



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:59 AM UTC

PDB ID : 5ERF / pdb_00005erf
Title : Ketosynthase from module 6 of the bacillaene synthase from *Bacillus subtilis* 168
Authors : Wagner, D.T.; Gay, D.C.; Keatinge-Clay, A.T.
Deposited on : 2015-11-14
Resolution : 3.10 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

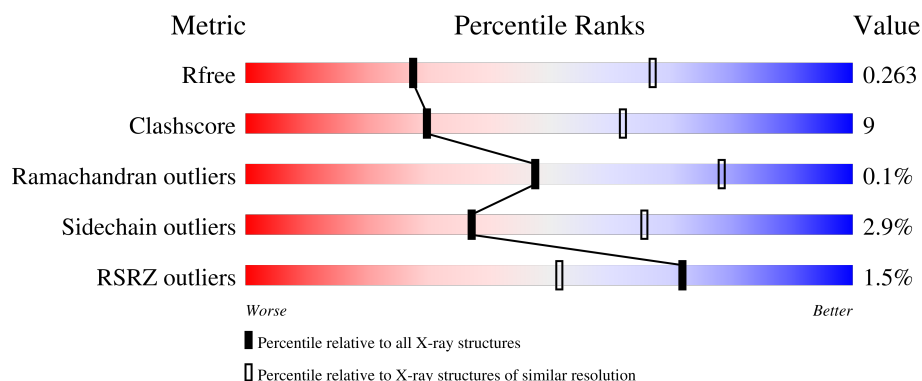
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

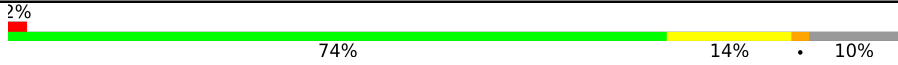
The reported resolution of this entry is 3.10 Å.

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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	617	
1	B	617	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8719 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 Polyketide synthase PksL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	553	Total	C	N	O	S	0	0	0
			4329	2748	726	829	26			
1	B	560	Total	C	N	O	S	0	0	0
			4390	2789	734	840	27			

There are 40 discrepancies between the modelled and reference sequences:

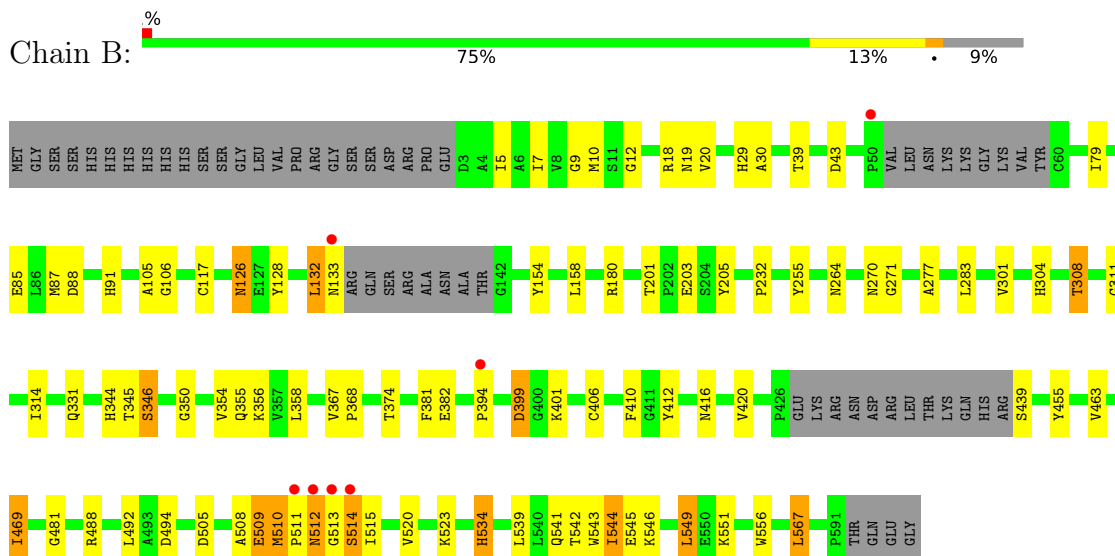
Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	initiating methionine	UNP Q05470
A	-20	GLY	-	expression tag	UNP Q05470
A	-19	SER	-	expression tag	UNP Q05470
A	-18	SER	-	expression tag	UNP Q05470
A	-17	HIS	-	expression tag	UNP Q05470
A	-16	HIS	-	expression tag	UNP Q05470
A	-15	HIS	-	expression tag	UNP Q05470
A	-14	HIS	-	expression tag	UNP Q05470
A	-13	HIS	-	expression tag	UNP Q05470
A	-12	HIS	-	expression tag	UNP Q05470
A	-11	SER	-	expression tag	UNP Q05470
A	-10	SER	-	expression tag	UNP Q05470
A	-9	GLY	-	expression tag	UNP Q05470
A	-8	LEU	-	expression tag	UNP Q05470
A	-7	VAL	-	expression tag	UNP Q05470
A	-6	PRO	-	expression tag	UNP Q05470
A	-5	ARG	-	expression tag	UNP Q05470
A	-4	GLY	-	expression tag	UNP Q05470
A	-3	SER	-	expression tag	UNP Q05470
A	-2	SER	-	expression tag	UNP Q05470
B	-21	MET	-	initiating methionine	UNP Q05470
B	-20	GLY	-	expression tag	UNP Q05470
B	-19	SER	-	expression tag	UNP Q05470
B	-18	SER	-	expression tag	UNP Q05470
B	-17	HIS	-	expression tag	UNP Q05470

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	HIS	-	expression tag	UNP Q05470
B	-15	HIS	-	expression tag	UNP Q05470
B	-14	HIS	-	expression tag	UNP Q05470
B	-13	HIS	-	expression tag	UNP Q05470
B	-12	HIS	-	expression tag	UNP Q05470
B	-11	SER	-	expression tag	UNP Q05470
B	-10	SER	-	expression tag	UNP Q05470
B	-9	GLY	-	expression tag	UNP Q05470
B	-8	LEU	-	expression tag	UNP Q05470
B	-7	VAL	-	expression tag	UNP Q05470
B	-6	PRO	-	expression tag	UNP Q05470
B	-5	ARG	-	expression tag	UNP Q05470
B	-4	GLY	-	expression tag	UNP Q05470
B	-3	SER	-	expression tag	UNP Q05470
B	-2	SER	-	expression tag	UNP Q05470

- Molecule 1: Polyketide synthase PksL



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	69.06Å 77.32Å 85.29Å 77.91° 70.26° 64.18°	Depositor
Resolution (Å)	39.73 – 3.10 39.73 – 3.10	Depositor EDS
% Data completeness (in resolution range)	94.1 (39.73-3.10) 94.1 (39.73-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.219 , 0.268 0.218 , 0.263	Depositor DCC
R_{free} test set	1285 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	71.8	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8719	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% 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	0.58	0/4424	0.94	12/5982 (0.2%)
1	B	0.58	0/4489	1.01	18/6072 (0.3%)
All	All	0.58	0/8913	0.98	30/12054 (0.2%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	512	ASN	CB-CA-C	-20.30	71.54	110.11
1	A	508	ALA	N-CA-C	-12.70	95.03	112.94
1	B	512	ASN	N-CA-C	11.95	127.66	113.20
1	B	515	ILE	N-CA-C	11.55	133.35	109.34
1	B	510	MET	N-CA-C	10.84	133.77	109.81
1	B	514	SER	CB-CA-C	9.52	129.37	110.42
1	B	510	MET	CB-CA-C	-8.73	92.97	110.17
1	B	515	ILE	N-CA-CB	-7.68	98.56	111.23
1	A	507	LEU	N-CA-C	7.49	126.74	110.80
1	B	549	LEU	N-CA-C	-7.11	104.74	113.41
1	A	502	ALA	N-CA-C	-7.07	101.86	112.04
1	B	271	GLY	N-CA-C	-6.93	100.19	110.87
1	A	501	LYS	CB-CA-C	-6.86	96.97	110.27
1	A	507	LEU	CB-CA-C	-6.82	96.85	110.42
1	A	515	ILE	N-CA-CB	6.32	122.75	111.63
1	A	456	ALA	N-CA-C	-6.29	103.56	111.11
1	B	514	SER	N-CA-C	-6.28	97.43	110.80
1	B	18	ARG	N-CA-C	-6.19	105.22	112.89
1	A	455	TYR	CA-C-N	5.81	128.38	120.54
1	A	455	TYR	C-N-CA	5.81	128.38	120.54
1	A	271	GLY	N-CA-C	-5.71	100.86	111.14
1	B	544	ILE	N-CA-C	-5.60	104.41	110.23
1	B	126	ASN	CB-CA-C	-5.54	102.11	111.02
1	B	304	HIS	N-CA-C	-5.54	104.94	110.97
1	A	544	ILE	N-CA-C	-5.51	104.49	110.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	549	LEU	N-CA-CB	5.48	118.48	110.47
1	B	132	LEU	N-CA-C	-5.45	101.29	109.79
1	B	133	ASN	N-CA-CB	5.28	119.48	110.50
1	A	18	ARG	N-CA-C	-5.26	105.60	112.23
1	B	399	ASP	CB-CA-C	-5.11	110.66	116.54

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	4329	0	4235	85	3
1	B	4390	0	4289	63	3
All	All	8719	0	8524	147	3

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 9.

All (147) 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:505:ASP:HB3	1:B:511:PRO:CG	1.34	1.54
1:B:505:ASP:CB	1:B:511:PRO:CG	2.18	1.20
1:A:505:ASP:CB	1:A:511:PRO:HB3	1.69	1.20
1:A:505:ASP:HB3	1:A:511:PRO:HB3	1.17	1.11
1:B:505:ASP:HB3	1:B:511:PRO:HG2	1.22	1.10
1:A:505:ASP:HB3	1:A:511:PRO:CB	1.81	1.09
1:B:505:ASP:HB3	1:B:511:PRO:CD	1.82	1.09
1:B:505:ASP:CG	1:B:511:PRO:HG3	1.79	1.07
1:B:505:ASP:CB	1:B:511:PRO:HG3	1.84	1.05
1:A:43:ASP:OD1	1:A:45:ASP:N	2.00	0.93
1:B:505:ASP:HB3	1:B:511:PRO:HG3	1.39	0.90
1:B:505:ASP:CB	1:B:511:PRO:HG2	1.96	0.85
1:A:131:MET:HE2	1:A:201:THR:HG22	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:ILE:HD13	1:A:87:MET:HE1	1.65	0.77
1:A:505:ASP:CG	1:A:511:PRO:HA	2.09	0.77
1:B:79:ILE:HD13	1:B:87:MET:HE1	1.68	0.74
1:A:331:GLN:OE1	1:A:387:TYR:HB3	1.88	0.72
1:A:505:ASP:CG	1:A:511:PRO:HB3	2.15	0.70
1:A:492:LEU:HD21	1:A:549:LEU:HB2	1.75	0.69
1:B:399:ASP:HB2	1:B:401:LYS:HG2	1.76	0.68
1:B:492:LEU:HD21	1:B:549:LEU:HB2	1.75	0.67
1:A:399:ASP:HB2	1:A:401:LYS:HG2	1.76	0.66
1:B:505:ASP:OD2	1:B:511:PRO:HG3	1.95	0.66
1:A:270:ASN:O	1:A:270:ASN:ND2	2.28	0.66
1:A:510:MET:N	1:A:511:PRO:CD	2.59	0.65
1:A:131:MET:CE	1:A:201:THR:CG2	2.75	0.64
1:A:301:VAL:HG22	1:A:406:CYS:HB2	1.79	0.64
1:B:301:VAL:HG22	1:B:406:CYS:HB2	1.80	0.64
1:B:88:ASP:HB3	1:B:91:HIS:CD2	2.33	0.64
1:A:505:ASP:CG	1:A:511:PRO:CA	2.71	0.63
1:B:270:ASN:O	1:B:270:ASN:ND2	2.30	0.63
1:A:505:ASP:HB3	1:A:511:PRO:CG	2.29	0.62
1:A:7:ILE:HD13	1:A:358:LEU:HD11	1.82	0.62
1:A:131:MET:HE2	1:A:201:THR:CG2	2.30	0.61
1:A:501:LYS:O	1:A:501:LYS:CG	2.46	0.61
1:B:7:ILE:HD13	1:B:358:LEU:HD11	1.82	0.61
1:A:504:ASP:O	1:A:507:LEU:HB2	2.01	0.61
1:A:131:MET:HE3	1:A:201:THR:HG23	1.82	0.61
1:A:88:ASP:HB3	1:A:91:HIS:CD2	2.37	0.60
1:A:455:TYR:O	1:A:458:ALA:HB3	2.02	0.60
1:A:510:MET:N	1:A:511:PRO:HD3	2.16	0.59
1:B:544:ILE:HG23	1:B:567:LEU:HD12	1.84	0.59
1:A:544:ILE:HG23	1:A:567:LEU:HD12	1.84	0.58
1:B:505:ASP:CB	1:B:511:PRO:CD	2.69	0.57
1:A:542:THR:O	1:A:546:LYS:HB2	2.03	0.57
1:A:492:LEU:HD12	1:A:492:LEU:N	2.19	0.57
1:A:505:ASP:CB	1:A:511:PRO:CB	2.54	0.57
1:B:492:LEU:HD12	1:B:492:LEU:N	2.19	0.57
1:A:508:ALA:O	1:A:509:GLU:HG2	2.05	0.56
1:A:505:ASP:O	1:A:509:GLU:HA	2.06	0.56
1:B:546:LYS:HG2	1:B:546:LYS:O	2.05	0.56
1:A:505:ASP:CG	1:A:511:PRO:CB	2.78	0.54
1:A:131:MET:HE3	1:A:201:THR:CG2	2.38	0.54
1:B:205:TYR:CZ	1:B:232:PRO:HG2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:508:ALA:C	1:A:509:GLU:HG2	2.32	0.53
1:B:264:ASN:ND2	1:B:283:LEU:HB2	2.23	0.53
1:B:201:THR:HB	1:B:203:GLU:OE2	2.08	0.53
1:A:201:THR:HB	1:A:203:GLU:OE2	2.08	0.52
1:A:424:TYR:CG	1:A:425:GLN:N	2.77	0.52
1:B:512:ASN:N	1:B:513:GLY:HA2	2.25	0.52
1:A:516:PHE:CD1	1:A:516:PHE:N	2.76	0.52
1:B:512:ASN:N	1:B:513:GLY:CA	2.73	0.51
1:A:205:TYR:CZ	1:A:232:PRO:HG2	2.45	0.51
1:A:406:CYS:SG	1:A:420:VAL:HG22	2.51	0.51
1:B:308:THR:CG2	1:B:311:GLY:HA3	2.41	0.51
1:B:12:GLY:O	1:B:19:ASN:HA	2.10	0.51
1:A:264:ASN:ND2	1:A:283:LEU:HB2	2.25	0.51
1:B:308:THR:HG23	1:B:311:GLY:H	1.75	0.50
1:B:505:ASP:HB3	1:B:511:PRO:HD3	1.82	0.50
1:A:12:GLY:O	1:A:19:ASN:HA	2.10	0.50
1:A:308:THR:CG2	1:A:311:GLY:HA3	2.42	0.50
1:A:508:ALA:O	1:A:510:MET:HG3	2.12	0.49
1:B:406:CYS:SG	1:B:420:VAL:HG22	2.52	0.49
1:A:264:ASN:OD1	1:A:264:ASN:C	2.55	0.49
1:B:39:THR:HG22	1:B:43:ASP:HA	1.94	0.49
1:B:277:ALA:HB2	1:B:314:ILE:HG23	1.93	0.49
1:A:308:THR:HG23	1:A:311:GLY:H	1.77	0.49
1:A:412:TYR:C	1:A:412:TYR:CD1	2.91	0.49
1:B:412:TYR:CD1	1:B:412:TYR:C	2.91	0.49
1:A:39:THR:HG22	1:A:43:ASP:HA	1.94	0.49
1:A:9:GLY:HA3	1:A:105:ALA:HB2	1.96	0.48
1:A:43:ASP:OD1	1:A:45:ASP:HB2	2.14	0.48
1:A:455:TYR:O	1:A:458:ALA:N	2.46	0.48
1:A:505:ASP:O	1:A:509:GLU:CA	2.62	0.48
1:A:508:ALA:O	1:A:509:GLU:CG	2.62	0.48
1:A:277:ALA:HB2	1:A:314:ILE:HG23	1.95	0.48
1:A:344:HIS:CD2	1:A:346:SER:H	2.32	0.47
1:B:9:GLY:HA3	1:B:105:ALA:HB2	1.95	0.47
1:B:344:HIS:CD2	1:B:346:SER:H	2.33	0.47
1:A:355:GLN:O	1:A:356:LYS:C	2.57	0.46
1:B:539:LEU:O	1:B:542:THR:HB	2.14	0.46
1:A:117:CYS:HB3	1:A:158:LEU:HD13	1.96	0.46
1:A:350:GLY:O	1:A:354:VAL:HG23	2.15	0.46
1:A:505:ASP:HB3	1:A:511:PRO:CD	2.44	0.46
1:B:510:MET:HB2	1:B:511:PRO:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:PHE:HD1	1:A:416:ASN:HB3	1.81	0.46
1:A:539:LEU:O	1:A:542:THR:HB	2.16	0.46
1:B:511:PRO:C	1:B:513:GLY:HA3	2.41	0.46
1:A:344:HIS:HD2	1:A:346:SER:H	1.64	0.46
1:A:29:HIS:O	1:A:30:ALA:C	2.59	0.45
1:A:423:GLU:HG2	1:A:424:TYR:O	2.15	0.45
1:B:132:LEU:HA	1:B:132:LEU:HD23	1.82	0.45
1:A:21:ARG:HE	1:A:462:PHE:HE1	1.63	0.45
1:A:131:MET:CE	1:A:201:THR:HG23	2.42	0.45
1:B:505:ASP:CB	1:B:511:PRO:HD3	2.43	0.45
1:B:350:GLY:O	1:B:354:VAL:HG23	2.16	0.45
1:B:117:CYS:HB3	1:B:158:LEU:HD13	1.98	0.45
1:B:270:ASN:C	1:B:270:ASN:HD22	2.22	0.45
1:A:169:CYS:HB2	1:A:411:GLY:HA2	1.98	0.44
1:B:106:GLY:O	1:B:481:GLY:HA2	2.18	0.44
1:B:355:GLN:O	1:B:356:LYS:C	2.60	0.44
1:A:424:TYR:C	1:A:425:GLN:HG3	2.42	0.44
1:A:270:ASN:ND2	1:A:270:ASN:C	2.76	0.44
1:A:21:ARG:O	1:A:22:GLU:C	2.60	0.44
1:B:488:ARG:HB3	1:B:556:TRP:CZ3	2.53	0.44
1:B:344:HIS:HD2	1:B:346:SER:H	1.66	0.43
1:A:270:ASN:C	1:A:270:ASN:HD22	2.19	0.43
1:B:508:ALA:O	1:B:509:GLU:HB2	2.18	0.43
1:A:106:GLY:O	1:A:481:GLY:HA2	2.19	0.43
1:A:488:ARG:HB3	1:A:556:TRP:CZ3	2.53	0.43
1:B:410:PHE:HD1	1:B:416:ASN:HB3	1.84	0.43
1:B:543:TRP:CE2	1:B:551:LYS:CD	3.02	0.43
1:A:501:LYS:O	1:A:501:LYS:HG2	2.19	0.43
1:B:29:HIS:O	1:B:30:ALA:C	2.61	0.43
1:A:534:HIS:ND1	1:A:534:HIS:C	2.77	0.43
1:B:10:MET:HE3	1:B:255:TYR:CD2	2.54	0.42
1:A:505:ASP:O	1:A:509:GLU:N	2.52	0.42
1:A:91:HIS:CE1	1:A:148:ALA:HB2	2.55	0.42
1:B:270:ASN:ND2	1:B:270:ASN:C	2.75	0.42
1:A:505:ASP:OD2	1:A:511:PRO:HA	2.20	0.42
1:A:10:MET:HE3	1:A:255:TYR:CD2	2.54	0.42
1:A:505:ASP:HB3	1:A:511:PRO:CA	2.46	0.42
1:B:463:VAL:HA	1:B:469:ILE:HD11	2.02	0.41
1:B:85:GLU:O	1:B:128:TYR:OH	2.26	0.41
1:A:410:PHE:CD1	1:A:416:ASN:HB3	2.56	0.41
1:A:24:TRP:HB2	1:A:359:LEU:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:PHE:O	1:A:382:GLU:C	2.63	0.41
1:B:541:GLN:O	1:B:545:GLU:HB2	2.21	0.41
1:B:439:SER:HA	1:B:494:ASP:HB3	2.03	0.41
1:B:534:HIS:ND1	1:B:534:HIS:C	2.79	0.40
1:B:5:ILE:HG13	1:B:180:ARG:HG3	2.04	0.40
1:A:424:TYR:O	1:A:425:GLN:HG3	2.22	0.40
1:A:580:ALA:O	1:A:581:TYR:C	2.64	0.40
1:B:381:PHE:O	1:B:382:GLU:C	2.64	0.40
1:B:264:ASN:OD1	1:B:264:ASN:C	2.64	0.40
1:B:367:VAL:HB	1:B:368:PRO:CD	2.52	0.40
1:A:272:ILE:HA	1:B:154:TYR:CE1	2.56	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:512:ASN:OD1	1:B:394:PRO:CB[1_456]	2.07	0.13
1:A:512:ASN:CG	1:B:394:PRO:CB[1_456]	2.17	0.03
1:A:512:ASN:C	1:B:394:PRO:CG[1_456]	2.18	0.02

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	545/617 (88%)	505 (93%)	39 (7%)	1 (0%)	43	73
1	B	552/617 (90%)	517 (94%)	35 (6%)	0	100	100
All	All	1097/1234 (89%)	1022 (93%)	74 (7%)	1 (0%)	48	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	514	SER

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	458/515 (89%)	446 (97%)	12 (3%)	40	68
1	B	465/515 (90%)	450 (97%)	15 (3%)	34	64
All	All	923/1030 (90%)	896 (97%)	27 (3%)	37	66

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	VAL
1	A	308	THR
1	A	345	THR
1	A	346	SER
1	A	455	TYR
1	A	469	ILE
1	A	510	MET
1	A	514	SER
1	A	520	VAL
1	A	523	LYS
1	A	534	HIS
1	A	567	LEU
1	B	20	VAL
1	B	126	ASN
1	B	308	THR
1	B	331	GLN
1	B	345	THR
1	B	346	SER
1	B	374	THR
1	B	455	TYR
1	B	469	ILE
1	B	509	GLU
1	B	514	SER

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Mol	Chain	Res	Type
1	B	520	VAL
1	B	523	LYS
1	B	534	HIS
1	B	567	LEU

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	91	HIS
1	A	181	GLN
1	A	344	HIS
1	A	378	HIS
1	A	416	ASN
1	A	541	GLN
1	B	91	HIS
1	B	126	ASN
1	B	181	GLN
1	B	344	HIS
1	B	378	HIS
1	B	416	ASN
1	B	519	HIS
1	B	541	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 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	553/617 (89%)	-0.24	10 (1%) 67 47	25, 73, 126, 164	0
1	B	560/617 (90%)	-0.32	7 (1%) 75 56	25, 67, 107, 144	0
All	All	1113/1234 (90%)	-0.28	17 (1%) 72 52	25, 70, 119, 164	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	512	ASN	9.5
1	B	512	ASN	7.1
1	A	511	PRO	6.2
1	A	513	GLY	5.7
1	B	513	GLY	5.6
1	B	514	SER	3.4
1	B	511	PRO	3.4
1	B	133	ASN	3.1
1	A	514	SER	2.6
1	A	133	ASN	2.5
1	B	394	PRO	2.2
1	A	425	GLN	2.1
1	B	50	PRO	2.1
1	A	510	MET	2.1
1	A	412	TYR	2.1
1	A	169	CYS	2.0
1	A	142	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 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.