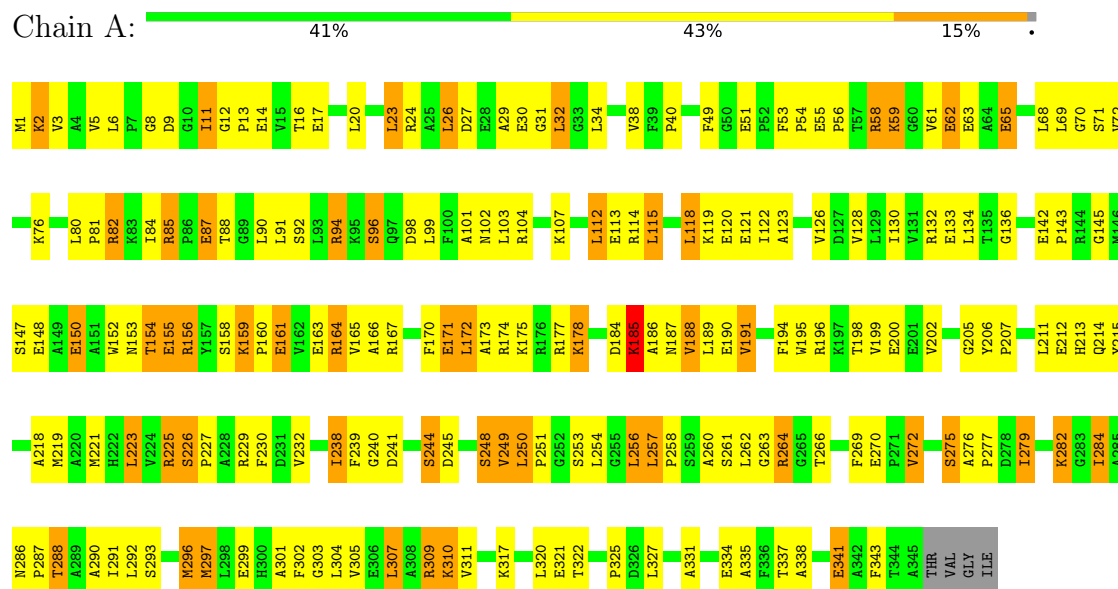
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	31	Total 31	O 31	0	0
2	B	38	Total 38	O 38	0	0

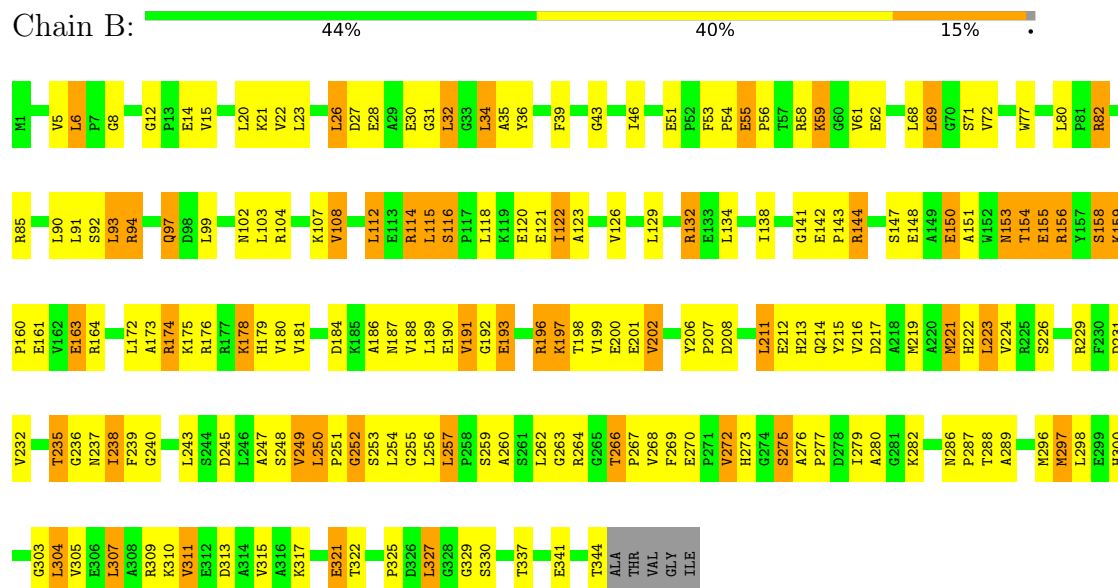
3 Residue-property plots

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: 3-ISOPROPYLMALATE DEHYDROGENASE



• Molecule 1: 3-ISOPROPYLMALATE DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.43Å 87.57Å 70.89Å 90.00° 100.50° 90.00°	Depositor
Resolution (Å)	6.80 – 2.80 6.80 – 2.80	Depositor EDS
% Data completeness (in resolution range)	86.6 (6.80-2.80) 80.3 (6.80-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.61 (at 2.20Å)	Xtriage
Refinement program	X-PLOR 98.1	Depositor
R, R_{free}	0.177 , 0.267 0.171 , 0.256	Depositor DCC
R_{free} test set	1925 reflections (9.86%)	wwPDB-VP
Wilson B-factor (Å ²)	17.7	Xtriage
Anisotropy	0.699	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.46 , 108.9	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	0.93	EDS
Total number of atoms	5248	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% 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	0.54	1/2647 (0.0%)	0.99	5/3594 (0.1%)
1	B	0.51	0/2642	0.98	10/3587 (0.3%)
All	All	0.53	1/5289 (0.0%)	0.99	15/7181 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	296	MET	SD-CE	-5.48	1.65	1.79

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	LEU	N-CA-C	7.96	124.30	109.10
1	B	224	VAL	N-CA-C	-7.06	107.00	113.71
1	B	266	THR	CA-C-N	6.31	126.26	119.76
1	B	266	THR	C-N-CA	6.31	126.26	119.76
1	B	257	LEU	N-CA-C	6.08	120.70	109.10
1	A	178	LYS	N-CA-C	6.03	119.86	111.90
1	B	300	HIS	N-CA-C	6.01	119.88	112.54
1	B	178	LYS	N-CA-C	5.85	119.62	111.90
1	B	237	ASN	N-CA-C	5.45	116.91	110.97
1	A	130	ILE	CB-CA-C	-5.43	105.36	111.23
1	A	29	ALA	N-CA-C	5.43	116.88	111.07
1	B	12	GLY	CA-C-N	5.23	124.41	118.97
1	B	12	GLY	C-N-CA	5.23	124.41	118.97
1	B	93	LEU	N-CA-C	-5.13	105.58	111.07
1	A	185	LYS	N-CA-C	-5.07	104.28	110.61

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	2592	0	2624	198	0
1	B	2587	0	2619	182	0
2	A	31	0	0	10	0
2	B	38	0	0	13	0
All	All	5248	0	5243	359	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 34.

All (359) 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:B:223:LEU:HB3	2:B:386:HOH:O	1.62	0.98
1:A:1:MET:HG2	1:A:34:LEU:HA	1.48	0.96
1:A:164:ARG:HA	1:A:167:ARG:HD2	1.47	0.95
1:A:69:LEU:HB3	1:A:270:GLU:HB3	1.47	0.93
1:A:309:ARG:HH11	1:A:309:ARG:HG2	1.31	0.92
1:B:123:ALA:O	1:B:126:VAL:HG23	1.70	0.92
1:A:297:MET:SD	2:A:373:HOH:O	2.28	0.90
1:A:221:MET:HB2	1:B:249:VAL:HG11	1.52	0.89
1:A:297:MET:HB2	2:A:373:HOH:O	1.74	0.88
1:B:138:ILE:HB	1:B:155:GLU:HG2	1.55	0.88
1:A:248:SER:HB2	1:A:257:LEU:HD12	1.57	0.87
1:A:80:LEU:HD22	1:A:84:ILE:HD11	1.58	0.83
1:B:215:TYR:CD2	2:B:373:HOH:O	2.32	0.83
1:A:257:LEU:O	1:A:272:VAL:HG23	1.80	0.82
1:B:53:PHE:CE2	1:B:58:ARG:HG3	2.14	0.82
1:B:187:ASN:HB3	1:B:215:TYR:CZ	2.13	0.82
1:B:26:LEU:HD21	1:B:344:THR:OG1	1.80	0.81
1:A:143:PRO:O	1:A:154:THR:HG23	1.79	0.81
1:B:108:VAL:CG2	1:B:126:VAL:HB	2.12	0.80
1:B:143:PRO:O	1:B:154:THR:HG22	1.82	0.79
1:B:184:ASP:O	1:B:215:TYR:HA	1.83	0.77
1:B:215:TYR:HB2	2:B:373:HOH:O	1.82	0.77
1:B:215:TYR:HD2	2:B:373:HOH:O	1.63	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:VAL:O	1:B:26:LEU:HD23	1.85	0.76
1:A:245:ASP:OD2	1:B:217:ASP:HB2	1.85	0.75
1:A:80:LEU:HD23	1:A:81:PRO:HD2	1.70	0.73
1:A:309:ARG:HG2	1:A:309:ARG:NH1	1.96	0.73
1:B:108:VAL:HG23	1:B:126:VAL:HB	1.68	0.73
1:B:235:THR:HG21	2:B:384:HOH:O	1.87	0.73
1:A:187:ASN:HB3	1:A:215:TYR:CZ	2.23	0.72
1:A:290:ALA:CB	2:A:353:HOH:O	2.38	0.71
1:B:298:LEU:HD22	1:B:304:LEU:HD12	1.70	0.71
1:A:20:LEU:O	1:A:24:ARG:HG3	1.89	0.71
1:B:54:PRO:HB2	1:B:56:PRO:HD2	1.72	0.71
1:B:26:LEU:HD12	1:B:307:LEU:HD21	1.71	0.71
1:B:118:LEU:HD13	1:B:122:ILE:HG22	1.73	0.71
1:B:193:GLU:O	1:B:197:LYS:HG2	1.91	0.70
1:A:297:MET:CB	2:A:373:HOH:O	2.33	0.70
1:A:248:SER:HB2	1:A:257:LEU:CD1	2.21	0.69
1:B:129:LEU:CD2	1:B:172:LEU:HD23	2.22	0.69
1:A:94:ARG:HA	1:A:99:LEU:HD12	1.74	0.69
1:A:297:MET:CG	2:A:373:HOH:O	2.38	0.69
1:A:3:VAL:HG21	1:A:297:MET:HE1	1.75	0.69
1:B:99:LEU:HA	1:B:263:GLY:HA3	1.74	0.69
1:B:235:THR:CG2	2:B:384:HOH:O	2.40	0.68
1:A:5:VAL:C	1:A:6:LEU:HD12	2.17	0.68
1:A:85:ARG:O	1:A:88:THR:HG22	1.93	0.68
1:B:341:GLU:HA	1:B:344:THR:HG22	1.76	0.67
1:A:53:PHE:CZ	1:A:58:ARG:HA	2.30	0.67
1:A:194:PHE:CG	1:B:151:ALA:HB2	2.29	0.67
1:A:282:LYS:HB3	1:A:284:ILE:HD13	1.77	0.67
1:A:221:MET:HE3	1:B:249:VAL:HG12	1.77	0.67
1:A:269:PHE:CE2	1:A:297:MET:HA	2.30	0.67
1:A:101:ALA:O	1:A:261:SER:HA	1.96	0.66
1:A:104:ARG:NH1	1:A:272:VAL:HG11	2.11	0.66
1:A:163:GLU:HB3	1:A:167:ARG:HE	1.61	0.66
1:B:159:LYS:O	1:B:163:GLU:HG2	1.94	0.66
1:A:164:ARG:CA	1:A:167:ARG:HD2	2.24	0.66
1:A:62:GLU:HG3	1:A:63:GLU:HG3	1.78	0.65
1:B:27:ASP:O	1:B:31:GLY:HA2	1.96	0.65
1:B:179:HIS:HE1	1:B:212:GLU:HG3	1.61	0.65
1:A:256:LEU:C	1:A:257:LEU:HD23	2.21	0.65
1:A:221:MET:HB2	1:B:249:VAL:CG1	2.26	0.65
1:B:115:LEU:HB3	1:B:252:GLY:HA3	1.76	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:ALA:HB2	1:A:296:MET:HE3	1.77	0.65
1:B:97:GLN:HB3	1:B:99:LEU:HD21	1.79	0.65
1:B:138:ILE:HB	1:B:155:GLU:CG	2.27	0.65
1:A:156:ARG:HD3	1:B:150:GLU:OE1	1.96	0.65
1:B:286:ASN:HD22	1:B:287:PRO:HD2	1.62	0.65
1:A:223:LEU:O	1:A:227:PRO:HG3	1.98	0.64
1:B:286:ASN:ND2	1:B:288:THR:H	1.97	0.63
1:B:223:LEU:HD13	2:B:386:HOH:O	1.98	0.63
1:A:276:ALA:HB1	1:A:279:ILE:HD11	1.80	0.63
1:A:337:THR:HG22	1:A:341:GLU:OE1	1.98	0.63
1:B:214:GLN:HE22	1:B:222:HIS:CD2	2.16	0.63
1:A:1:MET:HE2	1:A:32:LEU:HB2	1.80	0.63
1:A:198:THR:O	1:A:202:VAL:HG23	1.99	0.63
1:B:235:THR:HG23	1:B:236:GLY:N	2.13	0.63
1:B:211:LEU:HD23	1:B:212:GLU:H	1.65	0.62
1:A:256:LEU:HG	1:A:288:THR:HG21	1.82	0.62
1:A:304:LEU:HB3	1:A:307:LEU:HB2	1.82	0.62
1:B:286:ASN:HD22	1:B:287:PRO:CD	2.12	0.62
1:A:219:MET:HE3	1:A:223:LEU:HD22	1.83	0.60
1:A:256:LEU:HA	1:A:288:THR:HG22	1.83	0.60
1:A:99:LEU:HD22	1:A:261:SER:O	2.00	0.60
1:B:14:GLU:OE2	1:B:280:ALA:HA	2.02	0.60
1:A:133:GLU:HG2	1:A:165:VAL:HG11	1.82	0.59
1:A:62:GLU:HG3	1:A:63:GLU:N	2.15	0.59
1:A:159:LYS:N	1:A:160:PRO:HD2	2.17	0.59
1:A:240:GLY:O	1:A:244:SER:OG	2.13	0.59
1:B:181:VAL:HG11	1:B:219:MET:CE	2.32	0.59
1:B:260:ALA:HB2	1:B:296:MET:HE2	1.84	0.59
1:B:94:ARG:HD3	1:B:134:LEU:HD13	1.83	0.59
1:A:27:ASP:OD1	1:A:34:LEU:HB3	2.03	0.59
1:A:123:ALA:O	1:A:126:VAL:HG23	2.02	0.58
1:B:303:GLY:O	1:B:305:VAL:HG23	2.04	0.58
1:A:287:PRO:O	1:A:291:ILE:HG12	2.04	0.58
1:A:256:LEU:HA	1:A:288:THR:CG2	2.33	0.58
1:B:176:ARG:NH1	1:B:231:ASP:OD1	2.37	0.58
1:A:276:ALA:N	1:A:277:PRO:HD3	2.19	0.57
1:A:184:ASP:OD1	1:A:196:ARG:NH1	2.38	0.57
1:A:191:VAL:HG23	1:B:153:ASN:OD1	2.05	0.57
1:B:158:SER:OG	1:B:160:PRO:HD2	2.05	0.57
1:A:309:ARG:HH11	1:A:309:ARG:CG	2.12	0.57
1:B:69:LEU:HB3	1:B:270:GLU:HB3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:GLN:HG3	1:B:266:THR:N	2.20	0.57
1:A:26:LEU:HD22	1:A:307:LEU:HD13	1.86	0.56
1:B:129:LEU:HD21	1:B:172:LEU:HD23	1.85	0.56
1:A:158:SER:OG	1:A:161:GLU:HB2	2.05	0.56
1:A:196:ARG:NH2	1:A:213:HIS:HB3	2.20	0.56
1:A:102:ASN:HB2	1:A:134:LEU:HG	1.86	0.56
1:A:256:LEU:O	1:A:257:LEU:HD23	2.06	0.56
1:B:55:GLU:N	1:B:56:PRO:HD2	2.21	0.56
1:B:286:ASN:HD22	1:B:287:PRO:N	2.04	0.56
1:A:27:ASP:O	1:A:31:GLY:HA2	2.04	0.56
1:B:20:LEU:HD22	1:B:36:TYR:CE2	2.40	0.56
1:B:94:ARG:HD2	2:B:363:HOH:O	2.05	0.55
1:A:104:ARG:CZ	1:A:272:VAL:HG11	2.36	0.55
1:B:77:TRP:HA	1:B:80:LEU:HD13	1.88	0.55
1:A:191:VAL:CG2	1:B:153:ASN:HB2	2.35	0.55
1:A:221:MET:CE	1:B:249:VAL:HG12	2.37	0.55
1:B:26:LEU:HD12	1:B:307:LEU:CD2	2.37	0.55
1:A:187:ASN:OD1	1:A:188:VAL:HG22	2.06	0.55
1:A:142:GLU:OE2	1:A:143:PRO:HA	2.07	0.55
1:A:115:LEU:HD23	1:A:115:LEU:N	2.22	0.55
1:B:94:ARG:HD3	1:B:134:LEU:CD1	2.35	0.54
1:A:163:GLU:O	1:A:167:ARG:HG3	2.07	0.54
1:B:54:PRO:C	1:B:56:PRO:HD2	2.31	0.54
1:A:196:ARG:HG3	1:A:213:HIS:ND1	2.22	0.54
1:B:250:LEU:HB2	1:B:251:PRO:HD3	1.89	0.54
1:A:147:SER:OG	1:A:148:GLU:N	2.40	0.54
1:B:116:SER:HB2	1:B:250:LEU:O	2.07	0.54
1:A:153:ASN:ND2	1:B:189:LEU:HB3	2.23	0.54
1:A:219:MET:HE3	1:A:223:LEU:CD2	2.38	0.54
1:B:114:ARG:HG2	1:B:114:ARG:HH11	1.72	0.54
1:B:223:LEU:CB	2:B:386:HOH:O	2.34	0.54
1:B:235:THR:HG23	1:B:236:GLY:O	2.08	0.54
1:B:102:ASN:HB2	1:B:134:LEU:HG	1.90	0.54
1:B:269:PHE:CD1	1:B:297:MET:HG3	2.43	0.53
1:A:102:ASN:HD22	1:A:132:ARG:NH1	2.07	0.53
1:A:92:SER:O	1:A:96:SER:HB3	2.09	0.53
1:A:206:TYR:N	1:A:207:PRO:HD3	2.24	0.53
1:A:269:PHE:CZ	1:A:297:MET:HA	2.43	0.53
1:B:272:VAL:O	1:B:272:VAL:CG1	2.57	0.53
1:B:190:GLU:N	1:B:190:GLU:OE1	2.42	0.53
1:A:103:LEU:HD21	1:A:172:LEU:HD21	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:PHE:CD1	1:B:151:ALA:HB2	2.44	0.52
1:A:279:ILE:HG22	1:A:284:ILE:HD11	1.91	0.52
1:A:55:GLU:N	1:A:56:PRO:HD2	2.24	0.52
1:B:5:VAL:HG21	1:B:20:LEU:HD21	1.91	0.52
1:B:159:LYS:N	1:B:160:PRO:HD2	2.24	0.52
1:B:255:GLY:O	1:B:289:ALA:HB2	2.10	0.52
1:A:98:ASP:OD2	1:A:264:ARG:HD2	2.09	0.52
1:A:205:GLY:C	1:A:207:PRO:HD3	2.35	0.52
1:B:200:GLU:HG3	1:B:213:HIS:NE2	2.25	0.52
1:B:27:ASP:HA	1:B:32:LEU:HD12	1.90	0.52
1:B:198:THR:O	1:B:202:VAL:HG13	2.10	0.52
1:A:325:PRO:HD3	1:A:331:ALA:O	2.10	0.52
1:A:143:PRO:HD2	1:A:154:THR:HG21	1.92	0.52
1:A:166:ALA:O	1:A:170:PHE:HD2	1.92	0.52
1:A:99:LEU:HA	1:A:263:GLY:HA3	1.92	0.51
1:A:260:ALA:HB2	1:A:296:MET:CE	2.38	0.51
1:A:282:LYS:O	1:A:284:ILE:HG23	2.10	0.51
1:B:6:LEU:CD1	1:B:39:PHE:HB2	2.41	0.51
1:B:275:SER:O	1:B:276:ALA:C	2.53	0.51
1:A:174:ARG:HD2	1:A:206:TYR:HD2	1.74	0.51
1:A:165:VAL:HG23	1:A:166:ALA:N	2.26	0.51
1:A:167:ARG:HG2	1:A:167:ARG:HH11	1.76	0.51
1:A:187:ASN:HB3	1:A:215:TYR:OH	2.11	0.51
1:B:97:GLN:HB3	1:B:99:LEU:CD2	2.41	0.51
1:A:250:LEU:N	1:A:251:PRO:CD	2.74	0.51
1:B:214:GLN:HE22	1:B:222:HIS:HD2	1.57	0.51
1:A:185:LYS:HG3	1:A:188:VAL:CG2	2.40	0.50
1:A:213:HIS:C	1:A:214:GLN:HG2	2.35	0.50
1:B:97:GLN:HG3	1:B:266:THR:H	1.75	0.50
1:A:307:LEU:O	1:A:310:LYS:HB3	2.11	0.50
1:A:32:LEU:H	1:A:32:LEU:HD22	1.75	0.50
1:A:290:ALA:HA	2:A:353:HOH:O	2.12	0.50
1:B:206:TYR:N	1:B:207:PRO:HD3	2.26	0.50
1:B:286:ASN:HD22	1:B:286:ASN:C	2.20	0.50
1:A:213:HIS:O	1:A:214:GLN:HG2	2.12	0.50
1:A:177:ARG:H	1:A:177:ARG:HD3	1.77	0.50
1:B:34:LEU:HD22	1:B:35:ALA:N	2.27	0.50
1:A:54:PRO:HB2	1:A:56:PRO:HD2	1.92	0.49
1:A:133:GLU:CG	1:A:165:VAL:HG11	2.41	0.49
1:B:114:ARG:HH11	1:B:114:ARG:CG	2.25	0.49
1:A:1:MET:HA	1:A:302:PHE:CE1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:LEU:HD21	1:B:344:THR:CG2	2.43	0.49
1:B:90:LEU:HD22	1:B:94:ARG:NH2	2.26	0.49
1:B:248:SER:HB2	1:B:257:LEU:HD13	1.93	0.49
1:A:177:ARG:NH2	1:A:230:PHE:O	2.44	0.49
1:B:53:PHE:CZ	1:B:58:ARG:HA	2.48	0.49
1:A:343:PHE:N	1:A:343:PHE:CD1	2.79	0.49
1:B:187:ASN:OD1	1:B:188:VAL:HG23	2.13	0.49
1:A:82:ARG:HA	1:A:85:ARG:HD3	1.94	0.49
1:A:186:ALA:HB3	1:A:196:ARG:NH1	2.28	0.49
1:A:184:ASP:O	1:A:215:TYR:HA	2.12	0.49
1:B:59:LYS:HD3	1:B:59:LYS:HA	1.49	0.48
1:B:181:VAL:HG11	1:B:219:MET:HE1	1.94	0.48
1:B:132:ARG:HB2	1:B:240:GLY:HA3	1.95	0.48
1:B:142:GLU:OE1	1:B:144:ARG:HG3	2.14	0.48
1:A:190:GLU:N	1:A:190:GLU:OE1	2.46	0.48
1:B:174:ARG:HD3	1:B:208:ASP:OD2	2.13	0.48
1:B:26:LEU:CD2	1:B:344:THR:OG1	2.59	0.48
1:B:115:LEU:HD21	1:B:327:LEU:HG	1.95	0.48
1:A:221:MET:HG3	1:A:225:ARG:HH21	1.79	0.48
1:A:27:ASP:OD2	1:A:27:ASP:C	2.56	0.48
1:A:191:VAL:HG21	1:B:153:ASN:HB2	1.96	0.48
1:A:290:ALA:HB1	2:A:353:HOH:O	2.08	0.47
1:B:6:LEU:HD11	1:B:39:PHE:HB2	1.95	0.47
1:A:226:SER:N	1:A:227:PRO:HD3	2.27	0.47
1:B:26:LEU:CD2	1:B:26:LEU:N	2.77	0.47
1:B:155:GLU:OE2	1:B:191:VAL:HG11	2.14	0.47
1:B:99:LEU:HD11	1:B:268:VAL:HG23	1.96	0.47
1:A:310:LYS:CG	2:A:366:HOH:O	2.62	0.47
1:B:179:HIS:O	1:B:231:ASP:HB3	2.14	0.47
1:A:145:GLY:HA3	1:A:152:TRP:CZ2	2.50	0.47
1:A:118:LEU:HD13	1:B:118:LEU:HD23	1.96	0.47
1:A:164:ARG:CD	1:A:264:ARG:NH2	2.78	0.47
1:A:248:SER:O	1:A:251:PRO:HD2	2.15	0.47
1:A:309:ARG:HA	1:A:309:ARG:HD3	1.52	0.47
1:B:181:VAL:HG11	1:B:219:MET:HE2	1.96	0.47
1:A:12:GLY:O	1:A:16:THR:HG23	2.15	0.47
1:A:76:LYS:O	1:A:76:LYS:HG2	2.15	0.47
1:A:302:PHE:HB3	1:A:304:LEU:HD23	1.97	0.47
1:B:238:ILE:CG2	1:B:239:PHE:N	2.78	0.47
1:A:23:LEU:HD21	1:A:297:MET:HE2	1.98	0.46
1:A:113:GLU:CD	1:A:113:GLU:H	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:ARG:HG2	1:A:167:ARG:NH1	2.31	0.46
1:B:142:GLU:CG	1:B:143:PRO:HA	2.46	0.46
1:A:211:LEU:HD12	1:A:212:GLU:H	1.81	0.46
1:A:276:ALA:N	1:A:277:PRO:CD	2.79	0.46
1:B:141:GLY:HA3	1:B:154:THR:O	2.15	0.46
1:A:212:GLU:CD	1:A:229:ARG:HH22	2.24	0.45
1:A:258:PRO:HG3	1:A:292:LEU:HB3	1.97	0.45
1:B:121:GLU:H	1:B:121:GLU:CD	2.24	0.45
1:A:80:LEU:CD2	1:A:81:PRO:HD2	2.45	0.45
1:B:325:PRO:HA	1:B:329:GLY:O	2.15	0.45
1:B:142:GLU:CD	1:B:143:PRO:HA	2.41	0.45
1:A:249:VAL:C	1:A:251:PRO:HD2	2.40	0.45
1:B:26:LEU:HD11	1:B:344:THR:OG1	2.16	0.45
1:B:179:HIS:HB2	2:B:360:HOH:O	2.17	0.45
1:B:214:GLN:NE2	1:B:222:HIS:HD2	2.15	0.45
1:B:103:LEU:HD12	1:B:103:LEU:N	2.31	0.45
1:B:275:SER:HB2	1:B:277:PRO:HD3	1.97	0.45
1:A:153:ASN:HD22	1:B:189:LEU:HD22	1.81	0.45
1:B:138:ILE:CB	1:B:155:GLU:HG2	2.36	0.45
1:A:8:GLY:HA3	1:A:71:SER:O	2.17	0.45
1:A:276:ALA:O	1:A:279:ILE:HG13	2.17	0.45
1:A:27:ASP:HA	1:A:32:LEU:CD2	2.46	0.45
1:B:215:TYR:CB	2:B:373:HOH:O	2.53	0.45
1:B:298:LEU:CD2	1:B:304:LEU:HD12	2.44	0.45
1:B:199:VAL:HG12	1:B:211:LEU:HD11	1.98	0.44
1:A:13:PRO:O	1:A:17:GLU:HB2	2.17	0.44
1:A:40:PRO:CG	1:A:49:PHE:HE2	2.29	0.44
1:B:34:LEU:HD22	1:B:35:ALA:H	1.83	0.44
1:B:36:TYR:CD1	1:B:36:TYR:C	2.95	0.44
1:B:104:ARG:HE	1:B:104:ARG:HB3	1.71	0.44
1:A:6:LEU:O	1:A:69:LEU:HD12	2.18	0.44
1:B:143:PRO:N	1:B:154:THR:HG21	2.33	0.44
1:B:311:VAL:O	1:B:315:VAL:HG23	2.18	0.44
1:B:216:VAL:HG13	1:B:217:ASP:N	2.32	0.44
1:B:286:ASN:ND2	1:B:287:PRO:HD2	2.31	0.44
1:A:59:LYS:HA	1:A:62:GLU:HG2	1.99	0.44
1:A:195:TRP:O	1:A:199:VAL:HG23	2.18	0.44
1:A:189:LEU:CD2	1:B:138:ILE:HG12	2.47	0.44
1:B:159:LYS:H	1:B:160:PRO:HD2	1.82	0.44
1:B:304:LEU:HD12	1:B:307:LEU:HD12	2.00	0.44
1:A:185:LYS:HE2	1:A:187:ASN:OD1	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:LEU:HD12	1:B:114:ARG:CZ	2.48	0.43
1:A:276:ALA:CB	1:A:279:ILE:HD11	2.45	0.43
1:B:276:ALA:HB1	1:B:279:ILE:HG12	2.01	0.43
1:A:186:ALA:HB3	1:A:196:ARG:HH11	1.82	0.43
1:B:196:ARG:HE	1:B:196:ARG:HB2	1.50	0.43
1:B:55:GLU:N	1:B:56:PRO:CD	2.82	0.43
1:B:143:PRO:CD	1:B:154:THR:HG21	2.48	0.43
1:B:221:MET:HE3	1:B:221:MET:HB3	1.86	0.43
1:B:240:GLY:HA2	2:B:384:HOH:O	2.19	0.43
1:A:85:ARG:NH2	1:A:87:GLU:OE1	2.51	0.43
1:A:94:ARG:HH11	1:A:94:ARG:CG	2.31	0.43
1:B:43:GLY:HA3	1:B:77:TRP:CZ2	2.54	0.43
1:B:132:ARG:HH21	1:B:134:LEU:HD21	1.83	0.43
1:B:247:ALA:HA	1:B:250:LEU:HD22	2.01	0.43
1:A:196:ARG:O	1:A:200:GLU:HB2	2.18	0.43
1:A:102:ASN:ND2	1:A:132:ARG:NH1	2.67	0.43
1:B:266:THR:HA	1:B:267:PRO:HD2	1.87	0.43
1:B:273:HIS:ND1	1:B:273:HIS:C	2.76	0.43
1:B:303:GLY:O	1:B:305:VAL:N	2.51	0.43
1:A:248:SER:HA	1:A:257:LEU:HD11	2.00	0.43
1:A:290:ALA:O	1:A:293:SER:HB2	2.19	0.43
1:A:112:LEU:HD22	1:A:115:LEU:HG	2.01	0.42
1:A:194:PHE:CB	1:B:151:ALA:HB2	2.48	0.42
1:A:297:MET:O	1:A:301:ALA:HB3	2.18	0.42
1:A:290:ALA:CA	2:A:353:HOH:O	2.65	0.42
1:B:173:ALA:HB2	1:B:232:VAL:HG21	2.01	0.42
1:B:259:SER:HB3	1:B:272:VAL:HG23	2.01	0.42
1:B:8:GLY:HA3	1:B:71:SER:O	2.19	0.42
1:B:27:ASP:CG	1:B:34:LEU:H	2.27	0.42
1:B:196:ARG:CZ	1:B:213:HIS:HB3	2.49	0.42
1:A:11:ILE:HG12	1:A:276:ALA:HB3	2.00	0.42
1:A:164:ARG:HD2	1:A:264:ARG:NH2	2.35	0.42
1:B:43:GLY:O	1:B:46:ILE:HB	2.19	0.42
1:B:142:GLU:C	1:B:154:THR:HG21	2.45	0.42
1:B:158:SER:OG	1:B:160:PRO:HG2	2.19	0.42
1:A:173:ALA:HB2	1:A:232:VAL:HG21	2.02	0.42
1:B:141:GLY:O	1:B:144:ARG:HG2	2.20	0.42
1:A:6:LEU:HD12	1:A:6:LEU:N	2.35	0.42
1:A:174:ARG:HD2	1:A:206:TYR:CD2	2.53	0.42
1:B:82:ARG:O	1:B:82:ARG:HG2	2.15	0.42
1:B:272:VAL:O	1:B:272:VAL:HG12	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:SER:OG	1:B:160:PRO:CD	2.67	0.42
1:A:275:SER:O	1:A:276:ALA:C	2.62	0.42
1:A:69:LEU:HD12	1:A:70:GLY:H	1.85	0.42
1:A:238:ILE:HG23	1:A:239:PHE:N	2.34	0.42
1:A:310:LYS:CD	2:A:366:HOH:O	2.68	0.42
1:A:335:ALA:O	1:A:338:ALA:HB3	2.20	0.42
1:B:211:LEU:HD23	1:B:212:GLU:N	2.31	0.42
1:B:273:HIS:C	1:B:273:HIS:HD1	2.28	0.42
1:A:159:LYS:N	1:A:160:PRO:CD	2.82	0.41
1:B:243:LEU:HD12	2:B:384:HOH:O	2.20	0.41
1:B:304:LEU:N	1:B:304:LEU:HD23	2.35	0.41
1:A:2:LYS:H	1:A:65:GLU:HG3	1.84	0.41
1:A:287:PRO:O	1:A:290:ALA:HB3	2.19	0.41
1:B:142:GLU:HG3	1:B:143:PRO:HA	2.02	0.41
1:A:163:GLU:OE2	1:A:167:ARG:NH2	2.51	0.41
1:B:112:LEU:HD12	1:B:112:LEU:HA	1.79	0.41
1:A:118:LEU:CD1	1:B:118:LEU:HD23	2.50	0.41
1:B:103:LEU:N	1:B:103:LEU:CD1	2.84	0.41
1:A:49:PHE:CD2	1:A:54:PRO:HG3	2.56	0.41
1:A:5:VAL:O	1:A:6:LEU:HD12	2.21	0.41
1:B:307:LEU:O	1:B:310:LYS:HB3	2.21	0.41
1:A:27:ASP:HA	1:A:32:LEU:HD22	2.03	0.41
1:A:163:GLU:HB3	1:A:167:ARG:NE	2.34	0.41
1:A:174:ARG:O	1:A:178:LYS:HE3	2.21	0.41
1:A:189:LEU:HD22	1:B:138:ILE:HG12	2.02	0.41
1:B:213:HIS:O	1:B:214:GLN:HG2	2.21	0.41
1:A:1:MET:HE2	1:A:32:LEU:CB	2.49	0.41
1:A:27:ASP:O	1:A:31:GLY:N	2.53	0.41
1:A:164:ARG:HD3	1:A:264:ARG:NH2	2.36	0.41
1:A:171:GLU:OE1	1:A:206:TYR:HE2	2.04	0.41
1:B:238:ILE:HD13	1:B:238:ILE:O	2.20	0.41
1:B:297:MET:HE3	1:B:297:MET:HB3	1.86	0.41
1:A:150:GLU:CD	1:B:156:ARG:HH11	2.28	0.41
1:A:154:THR:HA	1:B:154:THR:HA	2.03	0.41
1:A:223:LEU:C	1:A:227:PRO:HG3	2.46	0.41
1:A:303:GLY:O	1:A:305:VAL:N	2.54	0.41
1:A:211:LEU:HD12	1:A:212:GLU:N	2.36	0.40
1:A:269:PHE:CD1	1:A:269:PHE:N	2.89	0.40
1:A:286:ASN:HD22	1:A:287:PRO:HD2	1.86	0.40
1:B:186:ALA:HB2	1:B:192:GLY:O	2.21	0.40
1:A:112:LEU:HD22	1:A:115:LEU:CG	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:LYS:HB2	1:A:122:ILE:HG13	2.04	0.40
1:A:218:ALA:O	1:A:221:MET:HG2	2.21	0.40
1:A:133:GLU:CD	1:A:136:GLY:H	2.29	0.40
1:B:143:PRO:HD2	1:B:154:THR:HG21	2.03	0.40
1:A:69:LEU:HB3	1:A:270:GLU:CB	2.35	0.40
1:B:143:PRO:O	1:B:154:THR:CG2	2.61	0.40
1:A:155:GLU:N	1:B:153:ASN:O	2.48	0.40
1:B:187:ASN:HB3	1:B:215:TYR:CE2	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	343/349 (98%)	316 (92%)	27 (8%)	0	100	100
1	B	342/349 (98%)	314 (92%)	26 (8%)	2 (1%)	21	51
All	All	685/698 (98%)	630 (92%)	53 (8%)	2 (0%)	36	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	252	GLY
1	B	321	GLU

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	266/269 (99%)	187 (70%)	79 (30%)	0	1
1	B	266/269 (99%)	180 (68%)	86 (32%)	0	1
All	All	532/538 (99%)	367 (69%)	165 (31%)	0	1

All (165) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	9	ASP
1	A	11	ILE
1	A	14	GLU
1	A	23	LEU
1	A	26	LEU
1	A	30	GLU
1	A	32	LEU
1	A	38	VAL
1	A	51	GLU
1	A	58	ARG
1	A	59	LYS
1	A	61	VAL
1	A	62	GLU
1	A	65	GLU
1	A	68	LEU
1	A	72	VAL
1	A	82	ARG
1	A	85	ARG
1	A	87	GLU
1	A	90	LEU
1	A	91	LEU
1	A	94	ARG
1	A	96	SER
1	A	107	LYS
1	A	112	LEU
1	A	114	ARG
1	A	115	LEU
1	A	118	LEU
1	A	120	GLU
1	A	121	GLU
1	A	128	VAL
1	A	150	GLU
1	A	154	THR

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Mol	Chain	Res	Type
1	A	155	GLU
1	A	156	ARG
1	A	159	LYS
1	A	161	GLU
1	A	164	ARG
1	A	171	GLU
1	A	172	LEU
1	A	175	LYS
1	A	185	LYS
1	A	188	VAL
1	A	191	VAL
1	A	223	LEU
1	A	225	ARG
1	A	226	SER
1	A	238	ILE
1	A	241	ASP
1	A	244	SER
1	A	248	SER
1	A	249	VAL
1	A	250	LEU
1	A	253	SER
1	A	254	LEU
1	A	256	LEU
1	A	262	LEU
1	A	264	ARG
1	A	266	THR
1	A	272	VAL
1	A	275	SER
1	A	279	ILE
1	A	282	LYS
1	A	284	ILE
1	A	288	THR
1	A	297	MET
1	A	299	GLU
1	A	307	LEU
1	A	309	ARG
1	A	310	LYS
1	A	311	VAL
1	A	317	LYS
1	A	320	LEU
1	A	321	GLU
1	A	322	THR

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Mol	Chain	Res	Type
1	A	327	LEU
1	A	334	GLU
1	A	341	GLU
1	B	6	LEU
1	B	15	VAL
1	B	21	LYS
1	B	23	LEU
1	B	26	LEU
1	B	28	GLU
1	B	30	GLU
1	B	32	LEU
1	B	34	LEU
1	B	51	GLU
1	B	55	GLU
1	B	59	LYS
1	B	61	VAL
1	B	62	GLU
1	B	68	LEU
1	B	69	LEU
1	B	72	VAL
1	B	82	ARG
1	B	85	ARG
1	B	91	LEU
1	B	92	SER
1	B	93	LEU
1	B	94	ARG
1	B	97	GLN
1	B	107	LYS
1	B	108	VAL
1	B	112	LEU
1	B	114	ARG
1	B	115	LEU
1	B	116	SER
1	B	120	GLU
1	B	122	ILE
1	B	132	ARG
1	B	144	ARG
1	B	147	SER
1	B	148	GLU
1	B	150	GLU
1	B	153	ASN
1	B	154	THR

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Mol	Chain	Res	Type
1	B	155	GLU
1	B	156	ARG
1	B	158	SER
1	B	159	LYS
1	B	161	GLU
1	B	163	GLU
1	B	164	ARG
1	B	174	ARG
1	B	175	LYS
1	B	178	LYS
1	B	180	VAL
1	B	191	VAL
1	B	193	GLU
1	B	196	ARG
1	B	197	LYS
1	B	201	GLU
1	B	202	VAL
1	B	211	LEU
1	B	221	MET
1	B	223	LEU
1	B	226	SER
1	B	229	ARG
1	B	235	THR
1	B	238	ILE
1	B	245	ASP
1	B	249	VAL
1	B	250	LEU
1	B	253	SER
1	B	254	LEU
1	B	256	LEU
1	B	262	LEU
1	B	264	ARG
1	B	272	VAL
1	B	275	SER
1	B	282	LYS
1	B	297	MET
1	B	304	LEU
1	B	307	LEU
1	B	309	ARG
1	B	311	VAL
1	B	313	ASP
1	B	317	LYS

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Mol	Chain	Res	Type
1	B	321	GLU
1	B	322	THR
1	B	327	LEU
1	B	330	SER
1	B	337	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	ASN
1	A	153	ASN
1	A	286	ASN
1	B	179	HIS
1	B	222	HIS
1	B	286	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 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 ⓘ

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	345/349 (98%)	-0.85	0 100 100	11, 39, 59, 83	0
1	B	344/349 (98%)	-0.90	0 100 100	11, 39, 54, 64	0
All	All	689/698 (98%)	-0.87	0 100 100	11, 39, 57, 83	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.