



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:10 AM UTC

PDB ID : 7W26 / pdb_00007w26
Title : monolignol ferulate transferase
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Deposited on : 2021-11-22
Resolution : 2.43 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

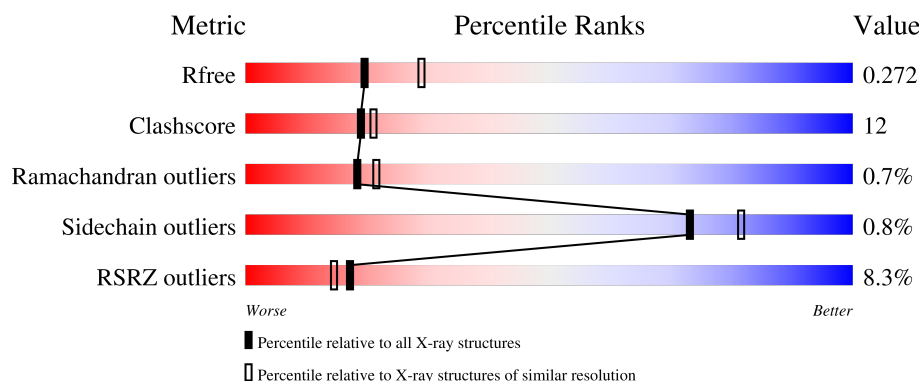
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.43 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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	449	6% 78% 18% ..
1	B	449	10% 73% 23% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7066 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 Ferulate monolignol transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	0	0
			3489	2237	578	657	17			
1	B	440	Total	C	N	O	S	0	0	0
			3482	2232	578	655	17			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	442	LEU	-	expression tag	UNP W8NXL9
A	443	GLU	-	expression tag	UNP W8NXL9
A	444	HIS	-	expression tag	UNP W8NXL9
A	445	HIS	-	expression tag	UNP W8NXL9
A	446	HIS	-	expression tag	UNP W8NXL9
A	447	HIS	-	expression tag	UNP W8NXL9
A	448	HIS	-	expression tag	UNP W8NXL9
A	449	HIS	-	expression tag	UNP W8NXL9
B	442	LEU	-	expression tag	UNP W8NXL9
B	443	GLU	-	expression tag	UNP W8NXL9
B	444	HIS	-	expression tag	UNP W8NXL9
B	445	HIS	-	expression tag	UNP W8NXL9
B	446	HIS	-	expression tag	UNP W8NXL9
B	447	HIS	-	expression tag	UNP W8NXL9
B	448	HIS	-	expression tag	UNP W8NXL9
B	449	HIS	-	expression tag	UNP W8NXL9

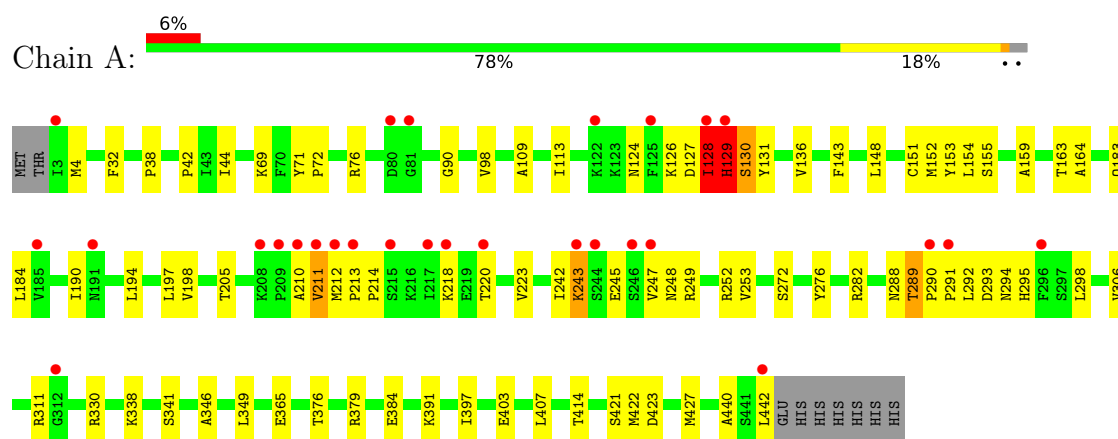
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	54	Total	O	0	0
			54	54		
2	B	41	Total	O	0	0
			41	41		

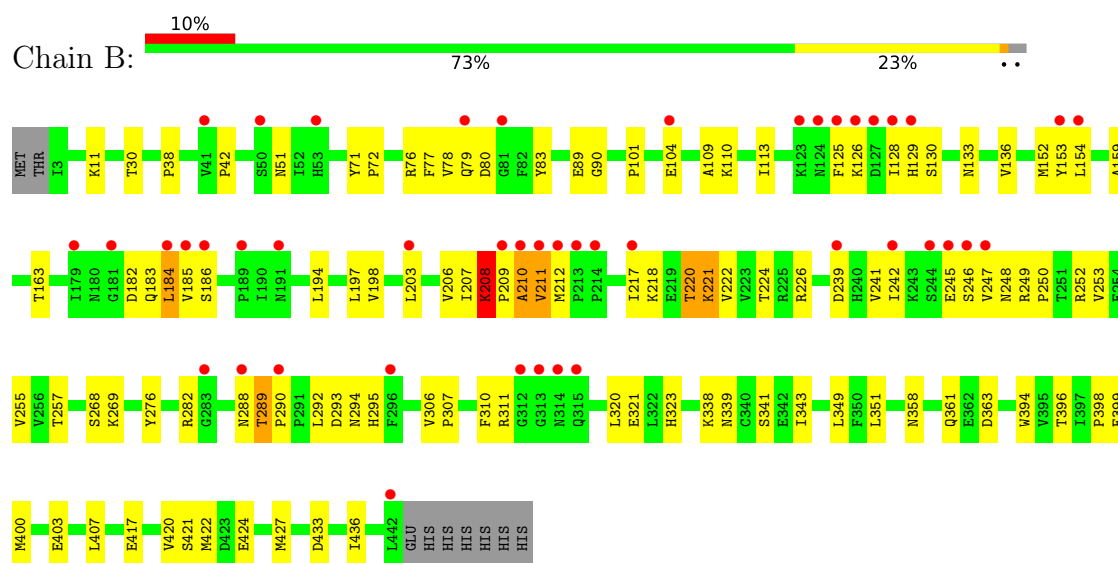
3 Residue-property plots [i](#)

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: Ferulate monolignol transferase



- Molecule 1: Ferulate monolignol transferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	136.42Å 136.42Å 102.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	96.46 – 2.43 96.46 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.7 (96.46-2.43) 99.7 (96.46-2.43)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 2.42Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.216 , 0.266 0.225 , 0.272	Depositor DCC
R_{free} test set	1847 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	50.7	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.5	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	0.94	EDS
Total number of atoms	7066	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.34	1/3567 (0.0%)	0.63	11/4845 (0.2%)
1	B	0.32	2/3560 (0.1%)	0.64	5/4836 (0.1%)
All	All	0.33	3/7127 (0.0%)	0.64	16/9681 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	221	LYS	C-O	-6.50	1.16	1.24
1	A	131	TYR	C-O	-5.97	1.16	1.24
1	B	208	LYS	C-O	-5.46	1.20	1.25

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	TYR	N-CA-C	-7.95	103.04	112.89
1	B	186	SER	N-CA-C	7.92	118.71	108.34
1	A	129	HIS	CA-CB-CG	7.08	120.88	113.80
1	A	249	ARG	CG-CD-NE	7.03	127.47	112.00
1	A	129	HIS	CB-CA-C	-6.99	96.50	110.42
1	B	182	ASP	N-CA-C	6.51	118.97	111.02
1	A	128	ILE	N-CA-C	-5.94	98.66	107.51
1	A	129	HIS	N-CA-CB	5.86	120.39	110.49
1	B	210	ALA	CA-C-N	5.61	132.07	121.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	210	ALA	C-N-CA	5.61	132.07	121.97
1	A	128	ILE	CA-C-N	5.29	131.65	121.54
1	A	128	ILE	C-N-CA	5.29	131.65	121.54
1	B	220	THR	N-CA-C	-5.19	99.32	108.20
1	A	249	ARG	CB-CG-CD	-5.10	99.57	111.30
1	A	243	LYS	CB-CA-C	-5.09	100.29	110.42
1	A	128	ILE	O-C-N	5.06	130.20	122.76

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	289	THR	Peptide
1	B	289	THR	Peptide

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	3489	0	3492	72	0
1	B	3482	0	3476	106	0
2	A	54	0	0	1	0
2	B	41	0	0	2	0
All	All	7066	0	6968	164	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 12.

All (164) 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:209:PRO:HA	1:B:361:GLN:HG2	1.25	1.13
1:A:129:HIS:CE1	1:B:252:ARG:HG3	1.85	1.11
1:B:221:LYS:NZ	1:B:424:GLU:OE1	1.82	1.11
1:B:217:ILE:HG23	1:B:220:THR:HG22	1.32	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:289:THR:HG22	1:B:290:PRO:HD3	1.28	1.10
1:A:289:THR:HG22	1:A:290:PRO:HD2	1.34	1.08
1:B:79:GLN:OE1	1:B:128:ILE:O	1.82	0.97
1:A:129:HIS:NE2	1:B:252:ARG:HG3	1.86	0.90
1:A:128:ILE:CD1	1:B:245:GLU:CD	2.51	0.83
1:B:217:ILE:HG23	1:B:220:THR:CG2	2.08	0.83
1:B:79:GLN:NE2	1:B:129:HIS:CE1	2.50	0.79
1:B:101:PRO:HD2	1:B:104:GLU:OE1	1.82	0.79
1:A:129:HIS:NE2	1:B:252:ARG:N	2.30	0.77
1:B:136:VAL:HG22	1:B:154:LEU:CD1	2.15	0.77
1:B:289:THR:HG22	1:B:290:PRO:CD	2.13	0.77
1:B:221:LYS:NZ	1:B:424:GLU:CD	2.44	0.75
1:B:136:VAL:HG22	1:B:154:LEU:HD11	1.67	0.75
1:B:253:VAL:O	1:B:257:THR:HG23	1.89	0.73
1:A:289:THR:HG22	1:A:290:PRO:CD	2.16	0.72
1:A:128:ILE:HD13	1:B:245:GLU:CD	2.17	0.69
1:B:194:LEU:HD22	1:B:198:VAL:HG22	1.75	0.69
1:A:183:GLN:HG2	1:A:184:LEU:HD22	1.75	0.68
1:B:241:VAL:HG12	1:B:242:ILE:HG13	1.75	0.68
1:B:76:ARG:NH1	1:B:90:GLY:O	2.27	0.68
1:B:77:PHE:HE2	1:B:79:GLN:OE1	1.78	0.67
1:B:217:ILE:CG2	1:B:220:THR:HG22	2.19	0.66
1:B:79:GLN:NE2	1:B:129:HIS:ND1	2.43	0.66
1:B:246:SER:HB2	1:B:248:ASN:HB2	1.78	0.65
1:A:197:LEU:HD12	1:A:292:LEU:HD23	1.77	0.65
1:A:194:LEU:HB2	1:A:298:LEU:HD11	1.78	0.64
1:A:126:LYS:HE3	1:B:249:ARG:HH22	1.62	0.64
1:A:346:ALA:HA	1:A:349:LEU:HD12	1.79	0.63
1:A:136:VAL:HG22	1:A:154:LEU:HD13	1.79	0.63
1:A:128:ILE:HD11	1:B:245:GLU:CD	2.22	0.63
1:B:321:GLU:OE1	1:B:323:HIS:NE2	2.26	0.62
1:A:212:MET:HG3	1:A:214:PRO:HD2	1.81	0.62
1:B:42:PRO:HG2	1:B:398:PRO:HG2	1.83	0.61
1:B:250:PRO:HB2	1:B:255:VAL:HG23	1.82	0.61
1:A:403:GLU:HG2	1:A:422:MET:HA	1.84	0.59
1:A:129:HIS:NE2	1:B:252:ARG:CG	2.63	0.59
1:A:293:ASP:OD1	1:A:294:ASN:N	2.36	0.59
1:A:128:ILE:HD11	1:B:245:GLU:OE1	2.03	0.58
1:A:252:ARG:NH1	1:A:414:THR:O	2.34	0.58
1:B:184:LEU:HD22	1:B:184:LEU:N	2.18	0.58
1:A:220:THR:HB	1:A:423:ASP:OD1	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:LYS:NZ	1:B:358:ASN:HB3	2.18	0.57
1:B:221:LYS:HZ2	1:B:424:GLU:CD	2.12	0.57
1:A:384:GLU:HA	1:A:391:LYS:HD2	1.87	0.56
1:A:282:ARG:O	1:A:295:HIS:ND1	2.39	0.56
1:B:224:THR:HG21	1:B:400:MET:H	1.71	0.56
1:B:42:PRO:HA	1:B:152:MET:O	2.05	0.56
1:A:210:ALA:HB1	1:A:365:GLU:HB2	1.88	0.55
1:B:209:PRO:HA	1:B:361:GLN:CG	2.17	0.55
1:A:98:VAL:HG13	1:A:143:PHE:HA	1.88	0.55
1:B:241:VAL:HG13	1:B:323:HIS:HB3	1.88	0.55
1:B:339:ASN:O	1:B:343:ILE:HG13	2.07	0.54
1:A:128:ILE:CD1	1:B:245:GLU:OE1	2.56	0.54
1:B:184:LEU:N	1:B:184:LEU:CD2	2.71	0.54
1:B:79:GLN:HE21	1:B:129:HIS:CE1	2.25	0.53
1:A:276:TYR:HA	1:A:306:VAL:O	2.07	0.53
1:B:30:THR:HG21	1:B:194:LEU:HB3	1.90	0.53
1:B:159:ALA:HB1	1:B:163:THR:HB	1.90	0.53
1:B:257:THR:HG22	1:B:407:LEU:HD11	1.91	0.53
1:B:422:MET:HG2	1:B:427:MET:HE2	1.90	0.53
1:B:276:TYR:HA	1:B:306:VAL:O	2.09	0.53
1:A:38:PRO:HD3	1:B:245:GLU:H	1.74	0.52
1:B:78:VAL:HG11	1:B:83:TYR:CZ	2.44	0.52
1:B:125:PHE:HD2	1:B:133:ASN:HA	1.75	0.51
1:A:126:LYS:HD3	1:A:127:ASP:N	2.25	0.51
1:B:310:PHE:CE2	1:B:320:LEU:HD21	2.45	0.51
1:B:282:ARG:O	1:B:295:HIS:ND1	2.44	0.50
1:B:311:ARG:NH2	1:B:363:ASP:OD1	2.45	0.50
1:B:269:LYS:NZ	2:B:506:HOH:O	2.44	0.50
1:B:224:THR:HG23	1:B:399:GLU:HG3	1.92	0.50
1:A:198:VAL:HG23	1:A:198:VAL:O	2.11	0.50
1:B:183:GLN:C	1:B:185:VAL:H	2.20	0.50
1:A:42:PRO:HB3	1:A:153:TYR:HD1	1.77	0.50
1:B:194:LEU:CD2	1:B:198:VAL:HG22	2.40	0.50
1:A:159:ALA:HB1	1:A:163:THR:HB	1.93	0.49
1:A:248:ASN:O	1:A:330:ARG:NH2	2.41	0.49
1:A:440:ALA:HB1	1:A:442:LEU:HG	1.94	0.49
1:A:288:ASN:N	1:A:288:ASN:OD1	2.45	0.49
1:A:218:LYS:O	1:A:220:THR:N	2.42	0.49
1:B:403:GLU:HG2	1:B:422:MET:HA	1.92	0.49
1:B:239:ASP:CG	1:B:249:ARG:HH21	2.21	0.49
1:B:184:LEU:CD2	1:B:184:LEU:H	2.26	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:LYS:HD2	1:B:249:ARG:HH12	1.78	0.48
1:B:79:GLN:CD	1:B:128:ILE:O	2.56	0.48
1:B:80:ASP:OD2	1:B:80:ASP:N	2.46	0.48
1:A:212:MET:SD	1:A:213:PRO:HD2	2.52	0.48
1:A:197:LEU:CD1	1:A:292:LEU:HD23	2.43	0.48
1:A:290:PRO:HG2	1:A:291:PRO:HD3	1.96	0.48
1:B:198:VAL:HG23	1:B:198:VAL:O	2.13	0.48
1:B:293:ASP:OD2	1:B:294:ASN:N	2.47	0.48
1:B:207:ILE:HA	1:B:358:ASN:OD1	2.14	0.47
1:A:253:VAL:HG22	1:A:379:ARG:HH21	1.80	0.47
1:B:30:THR:HB	1:B:198:VAL:HG23	1.97	0.47
1:B:42:PRO:HB3	1:B:153:TYR:HD1	1.80	0.47
1:B:79:GLN:HE21	1:B:129:HIS:HE1	1.63	0.47
1:A:338:LYS:O	1:A:341:SER:OG	2.29	0.47
1:B:343:ILE:HD12	1:B:349:LEU:HD23	1.95	0.47
1:A:44:ILE:HD12	1:A:151:CYS:HB2	1.96	0.47
1:B:320:LEU:HD23	1:B:320:LEU:HA	1.78	0.47
1:A:247:VAL:HG12	1:A:330:ARG:NH2	2.29	0.46
1:B:268:SER:O	1:B:269:LYS:HB2	2.15	0.46
1:A:129:HIS:CE1	1:B:252:ARG:CG	2.78	0.46
1:B:71:TYR:CD1	1:B:72:PRO:HD3	2.50	0.46
1:B:71:TYR:CG	1:B:72:PRO:HD3	2.50	0.46
1:B:338:LYS:O	1:B:341:SER:HB3	2.16	0.46
1:A:128:ILE:O	1:A:128:ILE:HG22	2.16	0.46
1:A:76:ARG:NH1	1:A:90:GLY:O	2.49	0.46
1:A:197:LEU:HD12	1:A:292:LEU:CD2	2.46	0.45
1:A:210:ALA:O	1:A:211:VAL:HB	2.17	0.45
1:A:197:LEU:HD11	1:A:291:PRO:HD2	1.99	0.45
1:A:109:ALA:O	1:A:113:ILE:HD12	2.16	0.45
1:B:218:LYS:HA	1:B:218:LYS:HD2	1.69	0.45
1:B:129:HIS:O	2:B:501:HOH:O	2.21	0.44
1:B:224:THR:HG22	1:B:421:SER:CB	2.47	0.44
1:B:101:PRO:CD	1:B:104:GLU:OE1	2.61	0.44
1:B:38:PRO:HB3	1:B:128:ILE:HD13	1.99	0.44
1:B:211:VAL:HG22	1:B:212:MET:N	2.33	0.44
1:A:440:ALA:CB	1:A:442:LEU:HG	2.47	0.44
1:A:4:MET:HE3	1:A:98:VAL:HG21	2.00	0.44
1:B:433:ASP:HB3	1:B:436:ILE:HG12	1.99	0.44
1:A:42:PRO:HA	1:A:152:MET:O	2.18	0.44
1:A:376:THR:HA	1:A:407:LEU:HB2	2.00	0.43
1:B:51:ASN:HD22	1:B:51:ASN:HA	1.68	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:LYS:O	1:B:130:SER:HB3	2.18	0.43
1:B:257:THR:HG22	1:B:407:LEU:CD1	2.48	0.43
1:A:243:LYS:HB3	1:A:243:LYS:HE2	1.28	0.43
1:B:128:ILE:HD12	1:B:128:ILE:HA	1.87	0.43
1:B:246:SER:O	1:B:247:VAL:HG12	2.18	0.43
1:A:422:MET:HG2	1:A:427:MET:HE2	2.01	0.43
1:A:32:PHE:CE2	1:A:198:VAL:HG21	2.53	0.43
1:B:224:THR:HA	1:B:420:VAL:O	2.19	0.43
1:A:126:LYS:HD3	1:A:127:ASP:H	1.84	0.42
1:B:110:LYS:HG2	1:B:394:TRP:CZ2	2.54	0.42
1:B:211:VAL:O	1:B:212:MET:HG2	2.19	0.42
1:A:154:LEU:HD23	1:A:164:ALA:HB2	2.02	0.42
1:B:394:TRP:CZ2	1:B:396:THR:HG21	2.54	0.42
1:A:223:VAL:O	1:A:421:SER:HA	2.20	0.41
1:B:109:ALA:O	1:B:113:ILE:HD13	2.21	0.41
1:A:124:ASN:ND2	2:A:501:HOH:O	2.22	0.41
1:B:224:THR:HG22	1:B:421:SER:HB3	2.01	0.41
1:B:154:LEU:HD12	1:B:154:LEU:HA	1.76	0.41
1:A:129:HIS:CE1	1:B:252:ARG:H	2.34	0.41
1:B:77:PHE:CE2	1:B:79:GLN:OE1	2.67	0.41
1:A:42:PRO:HB3	1:A:153:TYR:CD1	2.55	0.41
1:A:154:LEU:HD12	1:A:155:SER:H	1.86	0.41
1:A:242:ILE:O	1:A:245:GLU:N	2.49	0.41
1:B:207:ILE:O	1:B:207:ILE:HG13	2.21	0.41
1:A:148:LEU:HD23	1:A:148:LEU:HA	1.94	0.41
1:B:203:LEU:HD23	1:B:206:VAL:HG21	2.03	0.41
1:B:288:ASN:OD1	1:B:288:ASN:N	2.53	0.41
1:B:351:LEU:HD23	1:B:351:LEU:HA	1.89	0.41
1:A:71:TYR:CG	1:A:72:PRO:HD3	2.56	0.40
1:B:197:LEU:HD12	1:B:292:LEU:HD12	2.03	0.40
1:B:226:ARG:HE	1:B:417:GLU:CD	2.30	0.40
1:B:11:LYS:HB3	1:B:11:LYS:HE3	1.81	0.40
1:A:69:LYS:O	1:A:190:ILE:HD11	2.22	0.40
1:A:289:THR:O	1:A:290:PRO:C	2.64	0.40
1:A:42:PRO:O	1:A:397:ILE:HG23	2.21	0.40
1:A:272:SER:OG	1:A:311:ARG:HD3	2.21	0.40
1:B:76:ARG:NH2	1:B:89:GLU:OE1	2.54	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	438/449 (98%)	414 (94%)	21 (5%)	3 (1%)	18	21
1	B	438/449 (98%)	418 (95%)	17 (4%)	3 (1%)	18	21
All	All	876/898 (98%)	832 (95%)	38 (4%)	6 (1%)	18	21

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	211	VAL
1	A	129	HIS
1	B	210	ALA
1	A	130	SER
1	B	211	VAL
1	B	307	PRO

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	396/405 (98%)	393 (99%)	3 (1%)	73	81
1	B	394/405 (97%)	391 (99%)	3 (1%)	73	81
All	All	790/810 (98%)	784 (99%)	6 (1%)	73	81

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	ILE
1	A	130	SER
1	A	205	THR
1	B	184	LEU
1	B	208	LYS
1	B	222	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	40	GLN
1	A	124	ASN
1	A	248	ASN
1	A	294	ASN
1	B	51	ASN
1	B	79	GLN
1	B	183	GLN
1	B	401	HIS

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 ⓘ

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	440/449 (97%)	0.43	28 (6%) 25 22	33, 52, 99, 143	0
1	B	440/449 (97%)	0.56	45 (10%) 12 9	36, 58, 116, 154	0
All	All	880/898 (97%)	0.49	73 (8%) 17 14	33, 55, 109, 154	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	129	HIS	6.1
1	A	211	VAL	4.6
1	A	128	ILE	4.5
1	B	245	GLU	4.4
1	A	442	LEU	4.4
1	B	125	PHE	4.2
1	B	217	ILE	4.2
1	B	242	ILE	4.2
1	B	41	VAL	4.2
1	B	184	LEU	4.2
1	B	189	PRO	4.0
1	A	217	ILE	3.8
1	B	442	LEU	3.8
1	A	81	GLY	3.8
1	B	247	VAL	3.7
1	A	290	PRO	3.6
1	B	124	ASN	3.6
1	A	185	VAL	3.5
1	B	211	VAL	3.4
1	B	313	GLY	3.4
1	A	3	ILE	3.4
1	B	210	ALA	3.4
1	B	244	SER	3.4
1	A	247	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	209	PRO	3.2
1	B	129	HIS	3.1
1	B	127	ASP	3.1
1	B	191	ASN	3.1
1	B	123	LYS	3.0
1	B	126	LYS	3.0
1	A	246	SER	3.0
1	B	312	GLY	3.0
1	B	128	ILE	2.9
1	B	185	VAL	2.9
1	B	50	SER	2.9
1	B	283	GLY	2.9
1	A	215	SER	2.8
1	B	314	ASN	2.8
1	B	246	SER	2.7
1	A	210	ALA	2.7
1	A	212	MET	2.7
1	A	312	GLY	2.6
1	B	181	GLY	2.6
1	A	296	PHE	2.6
1	A	125	PHE	2.6
1	B	212	MET	2.5
1	A	122	LYS	2.5
1	A	191	ASN	2.5
1	B	53	HIS	2.5
1	A	218	LYS	2.5
1	B	79	GLN	2.4
1	B	290	PRO	2.4
1	B	153	TYR	2.4
1	B	296	PHE	2.4
1	B	209	PRO	2.4
1	A	208	LYS	2.4
1	B	104	GLU	2.4
1	A	213	PRO	2.4
1	B	203	LEU	2.4
1	B	154	LEU	2.3
1	A	80	ASP	2.3
1	B	315	GLN	2.3
1	B	186	SER	2.3
1	B	81	GLY	2.2
1	B	214	PRO	2.2
1	A	243	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	239	ASP	2.2
1	B	213	PRO	2.2
1	A	220	THR	2.1
1	B	179	ILE	2.1
1	B	288	ASN	2.1
1	A	291	PRO	2.0
1	A	244	SER	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.