



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2026 – 01:14 PM JST

PDB ID : 9WAR / pdb_00009war
Title : Open conformation of the EPSPS tetramer from *Amaranthus retroflexus*
Authors : Liu, X.X.; Zhao, K.H.; Luo, X.; Lu, T.T.; Li, X.Y.
Deposited on : 2025-08-12
Resolution : 2.08 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

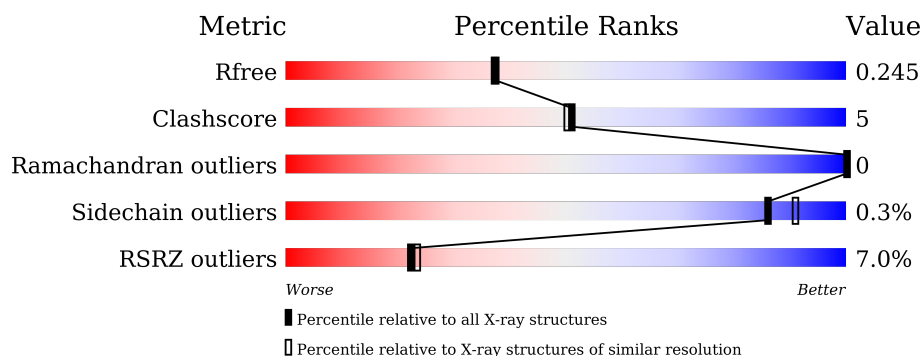
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	454	2% 87% 11% .
1	B	454	9% 84% 13% .
1	C	454	5% 89% 8% .
1	D	454	10% 81% 15% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13515 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 5-enolpyruvylshikimate-3-phosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	444	Total	C	N	O	S	0	0	0
			3329	2099	562	646	22			
1	B	440	Total	C	N	O	S	0	0	0
			3294	2079	555	638	22			
1	C	441	Total	C	N	O	S	0	0	0
			3302	2085	556	639	22			
1	D	442	Total	C	N	O	S	0	0	0
			3310	2089	558	641	22			

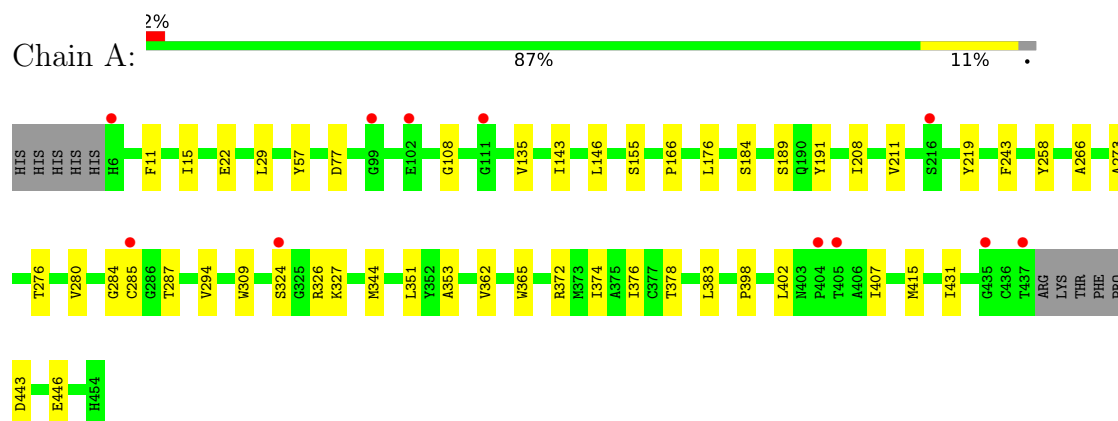
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	82	Total	O	0	0
			82	82		
2	B	58	Total	O	0	0
			58	58		
2	C	119	Total	O	0	0
			119	119		
2	D	21	Total	O	0	0
			21	21		

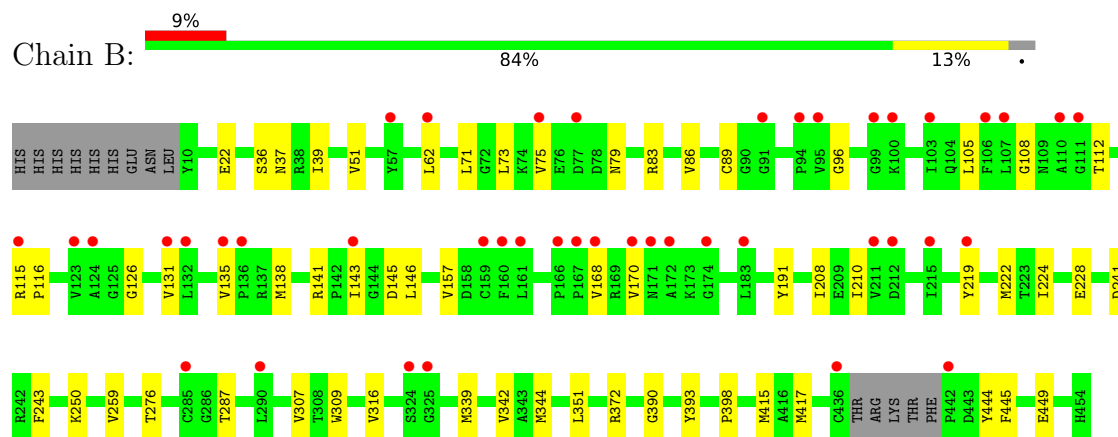
3 Residue-property plots

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.

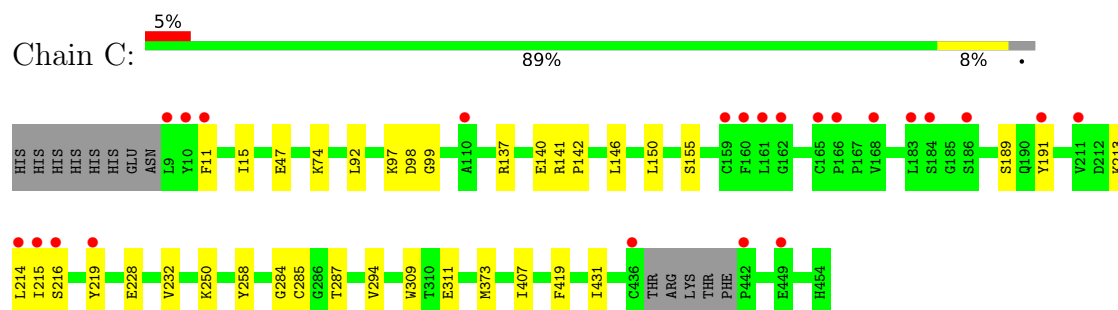
- Molecule 1: 5-enolpyruvylshikimate-3-phosphate synthase



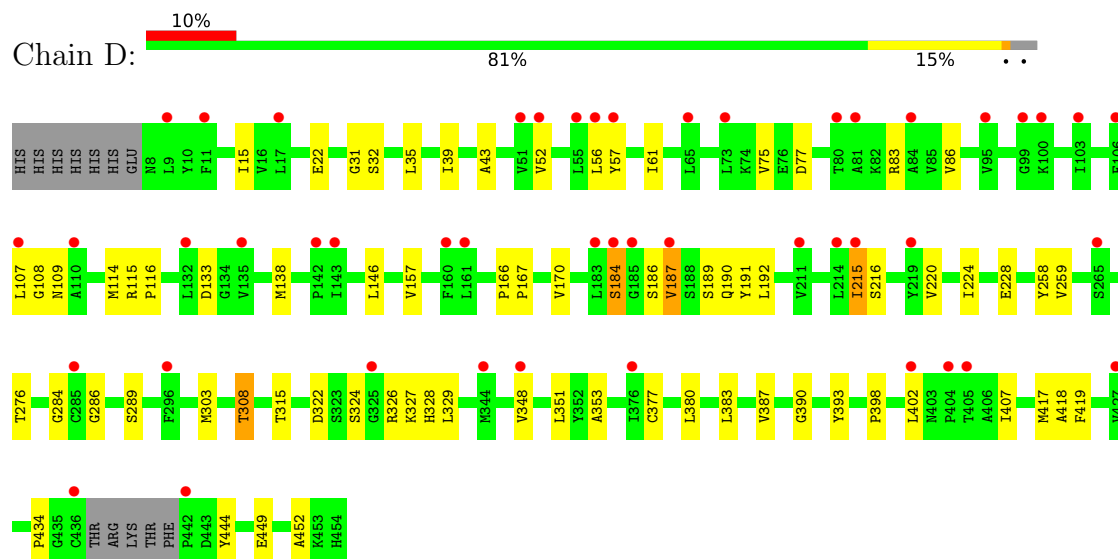
- Molecule 1: 5-enolpyruvylshikimate-3-phosphate synthase



- Molecule 1: 5-enolpyruvylshikimate-3-phosphate synthase



Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	162.86Å 120.00Å 93.02Å 90.00° 91.25° 90.00°	Depositor
Resolution (Å)	40.76 – 2.08 40.76 – 2.08	Depositor EDS
% Data completeness (in resolution range)	98.3 (40.76-2.08) 98.4 (40.76-2.08)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.08Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.209 , 0.245 0.209 , 0.245	Depositor DCC
R_{free} test set	5074 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	45.4	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13515	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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	0.13	0/3384	0.32	0/4587
1	B	0.16	0/3349	0.33	0/4539
1	C	0.15	0/3357	0.32	0/4550
1	D	0.19	0/3365	0.36	0/4561
All	All	0.16	0/13455	0.33	0/18237

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3329	0	3362	30	0
1	B	3294	0	3333	37	0
1	C	3302	0	3344	20	0
1	D	3310	0	3350	52	0
2	A	82	0	0	0	0
2	B	58	0	0	0	0
2	C	119	0	0	0	0
2	D	21	0	0	0	0
All	All	13515	0	13389	137	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 5.

All (137) 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:143:ILE:HD11	1:B:168:VAL:HG21	1.76	0.67
1:D:114:MET:HB3	1:D:115:ARG:HH12	1.63	0.64
1:A:383:LEU:HD21	1:A:407:ILE:HD11	1.81	0.61
1:B:108:GLY:HA2	1:B:135:VAL:HG13	1.84	0.60
1:B:344:MET:HE1	1:B:415:MET:HE2	1.84	0.60
1:D:133:ASP:OD2	1:D:167:PRO:HB3	2.02	0.59
1:C:11:PHE:CD1	1:D:449:GLU:HG3	2.37	0.59
1:D:322:ASP:H	1:D:328:HIS:CE1	2.20	0.59
1:B:344:MET:SD	1:B:372:ARG:HD3	2.43	0.58
1:A:11:PHE:CD1	1:B:449:GLU:HG3	2.38	0.58
1:D:57:TYR:OH	1:D:77:ASP:HB2	2.03	0.58
1:D:351:LEU:HD13	1:D:402:LEU:HD11	1.85	0.58
1:A:383:LEU:HB3	1:A:402:LEU:HD22	1.86	0.57
1:B:157:VAL:HG13	1:B:170:VAL:HG22	1.87	0.57
1:D:187:VAL:HG11	1:D:191:TYR:HD2	1.70	0.57
1:D:189:SER:HB3	1:D:216:SER:CB	2.35	0.57
1:A:15:ILE:HG13	1:A:431:ILE:HB	1.88	0.56
1:A:57:TYR:OH	1:A:77:ASP:HB2	2.06	0.56
1:D:56:LEU:O	1:D:61:ILE:HG13	2.06	0.56
1:B:307:VAL:HG22	1:B:316:VAL:HG22	1.88	0.54
1:C:311:GLU:H	1:C:311:GLU:CD	2.15	0.54
1:D:146:LEU:HA	1:D:191:TYR:CE1	2.43	0.54
1:B:210:ILE:HB	1:B:241:ASP:HA	1.90	0.54
1:C:47:GLU:OE2	1:C:250:LYS:HD3	2.08	0.53
1:C:189:SER:HB2	1:C:219:TYR:CD2	2.43	0.53
1:D:114:MET:HE3	1:D:138:MET:HE2	1.90	0.53
1:A:258:TYR:HE1	1:A:284:GLY:HA2	1.74	0.53
1:D:351:LEU:HA	1:D:398:PRO:HB3	1.90	0.53
1:B:417:MET:HE3	1:B:444:TYR:HB2	1.91	0.52
1:D:114:MET:HB3	1:D:115:ARG:NH1	2.25	0.51
1:A:376:ILE:HD13	1:A:415:MET:HE3	1.93	0.51
1:C:373:MET:HA	1:C:373:MET:HE2	1.92	0.51
1:B:37:ASN:HD22	1:B:112:THR:HB	1.75	0.51
1:A:108:GLY:HA2	1:A:135:VAL:HG13	1.93	0.50
1:C:92:LEU:HD21	1:C:97:LYS:HG3	1.92	0.50
1:A:184:SER:HA	1:A:211:VAL:HG22	1.94	0.50
1:D:39:ILE:HD12	1:D:259:VAL:HG22	1.94	0.50
1:B:73:LEU:HD11	1:B:89:CYS:HB3	1.93	0.50
1:C:285:CYS:SG	1:C:294:VAL:HG23	2.52	0.49
1:A:155:SER:HB2	1:A:176:LEU:HD12	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:326:ARG:HG3	1:D:328:HIS:CE1	2.46	0.49
1:D:43:ALA:HB1	1:D:86:VAL:HG21	1.94	0.49
1:B:71:LEU:HD21	1:B:105:LEU:HD11	1.95	0.48
1:C:213:LYS:HD3	1:C:214:LEU:N	2.28	0.48
1:D:308:THR:HG23	1:D:315:THR:HB	1.95	0.48
1:C:137:ARG:O	1:C:140:GLU:HG2	2.13	0.48
1:D:15:ILE:HG13	1:D:434:PRO:HG3	1.96	0.48
1:B:22:GLU:HA	1:B:276:THR:HB	1.96	0.48
1:B:219:TYR:HA	1:B:222:MET:HE2	1.94	0.48
1:C:407:ILE:HD12	1:C:419:PHE:CG	2.49	0.48
1:D:258:TYR:HE1	1:D:284:GLY:HA2	1.79	0.48
1:B:208:ILE:HB	1:B:243:PHE:HB2	1.96	0.48
1:D:186:SER:HA	1:D:215:ILE:HD13	1.95	0.48
1:D:115:ARG:HB2	1:D:116:PRO:HD3	1.96	0.47
1:D:108:GLY:O	1:D:138:MET:HG3	2.15	0.47
1:D:322:ASP:HB2	1:D:328:HIS:CE1	2.50	0.47
1:A:324:SER:HB3	1:A:326:ARG:HG2	1.96	0.47
1:B:339:MET:HB2	1:B:342:VAL:HG22	1.96	0.47
1:A:327:LYS:HD2	1:A:353:ALA:O	2.15	0.47
1:D:32:SER:OG	1:D:35:LEU:HB2	2.15	0.46
1:B:75:VAL:HG23	1:B:86:VAL:HG22	1.97	0.46
1:A:143:ILE:HG12	1:A:166:PRO:HD3	1.97	0.46
1:D:303:MET:HE2	1:D:329:LEU:HB3	1.97	0.46
1:A:258:TYR:CE1	1:A:284:GLY:HA2	2.49	0.46
1:A:146:LEU:HA	1:A:191:TYR:CE1	2.51	0.46
1:A:374:ILE:O	1:A:378:THR:HG23	2.15	0.46
1:D:377:CYS:HB3	1:D:387:VAL:HG11	1.98	0.46
1:A:189:SER:HB2	1:A:219:TYR:CD2	2.51	0.45
1:D:383:LEU:HD22	1:D:402:LEU:HD12	1.97	0.45
1:D:107:LEU:HD12	1:D:114:MET:HE2	1.98	0.45
1:B:145:ASP:HB3	1:B:191:TYR:CD2	2.51	0.45
1:B:146:LEU:HA	1:B:191:TYR:CE1	2.51	0.45
1:B:390:GLY:HA3	1:B:393:TYR:CE2	2.50	0.45
1:D:258:TYR:CE1	1:D:284:GLY:HA2	2.52	0.45
1:D:417:MET:HE3	1:D:444:TYR:HB2	1.97	0.45
1:B:351:LEU:HA	1:B:398:PRO:HB3	1.99	0.45
1:D:377:CYS:HA	1:D:380:LEU:HD12	1.99	0.45
1:B:96:GLY:O	1:B:126:GLY:HA3	2.17	0.45
1:D:157:VAL:HG22	1:D:170:VAL:HG22	1.99	0.45
1:D:115:ARG:HD2	1:D:190:GLN:HB3	1.99	0.45
1:A:287:THR:HG23	1:A:309:TRP:HB3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:ILE:HD12	1:B:259:VAL:HG22	1.99	0.44
1:C:258:TYR:HE1	1:C:284:GLY:HA2	1.81	0.44
1:B:224:ILE:O	1:B:228:GLU:HG3	2.16	0.44
1:B:131:VAL:HA	1:B:168:VAL:O	2.17	0.44
1:B:138:MET:HE1	1:B:141:ARG:HH11	1.83	0.44
1:B:51:VAL:HG11	1:B:83:ARG:HE	1.83	0.44
1:D:138:MET:CE	1:D:166:PRO:HG2	2.48	0.44
1:A:362:VAL:HG13	1:A:365:TRP:CE2	2.53	0.43
1:D:52:VAL:O	1:D:83:ARG:HG3	2.18	0.43
1:A:22:GLU:HA	1:A:276:THR:HB	2.00	0.43
1:D:115:ARG:CZ	1:D:115:ARG:H	2.31	0.43
1:D:407:ILE:HD13	1:D:419:PHE:CD1	2.54	0.43
1:A:285:CYS:SG	1:A:294:VAL:HG23	2.58	0.43
1:D:324:SER:HB3	1:D:326:ARG:HG2	2.01	0.43
1:A:273:ALA:HB2	1:A:280:VAL:HG23	2.01	0.43
1:B:62:LEU:HD21	1:B:79:ASN:HD21	1.84	0.43
1:D:286:GLY:O	1:D:289:SER:HB3	2.18	0.43
1:B:157:VAL:HG22	1:B:170:VAL:HG13	1.99	0.43
1:B:250:LYS:HE2	1:B:250:LYS:HB3	1.79	0.43
1:D:390:GLY:HA3	1:D:393:TYR:CE1	2.53	0.43
1:A:443:ASP:HB3	1:A:446:GLU:HB2	2.00	0.43
1:B:51:VAL:CG1	1:B:83:ARG:HE	2.32	0.43
1:C:98:ASP:CG	1:C:99:GLY:H	2.27	0.43
1:D:75:VAL:HG12	1:D:86:VAL:HG12	2.01	0.43
1:B:36:SER:OG	1:B:259:VAL:HG11	2.19	0.43
1:A:344:MET:HG3	1:A:376:ILE:HG13	2.01	0.43
1:C:150:LEU:HB3	1:C:155:SER:OG	2.19	0.43
1:D:348:VAL:HG21	1:D:418:ALA:O	2.19	0.43
1:C:215:ILE:HG13	1:C:216:SER:OG	2.19	0.42
1:D:166:PRO:HA	1:D:167:PRO:C	2.44	0.42
1:D:449:GLU:HA	1:D:452:ALA:HB3	2.01	0.42
1:D:322:ASP:HB2	1:D:328:HIS:HE1	1.84	0.42
1:B:62:LEU:HD21	1:B:79:ASN:ND2	2.35	0.42
1:D:192:LEU:HD23	1:D:220:VAL:HG21	2.01	0.42
1:A:208:ILE:HB	1:A:243:PHE:HB2	2.01	0.41
1:D:224:ILE:O	1:D:228:GLU:HG3	2.20	0.41
1:A:189:SER:HB2	1:A:219:TYR:HD2	1.85	0.41
1:B:445:PHE:O	1:B:449:GLU:HG2	2.20	0.41
1:D:31:GLY:O	1:D:56:LEU:HB2	2.19	0.41
1:A:351:LEU:HA	1:A:398:PRO:HB3	2.02	0.41
1:C:15:ILE:CG2	1:C:431:ILE:HB	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:ARG:HA	1:C:142:PRO:HD3	1.96	0.41
1:C:287:THR:HG23	1:C:309:TRP:HB3	2.01	0.41
1:B:287:THR:HG23	1:B:309:TRP:HB3	2.02	0.41
1:B:143:ILE:HD11	1:B:168:VAL:CG2	2.48	0.41
1:C:74:LYS:HD2	1:C:74:LYS:HA	1.68	0.41
1:D:109:ASN:HA	1:D:138:MET:HG2	2.03	0.41
1:D:327:LYS:HD2	1:D:353:ALA:O	2.21	0.41
1:D:184:SER:O	1:D:187:VAL:HB	2.21	0.41
1:A:372:ARG:O	1:A:376:ILE:HG12	2.21	0.40
1:A:376:ILE:HD13	1:A:376:ILE:HA	1.90	0.40
1:A:29:LEU:HD11	1:A:266:ALA:HB2	2.04	0.40
1:C:146:LEU:HA	1:C:191:TYR:CE1	2.56	0.40
1:C:228:GLU:HG2	1:C:232:VAL:O	2.21	0.40
1:B:115:ARG:HB2	1:B:116:PRO:HD3	2.04	0.40
1:D:22:GLU:HA	1:D:276:THR:HB	2.03	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	440/454 (97%)	433 (98%)	7 (2%)	0	100	100
1	B	436/454 (96%)	427 (98%)	9 (2%)	0	100	100
1	C	437/454 (96%)	432 (99%)	5 (1%)	0	100	100
1	D	438/454 (96%)	431 (98%)	7 (2%)	0	100	100
All	All	1751/1816 (96%)	1723 (98%)	28 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	362/372 (97%)	362 (100%)	0	100	100
1	B	358/372 (96%)	358 (100%)	0	100	100
1	C	359/372 (96%)	359 (100%)	0	100	100
1	D	360/372 (97%)	356 (99%)	4 (1%)	65	73
All	All	1439/1488 (97%)	1435 (100%)	4 (0%)	86	91

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	184	SER
1	D	187	VAL
1	D	215	ILE
1	D	308	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	28	GLN
1	B	12	GLN
1	B	328	HIS
1	B	413	HIS
1	C	12	GLN
1	C	18	GLN
1	C	104	GLN
1	C	171	ASN
1	C	312	ASN
1	D	12	GLN
1	D	190	GLN
1	D	328	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 [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	444/454 (97%)	0.33	11 (2%) 58 61	28, 51, 88, 137	0
1	B	440/454 (96%)	0.70	43 (9%) 13 13	35, 58, 112, 191	0
1	C	441/454 (97%)	0.38	23 (5%) 33 34	26, 48, 101, 130	0
1	D	442/454 (97%)	0.91	47 (10%) 11 12	43, 73, 112, 184	0
All	All	1767/1816 (97%)	0.58	124 (7%) 22 23	26, 57, 106, 191	0

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	9	LEU	5.0
1	C	11	PHE	4.7
1	B	161	LEU	4.2
1	B	62	LEU	3.6
1	D	442	PRO	3.6
1	B	110	ALA	3.5
1	A	435	GLY	3.5
1	C	184	SER	3.4
1	C	215	ILE	3.4
1	C	442	PRO	3.2
1	B	211	VAL	3.1
1	D	285	CYS	3.1
1	C	211	VAL	3.1
1	D	135	VAL	3.1
1	D	95	VAL	3.1
1	D	160	PHE	3.0
1	B	135	VAL	3.0
1	B	143	ILE	3.0
1	D	103	ILE	3.0
1	D	106	PHE	3.0
1	D	187	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	131	VAL	2.9
1	B	103	ILE	2.9
1	B	57	TYR	2.9
1	D	52	VAL	2.8
1	D	184	SER	2.8
1	C	162	GLY	2.8
1	B	107	LEU	2.8
1	C	219	TYR	2.8
1	B	170	VAL	2.8
1	C	186	SER	2.8
1	D	185	GLY	2.8
1	D	99	GLY	2.7
1	B	436	CYS	2.7
1	B	77	ASP	2.7
1	D	436	CYS	2.7
1	B	111	GLY	2.7
1	D	9	LEU	2.6
1	D	73	LEU	2.6
1	D	183	LEU	2.6
1	D	143	ILE	2.6
1	C	110	ALA	2.6
1	D	161	LEU	2.6
1	D	376	ILE	2.6
1	D	325	GLY	2.6
1	D	56	LEU	2.6
1	D	11	PHE	2.6
1	D	84	ALA	2.6
1	D	211	VAL	2.5
1	B	172	ALA	2.5
1	D	405	THR	2.5
1	A	102	GLU	2.5
1	B	215	ILE	2.5
1	C	10	TYR	2.5
1	C	159	CYS	2.5
1	A	324	SER	2.5
1	C	214	LEU	2.5
1	D	404	PRO	2.4
1	B	168	VAL	2.4
1	D	348	VAL	2.4
1	B	159	CYS	2.4
1	D	402	LEU	2.4
1	A	216	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	165	CYS	2.4
1	C	183	LEU	2.4
1	A	404	PRO	2.4
1	B	285	CYS	2.4
1	B	442	PRO	2.4
1	D	65	LEU	2.3
1	B	91	GLY	2.3
1	B	324	SER	2.3
1	D	215	ILE	2.3
1	B	132	LEU	2.3
1	D	55	LEU	2.3
1	B	325	GLY	2.3
1	B	100	LYS	2.3
1	C	168	VAL	2.3
1	C	161	LEU	2.3
1	D	142	PRO	2.2
1	B	183	LEU	2.2
1	D	107	LEU	2.2
1	B	219	TYR	2.2
1	C	216	SER	2.2
1	B	95	VAL	2.2
1	B	167	PRO	2.2
1	D	427	VAL	2.2
1	D	110	ALA	2.2
1	B	290	LEU	2.2
1	D	214	LEU	2.2
1	D	265	SER	2.2
1	A	405	THR	2.2
1	D	296	PHE	2.2
1	D	344	MET	2.2
1	B	174	GLY	2.2
1	C	166	PRO	2.1
1	B	106	PHE	2.1
1	B	160	PHE	2.1
1	B	75	VAL	2.1
1	A	99	GLY	2.1
1	D	17	LEU	2.1
1	D	100	LYS	2.1
1	D	219	TYR	2.1
1	A	6	HIS	2.1
1	D	51	VAL	2.1
1	C	436	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	99	GLY	2.1
1	C	160	PHE	2.1
1	D	132	LEU	2.1
1	C	449	GLU	2.1
1	A	437	THR	2.1
1	B	166	PRO	2.1
1	A	111	GLY	2.1
1	B	123	VAL	2.1
1	A	285	CYS	2.0
1	B	94	PRO	2.0
1	B	136	PRO	2.0
1	D	80	THR	2.0
1	D	57	TYR	2.0
1	B	124	ALA	2.0
1	D	81	ALA	2.0
1	B	171	ASN	2.0
1	B	212	ASP	2.0
1	B	115	ARG	2.0
1	C	191	TYR	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.