



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2023 – 05:11 PM JST

PDB ID : 7W26
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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

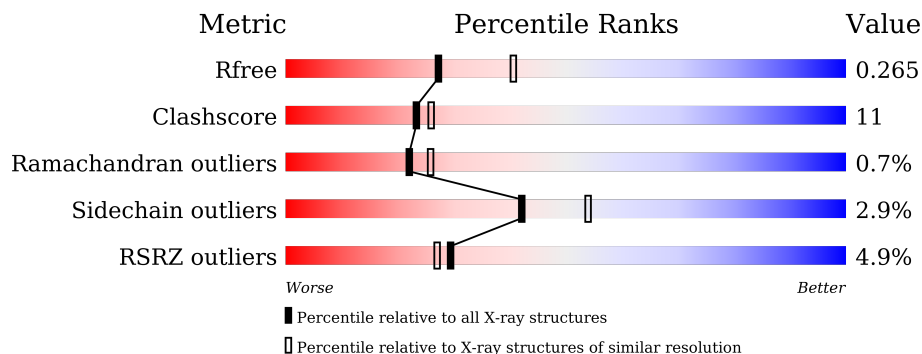
1 Overall quality at a glance

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}	130704	1564 (2.46-2.42)
Clashscore	141614	1631 (2.46-2.42)
Ramachandran outliers	138981	1617 (2.46-2.42)
Sidechain outliers	138945	1617 (2.46-2.42)
RSRZ outliers	127900	1547 (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	 4% (Poor fit) 78% (0 outliers) 18% (1 outlier) .. (2+ outliers)
1	B	449	 6% (Poor fit) 73% (0 outliers) 24% (1 outlier) .. (2+ outliers)

2 Entry composition i

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
			Total	C	N	O	S			
1	A	440	3489	2237	578	657	17	0	0	0
1	B	440	3482	2232	578	655	17	0	0	0

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

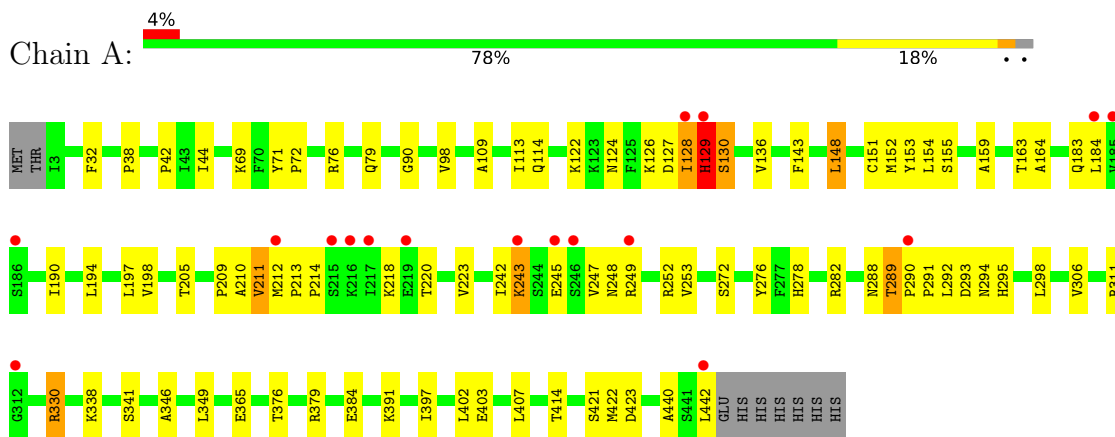
- Molecule 2 is water.

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

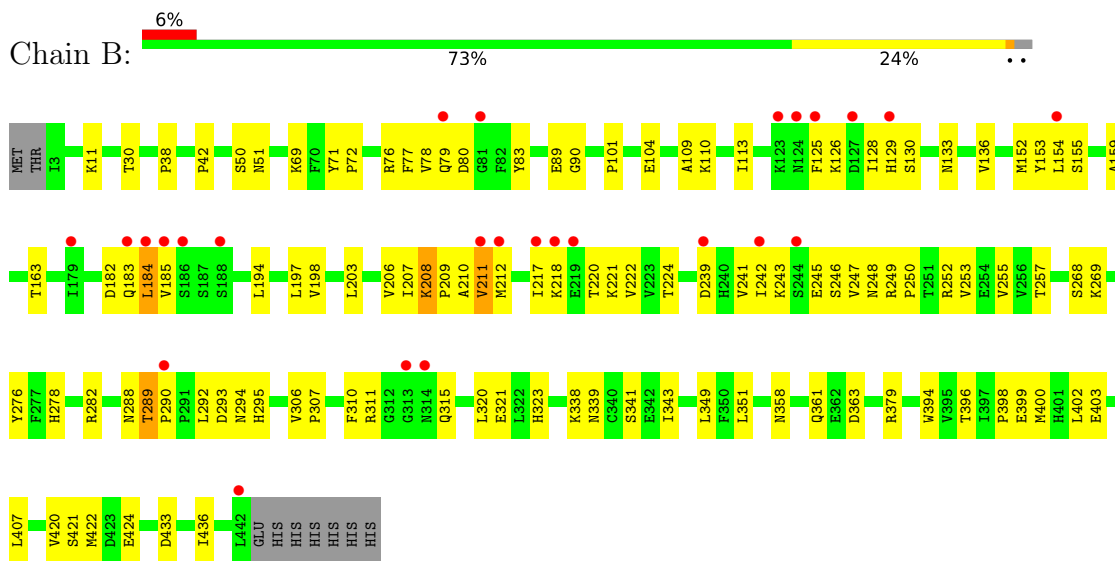
3 Residue-property plots

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) 99.8 (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)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.216 , 0.266 0.218 , 0.265	Depositor DCC
R_{free} test set	1847 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	50.7	Xtrriage
Anisotropy	0.097	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
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

Xtrriage'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 [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	1/3567 (0.0%)	0.61	4/4845 (0.1%)
1	B	0.33	0/3560	0.58	0/4836
All	All	0.35	1/7127 (0.0%)	0.59	4/9681 (0.0%)

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 (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	243	LYS	CD-CE	5.02	1.63	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	249	ARG	CG-CD-NE	7.46	127.47	111.80
1	A	129	HIS	CB-CA-C	-6.95	96.50	110.40
1	A	129	HIS	N-CA-CB	5.44	120.39	110.60
1	A	243	LYS	CB-CA-C	-5.05	100.29	110.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	289	THR	Peptide

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Mol	Chain	Res	Type	Group
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	70	0
1	B	3482	0	3476	102	0
2	A	54	0	0	1	0
2	B	41	0	0	1	0
All	All	7066	0	6968	158	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 11.

All (158) 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
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: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:A:128:ILE:CD1	1:B:245:GLU:CD	2.51	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:LYS:NZ	1:B:424:GLU:CD	2.44	0.71
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:128:ILE:HD13	1:B:245:GLU:CD	2.17	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: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:128:ILE:HD11	1:B:245:GLU:CD	2.22	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
1:B:208:LYS:NZ	1:B:358:ASN:HB3	2.18	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:221:LYS:HZ3	1:B:424:GLU:CD	2.11	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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: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:184:LEU:CD2	1:B:184:LEU:H	2.26	0.48
1:A:126:LYS:HD2	1:B:249:ARG:HH12	1.78	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:128:ILE:O	1:A:128:ILE:HG22	2.16	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: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:B:183:GLN:C	1:B:185:VAL:H	2.20	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:239:ASP:CG	1:B:249:ARG:HH21	2.21	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: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
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: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:128:ILE:O	1:A:128:ILE:CG2	2.68	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HG13	1:B:207:ILE:O	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: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: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 [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	438/449 (98%)	414 (94%)	21 (5%)	3 (1%)	22	26
1	B	438/449 (98%)	418 (95%)	17 (4%)	3 (1%)	22	26
All	All	876/898 (98%)	832 (95%)	38 (4%)	6 (1%)	22	26

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	211	VAL

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Mol	Chain	Res	Type
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 [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	396/405 (98%)	385 (97%)	11 (3%)	43 56
1	B	394/405 (97%)	382 (97%)	12 (3%)	41 53
All	All	790/810 (98%)	767 (97%)	23 (3%)	42 54

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	114	GLN
1	A	122	LYS
1	A	128	ILE
1	A	130	SER
1	A	148	LEU
1	A	205	THR
1	A	209	PRO
1	A	278	HIS
1	A	330	ARG
1	A	402	LEU
1	B	50	SER
1	B	69	LYS
1	B	155	SER
1	B	182	ASP
1	B	184	LEU
1	B	208	LYS
1	B	222	VAL

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Mol	Chain	Res	Type
1	B	243	LYS
1	B	278	HIS
1	B	315	GLN
1	B	379	ARG
1	B	402	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	A	124	ASN
1	B	51	ASN
1	B	79	GLN
1	B	183	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no 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	440/449 (97%)	0.20	17 (3%) 39 36	33, 52, 99, 143	0
1	B	440/449 (97%)	0.32	26 (5%) 22 18	36, 58, 116, 154	0
All	All	880/898 (97%)	0.26	43 (4%) 29 27	33, 55, 109, 154	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	186	SER	12.0
1	B	125	PHE	7.8
1	B	123	LYS	6.9
1	B	185	VAL	6.3
1	B	184	LEU	6.3
1	A	129	HIS	6.2
1	A	442	LEU	6.2
1	A	186	SER	5.7
1	B	442	LEU	5.6
1	B	217	ILE	4.5
1	B	127	ASP	4.5
1	A	185	VAL	4.0
1	B	124	ASN	4.0
1	B	218	LYS	3.9
1	B	219	GLU	3.9
1	B	290	PRO	3.8
1	A	128	ILE	3.7
1	B	183	GLN	3.7
1	A	215	SER	3.4
1	B	129	HIS	3.3
1	A	290	PRO	3.1
1	B	314	ASN	3.1
1	A	219	GLU	3.1
1	A	249	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	184	LEU	2.8
1	B	244	SER	2.6
1	A	216	LYS	2.5
1	B	79	GLN	2.4
1	A	212	MET	2.4
1	B	211	VAL	2.3
1	A	217	ILE	2.3
1	A	245	GLU	2.3
1	B	188	SER	2.3
1	B	242	ILE	2.3
1	A	246	SER	2.3
1	B	313	GLY	2.3
1	A	243	LYS	2.2
1	B	154	LEU	2.2
1	B	81	GLY	2.1
1	B	239	ASP	2.1
1	B	179	ILE	2.1
1	B	212	MET	2.1
1	A	312	GLY	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.