



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:41 PM UTC

PDB ID : 4FJQ / pdb_00004fjq
Title : Crystal Structure of an alpha-Bisabolol synthase
Authors : Jianxu, L.; Peng, Z.
Deposited on : 2012-06-12
Resolution : 2.00 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

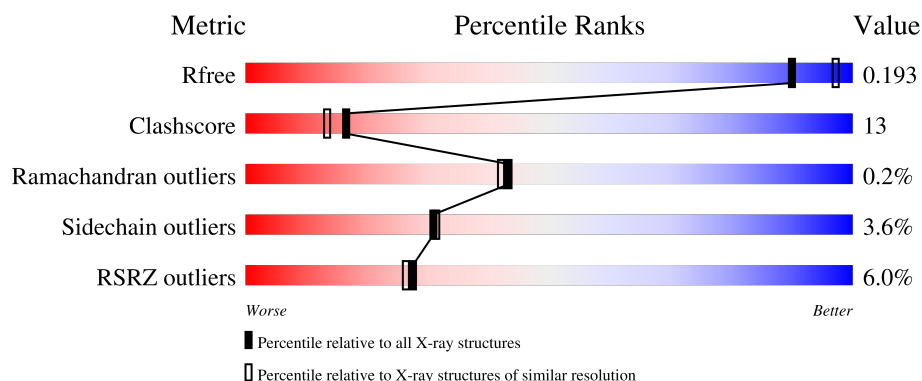
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	563	6% 70% 20% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4547 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 Amorpha-4,11-diene synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	521	Total	C	N	O	S	0	0	0
			4268	2762	698	791	17			

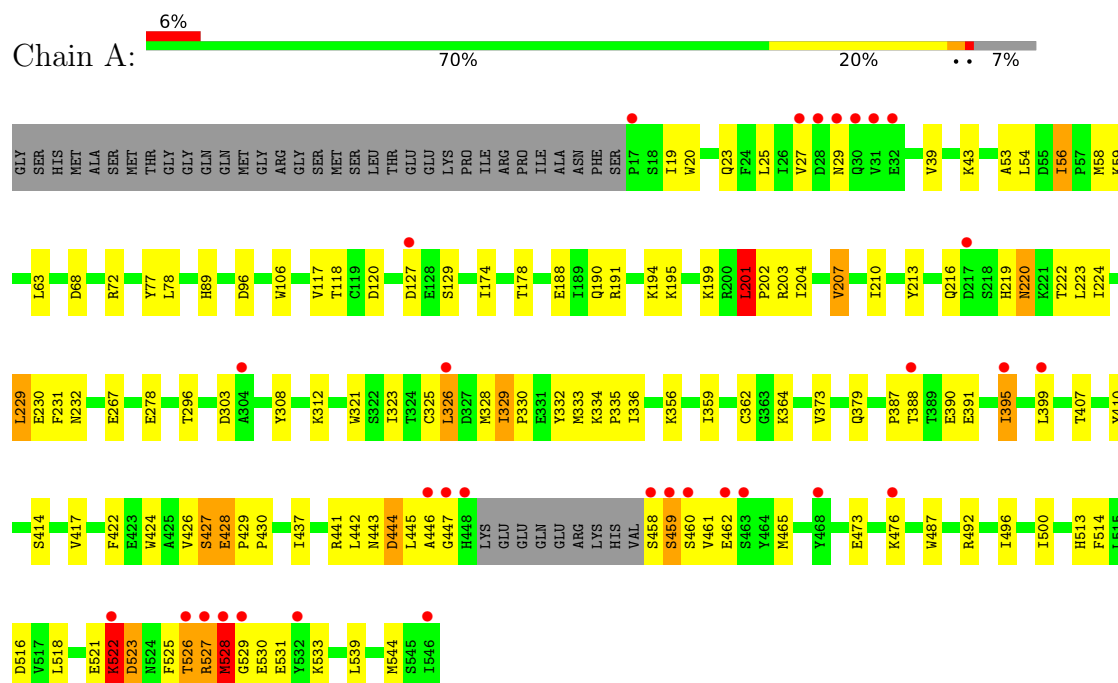
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	279	Total	O	0	0
			279	279		

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: Amorpha-4,11-diene synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.00Å 77.33Å 68.78Å 90.00° 102.50° 90.00°	Depositor
Resolution (Å)	39.22 – 2.00 39.22 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.5 (39.22-2.00) 97.1 (39.22-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.6.2_432	Depositor
R, R_{free}	0.181 , 0.223 (Not available) , 0.193	Depositor DCC
R_{free} test set	1754 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	21.9	Xtriage
Anisotropy	0.270	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.0	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	0.95	EDS
Total number of atoms	4547	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.07% 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	1.04	31/4362 (0.7%)	0.92	20/5904 (0.3%)

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	443	ASN	C-O	-15.18	1.06	1.24
1	A	430	PRO	C-O	-13.77	1.06	1.24
1	A	426	VAL	C-O	-13.48	1.07	1.24
1	A	527	ARG	CA-C	12.99	1.58	1.52
1	A	442	LEU	C-O	-12.66	1.09	1.24
1	A	429	PRO	C-O	-12.50	1.14	1.25
1	A	531	GLU	C-O	-12.42	1.09	1.24
1	A	529	GLY	C-O	-11.81	1.07	1.23
1	A	444	ASP	C-O	-10.90	1.11	1.24
1	A	528	MET	C-O	-10.70	1.10	1.24
1	A	428	GLU	C-O	-10.27	1.11	1.24
1	A	427	SER	CB-OG	-9.52	1.23	1.42
1	A	444	ASP	CA-C	-8.02	1.42	1.52
1	A	446	ALA	C-O	-7.86	1.13	1.24
1	A	523	ASP	C-O	-7.70	1.14	1.24
1	A	445	LEU	C-O	-7.62	1.14	1.24
1	A	530	GLU	CA-CB	-7.61	1.41	1.53
1	A	430	PRO	CB-CG	-7.57	1.11	1.49
1	A	429	PRO	CB-CG	-7.24	1.13	1.49
1	A	427	SER	C-O	-6.96	1.15	1.24
1	A	426	VAL	CB-CG2	-6.92	1.29	1.52
1	A	447	GLY	C-O	-6.90	1.14	1.23
1	A	530	GLU	C-O	-6.49	1.16	1.24
1	A	426	VAL	CA-CB	-6.44	1.45	1.54
1	A	530	GLU	C-N	-6.18	1.25	1.33
1	A	444	ASP	CG-OD2	-5.96	1.14	1.25
1	A	442	LEU	CG-CD1	-5.69	1.33	1.52
1	A	427	SER	CA-C	-5.63	1.44	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	522	LYS	CB-CG	-5.51	1.35	1.52
1	A	429	PRO	N-CA	-5.31	1.39	1.46
1	A	430	PRO	N-CD	-5.25	1.40	1.47

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	427	SER	CA-CB-OG	-14.03	83.05	111.10
1	A	530	GLU	N-CA-C	8.06	120.06	111.28
1	A	201	LEU	CA-C-N	7.10	126.62	119.24
1	A	201	LEU	C-N-CA	7.10	126.62	119.24
1	A	427	SER	N-CA-C	6.80	120.68	112.38
1	A	525	PHE	N-CA-C	6.43	117.95	111.07
1	A	531	GLU	CB-CA-C	6.42	121.45	110.79
1	A	329	ILE	N-CA-C	6.32	114.78	107.89
1	A	446	ALA	N-CA-C	5.93	120.53	112.88
1	A	522	LYS	CB-CG-CD	-5.64	98.33	111.30
1	A	529	GLY	CA-C-N	5.59	127.78	120.28
1	A	529	GLY	C-N-CA	5.59	127.78	120.28
1	A	459	SER	N-CA-C	5.47	117.24	111.28
1	A	428	GLU	CA-C-N	5.38	123.59	119.66
1	A	428	GLU	C-N-CA	5.38	123.59	119.66
1	A	443	ASN	N-CA-C	5.17	116.91	111.28
1	A	527	ARG	CA-C-N	5.08	131.24	121.54
1	A	527	ARG	C-N-CA	5.08	131.24	121.54
1	A	426	VAL	CA-C-N	5.05	128.69	120.60
1	A	426	VAL	C-N-CA	5.05	128.69	120.60

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	4268	0	4272	110	0
2	A	279	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4547	0	4272	110	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 13.

All (110) 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:A:462:GLU:HA	1:A:465:MET:HE2	1.46	0.97
1:A:188:GLU:OE1	1:A:191:ARG:HD3	1.67	0.95
1:A:118:THR:HG23	1:A:120:ASP:H	1.35	0.91
1:A:56:ILE:HD11	1:A:58:MET:HB2	1.53	0.90
1:A:232:ASN:HD21	1:A:544:MET:H	1.24	0.85
1:A:461:VAL:HG11	1:A:476:LYS:HE3	1.63	0.78
1:A:106:TRP:CH2	1:A:117:VAL:HG11	2.19	0.78
1:A:96:ASP:OD2	1:A:118:THR:HG22	1.83	0.77
1:A:523:ASP:CG	1:A:526:THR:HG23	2.10	0.76
1:A:207:VAL:HG12	1:A:231:PHE:CE1	2.25	0.72
1:A:424:TRP:O	1:A:427:SER:HB2	1.89	0.72
1:A:232:ASN:HD21	1:A:544:MET:N	1.86	0.72
1:A:53:ALA:HA	1:A:59:LYS:HE2	1.74	0.69
1:A:267:GLU:OE2	1:A:528:MET:HE2	1.95	0.67
1:A:19:ILE:HD11	1:A:528:MET:CE	2.24	0.67
1:A:513:HIS:HD2	2:A:637:HOH:O	1.77	0.67
1:A:308:TYR:CZ	1:A:312:LYS:HD2	2.29	0.66
1:A:462:GLU:HA	1:A:465:MET:CE	2.23	0.66
1:A:459:SER:HA	1:A:462:GLU:OE1	1.96	0.65
1:A:216:GLN:O	1:A:219:HIS:HD2	1.80	0.65
1:A:20:TRP:CZ2	1:A:528:MET:HE1	2.32	0.64
1:A:458:SER:HB3	1:A:461:VAL:HB	1.78	0.64
1:A:462:GLU:CA	1:A:465:MET:HE2	2.27	0.63
1:A:207:VAL:HG11	1:A:539:LEU:HD12	1.79	0.63
1:A:174:ILE:O	1:A:178:THR:HG23	1.99	0.62
1:A:461:VAL:O	1:A:465:MET:HG3	2.01	0.61
1:A:308:TYR:CE2	1:A:312:LYS:HD2	2.35	0.61
1:A:414:SER:O	1:A:417:VAL:HG22	2.04	0.58
1:A:220:ASN:HD22	1:A:223:LEU:H	1.52	0.58
1:A:388:THR:HG22	1:A:390:GLU:H	1.69	0.57
1:A:521:GLU:O	1:A:522:LYS:CD	2.53	0.57
1:A:188:GLU:HA	1:A:191:ARG:HG2	1.87	0.56
1:A:521:GLU:O	1:A:522:LYS:HD2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:LYS:NZ	1:A:487:TRP:CE2	2.74	0.56
1:A:188:GLU:CD	1:A:191:ARG:HD3	2.30	0.55
1:A:356:LYS:CD	1:A:414:SER:HB3	2.36	0.55
1:A:96:ASP:OD2	1:A:118:THR:CG2	2.53	0.54
1:A:207:VAL:HG12	1:A:231:PHE:HE1	1.71	0.54
1:A:19:ILE:CD1	1:A:528:MET:CE	2.85	0.54
1:A:321:TRP:CG	1:A:364:LYS:HE3	2.42	0.54
1:A:19:ILE:HD11	1:A:528:MET:HE1	1.89	0.54
1:A:527:ARG:O	1:A:528:MET:HB2	2.07	0.53
1:A:334:LYS:N	1:A:335:PRO:HD2	2.24	0.53
1:A:78:LEU:HD11	1:A:230:GLU:HG3	1.90	0.52
1:A:54:LEU:HD13	1:A:63:LEU:HD21	1.91	0.52
1:A:492:ARG:HD2	2:A:876:HOH:O	2.09	0.52
1:A:356:LYS:HD3	1:A:414:SER:HB3	1.90	0.52
1:A:19:ILE:CD1	1:A:528:MET:HE3	2.40	0.52
1:A:444:ASP:HB3	1:A:460:SER:HB2	1.90	0.52
1:A:332:TYR:OH	1:A:333:MET:HE3	2.10	0.52
1:A:513:HIS:HE1	2:A:630:HOH:O	1.93	0.52
1:A:523:ASP:OD1	1:A:526:THR:HG23	2.09	0.51
1:A:203:ARG:O	1:A:207:VAL:HG13	2.11	0.51
1:A:190:GLN:HG3	1:A:194:LYS:HE3	1.92	0.51
1:A:188:GLU:O	1:A:191:ARG:HG2	2.12	0.50
1:A:201:LEU:HD21	1:A:514:PHE:HD1	1.77	0.49
1:A:326:LEU:O	1:A:334:LYS:HE2	2.13	0.49
1:A:325:CYS:O	1:A:328:MET:HG2	2.13	0.49
1:A:387:PRO:HA	1:A:391:GLU:OE1	2.13	0.49
1:A:521:GLU:C	1:A:522:LYS:CD	2.86	0.49
1:A:118:THR:HG23	1:A:120:ASP:N	2.16	0.48
1:A:220:ASN:HD21	1:A:222:THR:HB	1.77	0.48
1:A:127:ASP:HB3	1:A:129:SER:H	1.79	0.48
1:A:106:TRP:CH2	1:A:117:VAL:CG1	2.95	0.47
1:A:207:VAL:HG12	1:A:231:PHE:CZ	2.50	0.47
1:A:308:TYR:OH	1:A:379:GLN:NE2	2.46	0.47
1:A:459:SER:CA	1:A:462:GLU:OE1	2.62	0.47
1:A:27:VAL:CG1	1:A:29:ASN:HD21	2.27	0.47
1:A:77:TYR:HA	2:A:669:HOH:O	2.15	0.47
1:A:77:TYR:HB3	1:A:278:GLU:HG2	1.96	0.47
1:A:195:LYS:NZ	1:A:516:ASP:OD2	2.48	0.46
1:A:359:ILE:HD13	1:A:417:VAL:CG2	2.45	0.46
1:A:428:GLU:OE2	2:A:787:HOH:O	2.21	0.46
1:A:23:GLN:HB3	2:A:878:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:459:SER:C	1:A:462:GLU:OE1	2.59	0.46
1:A:20:TRP:CE2	1:A:528:MET:HE1	2.51	0.46
1:A:216:GLN:O	1:A:219:HIS:CD2	2.65	0.45
1:A:201:LEU:HB2	1:A:204:ILE:HD12	1.98	0.45
1:A:39:VAL:O	1:A:43:LYS:HG3	2.16	0.45
1:A:56:ILE:O	1:A:56:ILE:HG13	2.16	0.45
1:A:210:ILE:HG23	1:A:224:ILE:HD11	1.98	0.45
1:A:68:ASP:O	1:A:72:ARG:HG3	2.17	0.45
1:A:201:LEU:HA	1:A:202:PRO:HD3	1.86	0.44
1:A:220:ASN:ND2	1:A:223:LEU:H	2.15	0.44
1:A:296:THR:HB	2:A:732:HOH:O	2.17	0.44
1:A:329:ILE:HG22	1:A:330:PRO:O	2.17	0.44
1:A:399:LEU:HD13	1:A:399:LEU:O	2.18	0.44
1:A:444:ASP:HB3	1:A:460:SER:CB	2.48	0.44
1:A:54:LEU:HD13	1:A:63:LEU:CD2	2.48	0.43
1:A:521:GLU:O	1:A:522:LYS:HD3	2.16	0.43
1:A:395:ILE:HG23	1:A:399:LEU:HB2	2.00	0.43
1:A:204:ILE:O	1:A:207:VAL:HG22	2.18	0.43
1:A:461:VAL:HG21	1:A:476:LYS:HE2	2.00	0.43
1:A:106:TRP:CZ3	1:A:117:VAL:HG11	2.54	0.42
1:A:323:ILE:O	1:A:326:LEU:HB2	2.19	0.42
1:A:359:ILE:HD13	1:A:417:VAL:HG21	2.02	0.42
1:A:373:VAL:HG12	1:A:399:LEU:HD23	2.01	0.42
1:A:56:ILE:HD12	1:A:58:MET:H	1.84	0.42
1:A:527:ARG:O	1:A:528:MET:CB	2.68	0.42
1:A:458:SER:O	1:A:462:GLU:OE1	2.39	0.41
1:A:441:ARG:HD2	2:A:737:HOH:O	2.19	0.41
1:A:437:ILE:O	1:A:441:ARG:HG2	2.20	0.41
1:A:54:LEU:HD11	1:A:89:HIS:CD2	2.56	0.41
1:A:213:TYR:CE2	1:A:219:HIS:HB2	2.56	0.41
1:A:229:LEU:HD13	2:A:806:HOH:O	2.19	0.41
1:A:362:CYS:HB3	1:A:422:PHE:CD1	2.56	0.41
1:A:522:LYS:HD2	1:A:522:LYS:HA	1.58	0.41
1:A:336:ILE:HD12	1:A:336:ILE:N	2.36	0.41
1:A:407:THR:O	1:A:410:TYR:HB2	2.20	0.41
1:A:54:LEU:HA	1:A:63:LEU:HD22	2.01	0.41

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	517/563 (92%)	507 (98%)	9 (2%)	1 (0%)	43	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	528	MET

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	467/507 (92%)	450 (96%)	17 (4%)	31	31

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LEU
1	A	56	ILE
1	A	201	LEU
1	A	207	VAL
1	A	220	ASN
1	A	229	LEU
1	A	303	ASP
1	A	326	LEU
1	A	395	ILE
1	A	473	GLU

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Mol	Chain	Res	Type
1	A	496	ILE
1	A	500	ILE
1	A	518	LEU
1	A	522	LYS
1	A	526	THR
1	A	528	MET
1	A	533	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	33	GLN
1	A	37	GLN
1	A	85	HIS
1	A	125	HIS
1	A	134	GLN
1	A	215	GLN
1	A	219	HIS
1	A	220	ASN
1	A	232	ASN
1	A	235	GLN
1	A	379	GLN
1	A	385	HIS
1	A	513	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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	521/563 (92%)	0.15	31 (5%) 27 26	12, 24, 52, 71	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	528	MET	6.1
1	A	304	ALA	4.1
1	A	32	GLU	3.9
1	A	546	ILE	3.9
1	A	446	ALA	3.9
1	A	448	HIS	3.8
1	A	462	GLU	3.3
1	A	29	ASN	3.3
1	A	527	ARG	3.1
1	A	463	SER	3.0
1	A	447	GLY	2.9
1	A	529	GLY	2.8
1	A	460	SER	2.7
1	A	31	VAL	2.7
1	A	217	ASP	2.7
1	A	388	THR	2.7
1	A	526	THR	2.7
1	A	27	VAL	2.7
1	A	127	ASP	2.7
1	A	399	LEU	2.6
1	A	468	TYR	2.5
1	A	17	PRO	2.4
1	A	326	LEU	2.4
1	A	459	SER	2.4
1	A	476	LYS	2.2
1	A	458	SER	2.2
1	A	532	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	28	ASP	2.1
1	A	395	ILE	2.1
1	A	30	GLN	2.0
1	A	522	LYS	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.