



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:27 PM UTC

PDB ID : 1SEV / pdb_00001sev
Title : Mature and translocatable forms of glyoxysomal malate dehydrogenase have different activities and stabilities but similar crystal structures
Authors : Cox, B.R.; Chit, M.M.; Weaver, T.M.; Bailey, J.; Gietl, C.; Bell, E.; Banaszak, L.J.
Deposited on : 2004-02-18
Resolution : 2.55 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

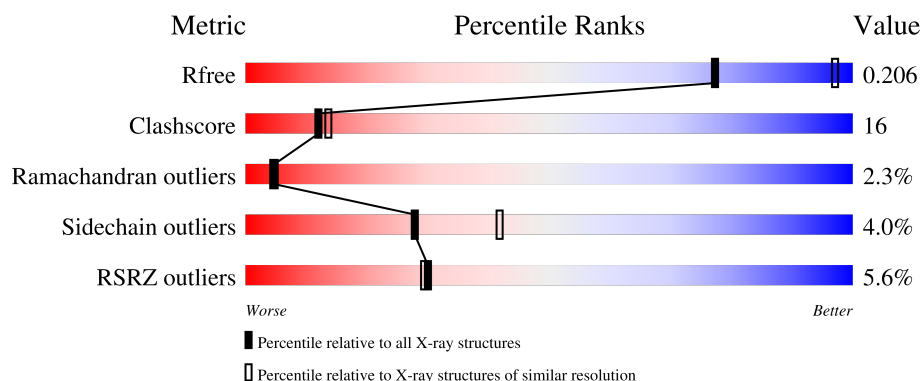
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	362	5% 65% 18% • • 14%
1	B	362	5% 64% 18% • • 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4911 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 Malate dehydrogenase, glyoxysomal precursor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2309	1469	396	431	13			
1	B	313	Total	C	N	O	S	0	0	0
			2309	1469	396	431	13			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	357	HIS	-	expression tag	UNP P19446
A	358	HIS	-	expression tag	UNP P19446
A	359	HIS	-	expression tag	UNP P19446
A	360	HIS	-	expression tag	UNP P19446
A	361	HIS	-	expression tag	UNP P19446
A	362	HIS	-	expression tag	UNP P19446
B	357	HIS	-	expression tag	UNP P19446
B	358	HIS	-	expression tag	UNP P19446
B	359	HIS	-	expression tag	UNP P19446
B	360	HIS	-	expression tag	UNP P19446
B	361	HIS	-	expression tag	UNP P19446
B	362	HIS	-	expression tag	UNP P19446

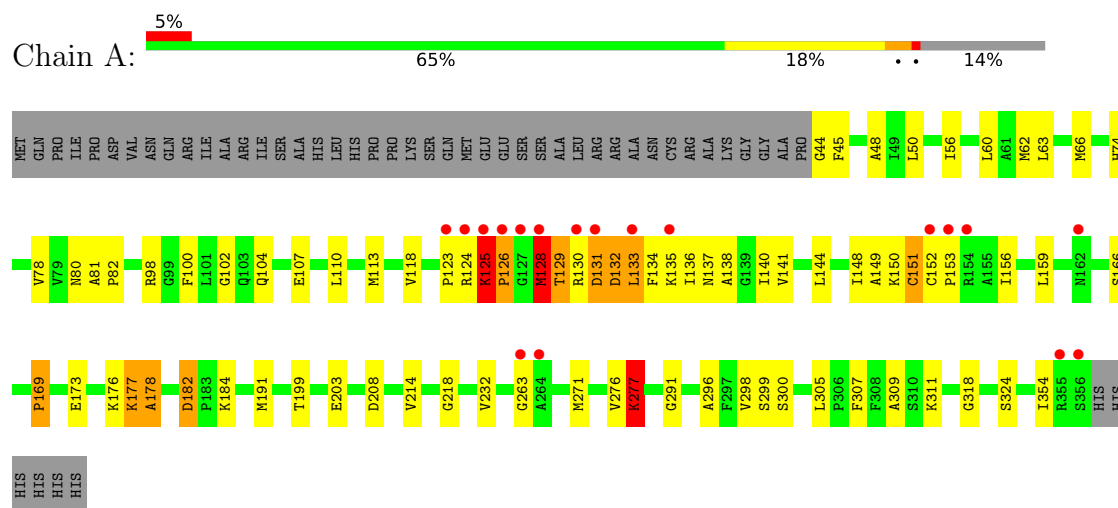
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	151	Total	O	0	0
			151	151		
2	B	142	Total	O	0	0
			142	142		

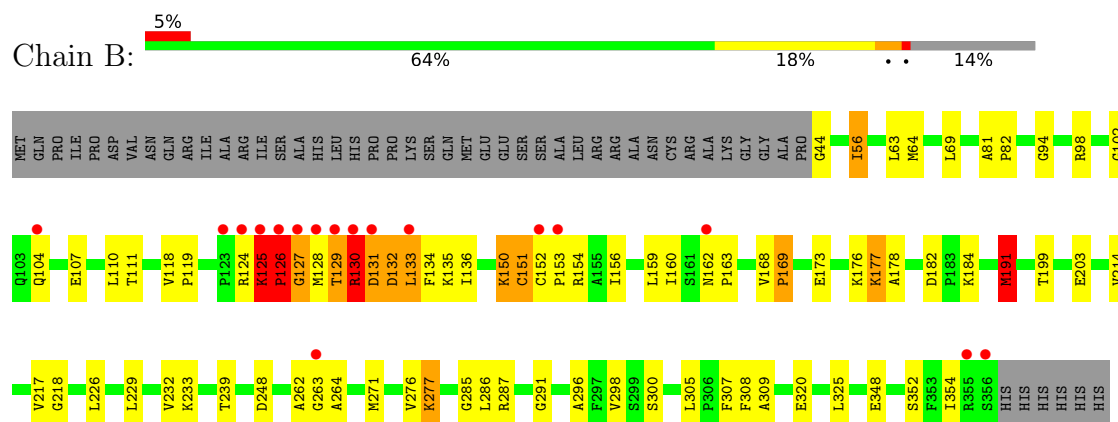
3 Residue-property plots

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: Malate dehydrogenase, glyoxysomal precursor



- Molecule 1: Malate dehydrogenase, glyoxysomal precursor



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	96.80Å 96.80Å 213.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.55 50.00 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.55) 99.8 (50.00-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.96 (at 2.51Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.207 , 0.258 0.206 , 0.206	Depositor DCC
R_{free} test set	1758 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 52.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4911	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.06	4/2345 (0.2%)	1.34	21/3179 (0.7%)
1	B	1.10	4/2345 (0.2%)	1.35	21/3179 (0.7%)
All	All	1.08	8/4690 (0.2%)	1.35	42/6358 (0.7%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	64	MET	SD-CE	-7.12	1.61	1.79
1	A	44	GLY	N-CA	6.73	1.56	1.45
1	B	226	LEU	CA-C	-6.62	1.46	1.52
1	A	156	ILE	CA-CB	-6.42	1.46	1.54
1	A	140	ILE	CA-CB	-5.48	1.48	1.54
1	B	191	MET	SD-CE	5.24	1.92	1.79
1	B	44	GLY	N-CA	5.15	1.53	1.45
1	A	66	MET	SD-CE	-5.08	1.66	1.79

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	125	LYS	CA-C-N	13.19	158.66	127.00
1	B	125	LYS	C-N-CA	13.19	158.66	127.00
1	A	132	ASP	N-CA-C	-12.62	97.34	113.12
1	B	125	LYS	C-N-CD	-12.05	94.09	120.60
1	B	151	CYS	N-CA-C	10.05	125.54	113.28
1	A	151	CYS	N-CA-C	8.98	124.23	113.28
1	A	138	ALA	N-CA-C	-8.54	102.06	111.36
1	B	130	ARG	N-CA-C	-7.82	101.98	114.09
1	B	132	ASP	N-CA-C	-7.45	102.32	111.33
1	B	126	PRO	N-CA-C	7.37	131.27	112.10
1	A	208	ASP	N-CA-C	-7.16	100.77	109.72
1	B	305	LEU	CA-C-N	6.56	126.14	119.19
1	B	305	LEU	C-N-CA	6.56	126.14	119.19
1	B	94	GLY	N-CA-C	6.33	120.33	112.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	45	PHE	N-CA-C	6.24	119.81	110.14
1	A	218	GLY	N-CA-C	6.11	117.29	111.67
1	B	277	LYS	N-CA-C	-6.08	104.65	111.28
1	A	133	LEU	N-CA-C	-6.04	104.39	110.97
1	A	291	GLY	N-CA-C	6.02	123.19	114.10
1	A	277	LYS	N-CA-C	-5.82	104.84	111.07
1	A	78	VAL	N-CA-C	-5.74	106.16	113.22
1	B	56	ILE	CB-CA-C	-5.67	104.61	112.04
1	B	300	SER	N-CA-C	5.65	117.65	109.07
1	A	178	ALA	N-CA-C	-5.63	105.83	114.16
1	B	325	LEU	N-CA-C	5.57	119.33	112.54
1	A	300	SER	N-CA-C	5.53	117.48	109.07
1	B	111	THR	N-CA-C	5.51	118.32	110.10
1	B	239	THR	N-CA-C	-5.49	103.45	110.53
1	B	110	LEU	N-CA-C	5.43	118.12	111.82
1	A	305	LEU	CA-C-N	5.42	125.06	119.05
1	A	305	LEU	C-N-CA	5.42	125.06	119.05
1	B	298	VAL	N-CA-C	5.41	116.69	108.80
1	B	285	GLY	N-CA-C	-5.37	106.00	112.49
1	A	125	LYS	N-CA-C	5.36	121.66	109.81
1	B	291	GLY	N-CA-C	5.33	122.55	114.61
1	A	318	GLY	N-CA-C	-5.26	100.70	113.18
1	A	298	VAL	N-CA-C	5.24	116.25	108.65
1	A	299	SER	N-CA-C	-5.14	102.30	109.96
1	A	60	LEU	N-CA-C	-5.12	105.62	111.14
1	A	324	SER	N-CA-C	-5.10	103.95	110.53
1	B	320	GLU	N-CA-C	-5.08	107.25	113.50
1	A	182	ASP	N-CA-C	-5.04	99.37	108.85

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	2309	0	2393	75	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2309	0	2393	74	0
2	A	151	0	0	3	0
2	B	142	0	0	3	0
All	All	4911	0	4786	148	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 16.

All (148) 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:129:THR:CG2	1:B:132:ASP:H	1.58	1.16
1:A:129:THR:HG23	1:A:131:ASP:H	1.09	1.07
1:B:129:THR:HG23	1:B:131:ASP:H	1.00	1.07
1:B:129:THR:HG23	1:B:131:ASP:N	1.81	0.94
1:B:107:GLU:HG3	1:B:151:CYS:SG	2.12	0.89
1:A:107:GLU:HG3	1:A:151:CYS:SG	2.12	0.89
1:A:129:THR:CG2	1:A:132:ASP:H	1.87	0.88
1:B:129:THR:CG2	1:B:131:ASP:H	1.87	0.87
1:A:173:GLU:HB3	1:A:354:ILE:HG21	1.54	0.86
1:B:129:THR:HG21	1:B:132:ASP:H	1.40	0.86
1:A:129:THR:HG23	1:A:131:ASP:N	1.90	0.85
1:B:129:THR:HG23	1:B:130:ARG:N	1.86	0.85
1:A:124:ARG:NH2	1:A:136:ILE:HD12	1.91	0.84
1:B:129:THR:CG2	1:B:132:ASP:N	2.42	0.82
1:B:129:THR:HG22	1:B:132:ASP:H	1.45	0.81
1:A:311:LYS:HE3	2:A:716:HOH:O	1.79	0.81
1:A:124:ARG:CD	1:A:128:MET:HA	2.12	0.80
1:B:129:THR:HG22	1:B:132:ASP:CG	2.07	0.80
1:A:129:THR:HG21	1:A:132:ASP:N	2.00	0.76
1:B:129:THR:HG23	1:B:130:ARG:H	1.49	0.74
1:B:56:ILE:HD13	1:B:271:MET:HG2	1.69	0.74
1:A:129:THR:HG21	1:A:132:ASP:CG	2.12	0.73
1:B:128:MET:HE3	1:B:133:LEU:HB2	1.69	0.73
1:A:129:THR:CG2	1:A:132:ASP:N	2.51	0.72
1:A:129:THR:HG23	1:A:132:ASP:H	1.54	0.72
1:B:124:ARG:NH1	2:B:893:HOH:O	2.22	0.72
1:A:177:LYS:O	1:A:177:LYS:HG2	1.90	0.71
1:B:63:LEU:HB2	1:B:276:VAL:HG21	1.74	0.70
1:A:129:THR:HG21	1:A:132:ASP:CB	2.21	0.70
1:B:124:ARG:O	1:B:128:MET:HB2	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:MET:HG3	1:A:129:THR:N	2.08	0.69
1:A:124:ARG:HB2	1:A:124:ARG:NH1	2.07	0.69
1:A:128:MET:HG3	1:A:129:THR:H	1.57	0.68
1:A:128:MET:CG	1:A:129:THR:H	2.06	0.68
1:A:169:PRO:HG3	1:A:307:PHE:CD2	2.28	0.68
1:A:102:GLY:HA3	1:A:104:GLN:NE2	2.07	0.68
1:A:124:ARG:HD3	1:A:128:MET:HA	1.77	0.66
1:A:129:THR:HG23	1:A:130:ARG:N	2.08	0.66
1:B:133:LEU:C	1:B:133:LEU:HD23	2.20	0.66
1:A:124:ARG:O	1:A:126:PRO:HD2	1.96	0.66
1:A:104:GLN:CD	1:A:104:GLN:H	2.04	0.66
1:A:124:ARG:HD2	1:A:128:MET:HA	1.76	0.66
1:A:56:ILE:HD13	1:A:271:MET:HG2	1.78	0.66
1:A:63:LEU:HB2	1:A:276:VAL:HG21	1.78	0.64
1:B:133:LEU:HD23	1:B:134:PHE:N	2.12	0.64
1:B:129:THR:CG2	1:B:132:ASP:CG	2.69	0.64
1:B:173:GLU:HB3	1:B:354:ILE:HG21	1.80	0.63
1:B:177:LYS:O	1:B:177:LYS:HG2	1.99	0.62
1:B:156:ILE:HD11	1:B:286:LEU:HD21	1.84	0.60
1:A:177:LYS:O	1:A:177:LYS:CG	2.49	0.60
1:A:62:MET:HE3	1:B:63:LEU:HD21	1.84	0.59
1:B:102:GLY:HA3	1:B:104:GLN:NE2	2.17	0.59
1:A:124:ARG:HB2	1:A:124:ARG:CZ	2.33	0.59
1:A:102:GLY:HA3	1:A:104:GLN:HE22	1.66	0.58
1:A:176:LYS:C	1:A:178:ALA:H	2.12	0.58
1:A:124:ARG:NH1	1:A:124:ARG:CB	2.66	0.58
1:A:173:GLU:HB3	1:A:354:ILE:CG2	2.30	0.57
1:B:156:ILE:CD1	1:B:286:LEU:HD21	2.34	0.57
1:B:199:THR:O	1:B:203:GLU:HG3	2.04	0.57
1:A:129:THR:CG2	1:A:130:ARG:N	2.68	0.57
1:B:104:GLN:H	1:B:104:GLN:CD	2.12	0.56
1:B:129:THR:HG22	1:B:132:ASP:CB	2.35	0.56
1:B:169:PRO:HG3	1:B:307:PHE:CD2	2.40	0.56
1:B:214:VAL:HG22	1:B:232:VAL:HG22	1.88	0.56
1:A:124:ARG:NH1	1:A:133:LEU:HB2	2.21	0.55
1:B:129:THR:HG22	1:B:132:ASP:N	2.18	0.55
1:A:166:SER:C	1:A:169:PRO:HD2	2.32	0.55
1:B:125:LYS:HB2	1:B:128:MET:CG	2.37	0.55
1:B:125:LYS:CG	1:B:128:MET:HG2	2.37	0.54
1:A:277:LYS:NZ	2:A:914:HOH:O	2.36	0.54
1:B:129:THR:HG22	1:B:132:ASP:OD2	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:CYS:N	1:B:153:PRO:HD3	2.22	0.54
1:A:128:MET:O	1:A:129:THR:HB	2.08	0.54
1:B:271:MET:O	1:B:271:MET:HE3	2.08	0.53
1:A:214:VAL:HG22	1:A:232:VAL:HG22	1.89	0.53
1:A:124:ARG:HH22	1:A:136:ILE:HD12	1.73	0.53
1:B:177:LYS:O	1:B:177:LYS:CG	2.55	0.53
1:B:125:LYS:HB2	1:B:128:MET:HG2	1.89	0.53
1:A:124:ARG:NH2	1:A:133:LEU:HA	2.24	0.53
1:A:152:CYS:N	1:A:153:PRO:HD3	2.23	0.53
1:A:81:ALA:HB3	1:A:82:PRO:HD3	1.91	0.52
1:A:48:ALA:HB2	1:A:113:MET:HG3	1.91	0.52
1:B:125:LYS:HG3	1:B:128:MET:CG	2.39	0.52
1:B:129:THR:HB	1:B:132:ASP:OD2	2.08	0.52
1:A:166:SER:O	1:A:169:PRO:HD2	2.10	0.51
1:A:144:LEU:O	1:A:148:ILE:HG13	2.09	0.51
1:A:124:ARG:CB	1:A:124:ARG:HH11	2.24	0.50
1:B:229:LEU:HB2	2:B:826:HOH:O	2.11	0.50
1:B:125:LYS:HG3	1:B:128:MET:HG2	1.94	0.49
1:B:218:GLY:HA2	1:B:308:PHE:CZ	2.46	0.49
1:A:129:THR:O	1:A:130:ARG:HB3	2.13	0.49
1:B:129:THR:CG2	1:B:132:ASP:OD2	2.61	0.49
1:B:176:LYS:C	1:B:178:ALA:H	2.20	0.48
1:A:128:MET:O	1:A:129:THR:CB	2.61	0.48
1:B:296:ALA:O	1:B:309:ALA:HA	2.14	0.48
1:A:130:ARG:HG3	2:A:778:HOH:O	2.13	0.48
1:A:182:ASP:OD1	1:A:184:LYS:HB2	2.14	0.48
1:A:149:ALA:O	1:A:153:PRO:HG3	2.13	0.47
1:B:129:THR:CB	1:B:132:ASP:OD2	2.62	0.47
1:B:102:GLY:HA3	1:B:104:GLN:HE22	1.79	0.47
1:A:125:LYS:HB2	1:A:126:PRO:CD	2.44	0.47
1:B:129:THR:CG2	1:B:130:ARG:H	2.21	0.47
1:B:182:ASP:OD1	1:B:184:LYS:HB2	2.15	0.47
1:A:124:ARG:O	1:A:126:PRO:CD	2.63	0.46
1:B:150:LYS:CB	1:B:150:LYS:NZ	2.78	0.46
1:A:129:THR:HG23	1:A:132:ASP:N	2.26	0.45
1:A:129:THR:HG21	1:A:132:ASP:OD2	2.17	0.45
1:B:129:THR:CG2	1:B:130:ARG:N	2.61	0.45
1:A:50:LEU:HD11	1:A:110:LEU:HD21	1.99	0.45
1:A:125:LYS:CE	1:A:125:LYS:HA	2.47	0.44
1:A:133:LEU:O	1:A:134:PHE:C	2.60	0.44
1:A:125:LYS:HB2	1:A:126:PRO:HD2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:LYS:NZ	1:B:150:LYS:HB2	2.32	0.44
1:B:126:PRO:O	1:B:127:GLY:C	2.61	0.44
1:B:217:VAL:HB	1:B:309:ALA:O	2.17	0.44
1:A:296:ALA:O	1:A:309:ALA:HA	2.17	0.44
1:A:118:VAL:HB	1:A:159:LEU:HD12	2.00	0.44
1:A:124:ARG:HH21	1:A:136:ILE:HD12	1.73	0.44
1:B:262:ALA:C	1:B:264:ALA:H	2.26	0.44
1:B:119:PRO:HA	1:B:160:ILE:HB	1.99	0.43
1:B:262:ALA:O	1:B:264:ALA:N	2.50	0.43
1:A:199:THR:O	1:A:203:GLU:HG3	2.18	0.43
1:B:129:THR:HG21	1:B:132:ASP:N	2.20	0.43
1:B:162:ASN:OD1	1:B:163:PRO:HA	2.19	0.43
1:A:129:THR:O	1:A:130:ARG:CB	2.67	0.43
1:A:124:ARG:HH22	1:A:133:LEU:HA	1.83	0.43
1:B:133:LEU:C	1:B:133:LEU:CD2	2.92	0.43
1:B:232:VAL:CG1	1:B:233:LYS:N	2.82	0.43
1:B:125:LYS:CB	1:B:128:MET:HG2	2.48	0.43
1:A:125:LYS:HA	1:A:125:LYS:HE3	2.01	0.42
1:A:128:MET:C	1:A:129:THR:HG22	2.44	0.42
1:A:176:LYS:O	1:A:178:ALA:N	2.49	0.42
1:B:132:ASP:O	1:B:136:ILE:HG13	2.18	0.42
1:B:118:VAL:HB	1:B:159:LEU:HD12	1.99	0.42
1:B:354:ILE:HD13	1:B:354:ILE:HA	1.85	0.42
1:A:74:HIS:CG	1:A:100:PHE:HE2	2.38	0.42
1:A:137:ASN:O	1:A:141:VAL:HG23	2.19	0.42
1:B:69:LEU:HD21	1:B:287:ARG:NH1	2.35	0.41
1:B:348:GLU:OE2	1:B:352:SER:HB3	2.21	0.41
1:B:124:ARG:C	1:B:128:MET:HG3	2.46	0.41
1:A:63:LEU:CB	1:A:276:VAL:HG21	2.48	0.41
1:A:81:ALA:N	1:A:82:PRO:CD	2.84	0.41
1:B:81:ALA:N	1:B:82:PRO:CD	2.83	0.41
1:B:154:ARG:HB3	1:B:154:ARG:NH1	2.36	0.41
1:A:176:LYS:C	1:A:178:ALA:N	2.75	0.41
1:B:248:ASP:HB2	2:B:902:HOH:O	2.21	0.41
1:B:191:MET:O	1:B:191:MET:HG3	2.20	0.40
1:B:131:ASP:O	1:B:132:ASP:C	2.65	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	311/362 (86%)	295 (95%)	8 (3%)	8 (3%)	4	3
1	B	311/362 (86%)	297 (96%)	8 (3%)	6 (2%)	6	7
All	All	622/724 (86%)	592 (95%)	16 (3%)	14 (2%)	5	4

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	125	LYS
1	A	126	PRO
1	A	128	MET
1	B	125	LYS
1	B	126	PRO
1	A	129	THR
1	A	177	LYS
1	B	129	THR
1	B	263	GLY
1	A	80	ASN
1	A	123	PRO
1	A	263	GLY
1	B	177	LYS
1	B	127	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	247/288 (86%)	238 (96%)	9 (4%)	31	46
1	B	247/288 (86%)	236 (96%)	11 (4%)	24	38
All	All	494/576 (86%)	474 (96%)	20 (4%)	28	42

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	ARG
1	A	125	LYS
1	A	128	MET
1	A	131	ASP
1	A	135	LYS
1	A	150	LYS
1	A	169	PRO
1	A	191	MET
1	A	277	LYS
1	B	98	ARG
1	B	126	PRO
1	B	130	ARG
1	B	131	ASP
1	B	133	LEU
1	B	135	LYS
1	B	150	LYS
1	B	168	VAL
1	B	169	PRO
1	B	191	MET
1	B	277	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	A	301	GLN
1	B	158	ASN
1	B	301	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	313/362 (86%)	-0.01	18 (5%) 29 28	19, 33, 67, 120	0
1	B	313/362 (86%)	-0.18	17 (5%) 31 31	18, 30, 65, 120	0
All	All	626/724 (86%)	-0.10	35 (5%) 30 29	18, 32, 67, 120	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	125	LYS	6.3
1	B	124	ARG	5.6
1	B	127	GLY	5.5
1	A	128	MET	5.4
1	B	126	PRO	5.4
1	A	126	PRO	5.3
1	A	125	LYS	4.9
1	B	356	SER	4.8
1	A	130	ARG	4.7
1	A	124	ARG	4.6
1	B	130	ARG	3.7
1	A	356	SER	3.6
1	A	127	GLY	3.4
1	A	153	PRO	3.3
1	A	152	CYS	3.1
1	B	153	PRO	3.0
1	B	133	LEU	3.0
1	B	162	ASN	2.9
1	B	263	GLY	2.7
1	A	355	ARG	2.7
1	A	263	GLY	2.7
1	A	133	LEU	2.6
1	B	131	ASP	2.5
1	B	128	MET	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	123	PRO	2.4
1	B	152	CYS	2.3
1	B	355	ARG	2.3
1	A	131	ASP	2.2
1	A	162	ASN	2.2
1	A	154	ARG	2.2
1	B	104	GLN	2.1
1	B	129	THR	2.1
1	A	264	ALA	2.1
1	A	135	LYS	2.0
1	B	123	PRO	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 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.