



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 12:21 PM UTC

PDB ID : 2DZA / pdb_00002dza
Title : Crystal Structure of Dihydropteroate Synthase from *Thermus thermophilus* HB8 in Complex with 4-aminobenzoate
Authors : Bagautdinov, B.; Kunishima, N.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2006-09-27
Resolution : 1.90 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

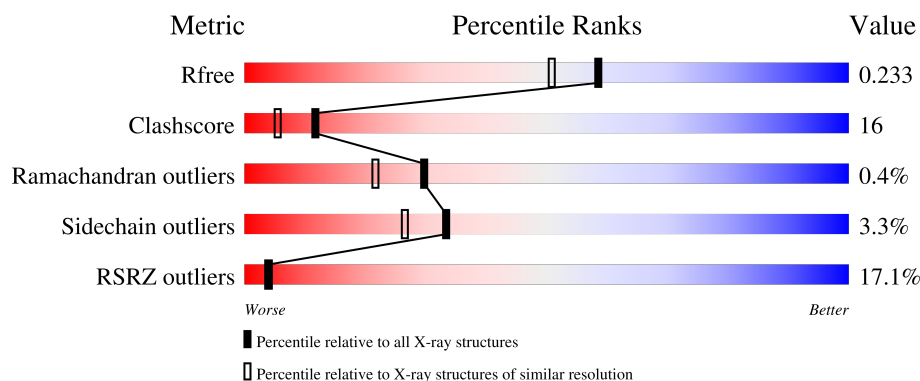
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	294	12% 63% 17% • 17%
1	B	294	17% 62% 23% • 13%

2 Entry composition [i](#)

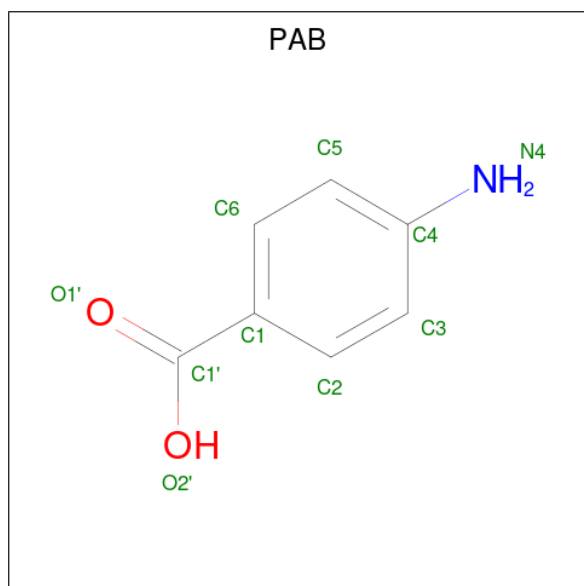
There are 3 unique types of molecules in this entry. The entry contains 4269 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dihydropteroate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1874	1185	354	328	7			
1	B	257	Total	C	N	O	S	0	0	0
			1977	1247	373	350	7			

- Molecule 2 is 4-AMINO BENZOIC ACID (CCD ID: PAB) (formula: $C_7H_7NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			10	7	1	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	247	Total	O	0	0
			247	247		

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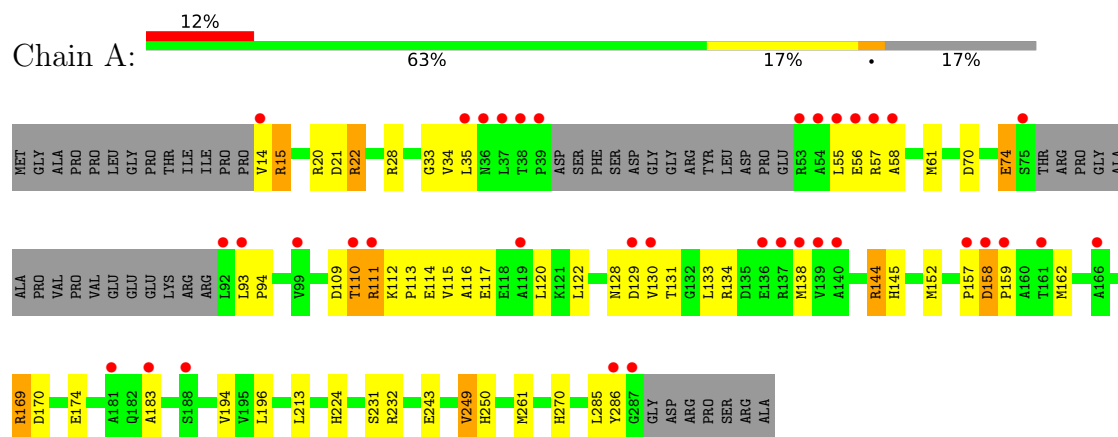
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	161	Total 161	O 161	0	0

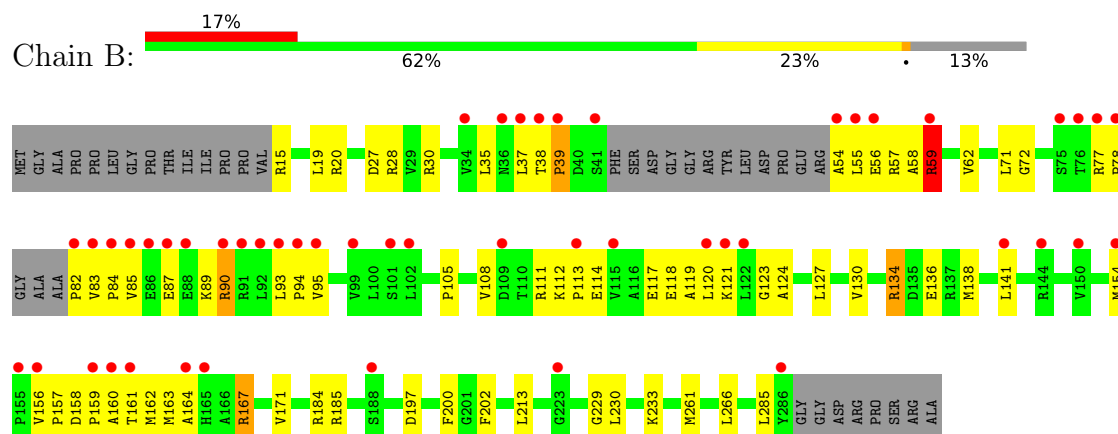
3 Residue-property plots [i](#)

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: Dihydropteroate synthase



• Molecule 1: Dihydropteroate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	110.60Å 110.60Å 88.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.20 – 1.90 36.20 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (36.20-1.90) 99.8 (36.20-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 1.89Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.215 , 0.233 0.215 , 0.233	Depositor DCC
R_{free} test set	2410 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	29.7	Xtriage
Anisotropy	0.321	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.038 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4269	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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

Bond lengths and bond angles in the following residue types are not validated in this section: PAB

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	0.52	0/1902	0.97	7/2577 (0.3%)
1	B	0.53	0/2008	1.02	9/2721 (0.3%)
All	All	0.53	0/3910	1.00	16/5298 (0.3%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	ASP	CA-C-N	6.48	126.11	119.56
1	A	158	ASP	C-N-CA	6.48	126.11	119.56
1	B	171	VAL	N-CA-C	5.79	117.22	110.62
1	B	84	PRO	N-CA-CB	5.63	108.31	103.36
1	B	134	ARG	CD-NE-CZ	-5.56	116.62	124.40
1	B	197	ASP	CA-C-N	5.54	125.21	119.56
1	B	197	ASP	C-N-CA	5.54	125.21	119.56
1	A	34	VAL	N-CA-C	5.45	116.08	108.89
1	B	90	ARG	CD-NE-CZ	-5.13	117.22	124.40
1	A	28	ARG	CD-NE-CZ	-5.12	117.23	124.40
1	B	28	ARG	CD-NE-CZ	-5.10	117.25	124.40
1	A	249	VAL	N-CA-C	5.10	115.82	110.62
1	A	144	ARG	CD-NE-CZ	-5.09	117.27	124.40
1	B	59	ARG	CD-NE-CZ	-5.07	117.30	124.40
1	A	111	ARG	CD-NE-CZ	-5.05	117.33	124.40
1	B	57	ARG	CD-NE-CZ	-5.01	117.39	124.40

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	1874	0	1953	57	0
1	B	1977	0	2054	74	0
2	B	10	0	6	0	0
3	A	247	0	0	8	0
3	B	161	0	0	10	0
All	All	4269	0	4013	125	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 16.

All (125) 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:116:ALA:HB2	1:A:138:MET:HE2	1.43	0.97
1:B:59:ARG:HH11	1:B:59:ARG:HB3	1.30	0.93
1:A:116:ALA:CB	1:A:138:MET:HE2	1.99	0.91
1:A:14:VAL:HG22	1:A:15:ARG:H	1.35	0.91
1:B:77:ARG:HG3	1:B:78:PRO:HD2	1.59	0.84
1:A:134:ARG:HD2	1:A:157:PRO:HG3	1.66	0.77
1:B:185:ARG:HD3	3:B:1394:HOH:O	1.85	0.76
1:A:110:THR:HG21	1:A:115:VAL:HG11	1.67	0.76
1:A:113:PRO:HA	1:A:138:MET:HE3	1.69	0.74
1:B:134:ARG:HD2	1:B:157:PRO:HG3	1.70	0.73
1:B:167:ARG:HA	1:B:167:ARG:HE	1.54	0.72
1:B:130:VAL:HB	1:B:154:MET:HE1	1.71	0.72
1:B:20:ARG:NH1	1:B:123:GLY:HA2	2.05	0.71
1:B:59:ARG:HH11	1:B:59:ARG:CB	2.03	0.71
1:A:113:PRO:CA	1:A:138:MET:HE3	2.22	0.70
1:B:185:ARG:HD2	3:B:1363:HOH:O	1.91	0.69
1:A:14:VAL:HG22	1:A:15:ARG:N	2.05	0.68
1:A:116:ALA:HB2	1:A:138:MET:CE	2.22	0.67
1:B:54:ALA:CB	1:B:95:VAL:HG13	2.25	0.67
1:B:20:ARG:HH12	1:B:123:GLY:HA2	1.58	0.67
1:A:56:GLU:HG3	1:A:57:ARG:N	2.13	0.64
1:B:82:PRO:HB2	1:B:160:ALA:HB2	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:ARG:HG3	1:B:78:PRO:CD	2.29	0.63
1:B:56:GLU:HA	1:B:56:GLU:OE1	1.97	0.62
1:A:261:MET:SD	1:B:261:MET:SD	2.97	0.62
1:A:113:PRO:HA	1:A:138:MET:CE	2.30	0.62
1:B:77:ARG:NH1	3:B:1365:HOH:O	2.29	0.62
1:A:128:ASN:C	1:A:128:ASN:HD22	2.08	0.61
1:A:15:ARG:NH2	3:A:538:HOH:O	2.34	0.61
1:A:113:PRO:N	1:A:138:MET:HE3	2.14	0.61
1:A:285:LEU:HD21	1:B:213:LEU:HD13	1.82	0.61
1:A:74:GLU:OE1	1:A:110:THR:HG22	2.01	0.60
1:B:117:GLU:HB2	1:B:141:LEU:HD11	1.83	0.60
1:B:229:GLY:O	1:B:233:LYS:HE2	2.02	0.60
1:B:54:ALA:HB1	1:B:95:VAL:CG1	2.32	0.60
1:A:55:LEU:H	1:A:55:LEU:HD23	1.68	0.59
1:A:232:ARG:HG2	3:A:356:HOH:O	2.01	0.58
1:B:167:ARG:HE	1:B:167:ARG:CA	2.16	0.58
1:A:162:MET:HE3	3:A:327:HOH:O	2.03	0.58
1:B:54:ALA:HB1	1:B:95:VAL:HG13	1.85	0.58
1:B:108:VAL:HG21	1:B:119:ALA:HB1	1.86	0.57
1:B:117:GLU:OE1	1:B:141:LEU:HD21	2.04	0.57
1:A:110:THR:HG21	1:A:115:VAL:CG1	2.35	0.57
1:A:93:LEU:HB2	1:A:94:PRO:HD3	1.87	0.57
1:B:55:LEU:O	1:B:59:ARG:HG3	2.05	0.56
1:B:167:ARG:HA	1:B:167:ARG:NE	2.21	0.56
1:B:130:VAL:HB	1:B:154:MET:CE	2.36	0.55
1:A:15:ARG:HG2	3:A:305:HOH:O	2.06	0.55
1:B:111:ARG:C	1:B:138:MET:HE2	2.31	0.55
1:A:56:GLU:HG3	1:A:57:ARG:H	1.71	0.54
1:B:77:ARG:CG	1:B:78:PRO:HD2	2.36	0.54
1:A:15:ARG:HB3	3:A:421:HOH:O	2.08	0.53
1:B:58:ALA:O	1:B:62:VAL:HG23	2.09	0.53
1:B:59:ARG:HB3	1:B:59:ARG:NH1	2.12	0.53
1:B:38:THR:HB	1:B:39:PRO:HD2	1.90	0.53
1:B:89:LYS:HE2	1:B:118:GLU:HG3	1.90	0.52
1:B:154:MET:HE2	1:B:202:PHE:CD2	2.46	0.51
1:A:14:VAL:CG2	1:A:15:ARG:H	2.14	0.51
1:A:112:LYS:C	1:A:138:MET:HE3	2.36	0.51
1:B:233:LYS:HE3	3:B:1457:HOH:O	2.10	0.51
1:B:35:LEU:HD23	1:B:71:LEU:HG	1.93	0.51
1:A:33:GLY:O	1:A:61:MET:HE1	2.11	0.50
1:B:83:VAL:HG23	1:B:83:VAL:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:GLY:C	1:A:61:MET:HE1	2.36	0.50
1:A:113:PRO:HG3	1:A:138:MET:HG3	1.93	0.50
1:A:70:ASP:OD2	1:A:109:ASP:OD2	2.30	0.50
1:A:109:ASP:OD1	1:A:128:ASN:HB3	2.12	0.50
1:A:129:ASP:OD2	1:A:133:LEU:HA	2.13	0.49
1:B:54:ALA:CB	1:B:95:VAL:CG1	2.90	0.49
1:A:21:ASP:OD1	1:A:22:ARG:HD2	2.13	0.48
1:B:134:ARG:CD	1:B:157:PRO:HG3	2.41	0.48
1:A:261:MET:HG3	1:A:286:TYR:CE2	2.49	0.48
1:B:85:VAL:O	1:B:112:LYS:HE3	2.13	0.48
1:B:93:LEU:N	1:B:94:PRO:CD	2.76	0.48
1:A:170:ASP:O	1:A:174:GLU:HG3	2.13	0.48
1:A:231:SER:HB3	1:A:270:HIS:HD2	1.79	0.48
1:A:286:TYR:CD2	1:B:261:MET:HE1	2.49	0.48
1:A:116:ALA:HB3	1:A:138:MET:HE2	1.88	0.47
1:B:184:ARG:HG3	1:B:184:ARG:HH11	1.79	0.47
1:A:152:MET:HE1	3:A:511:HOH:O	2.15	0.46
1:B:90:ARG:O	1:B:94:PRO:HG2	2.16	0.46
1:B:113:PRO:HG3	1:B:138:MET:HG2	1.97	0.46
1:A:117:GLU:OE1	1:A:145:HIS:NE2	2.43	0.46
1:A:169:ARG:CG	1:A:169:ARG:HH11	2.27	0.46
1:B:117:GLU:O	1:B:121:LYS:HG3	2.15	0.45
1:B:59:ARG:HH11	1:B:59:ARG:CG	2.28	0.45
1:A:128:ASN:C	1:A:128:ASN:ND2	2.74	0.45
1:B:27:ASP:HB3	3:B:1384:HOH:O	2.17	0.45
1:A:213:LEU:HD23	1:B:285:LEU:CD2	2.47	0.45
1:B:78:PRO:O	1:B:82:PRO:HD3	2.17	0.44
1:B:230:LEU:HD12	1:B:230:LEU:N	2.32	0.44
1:B:158:ASP:OD2	1:B:160:ALA:HB3	2.17	0.44
1:A:144:ARG:HD3	3:A:510:HOH:O	2.16	0.44
1:A:183:ALA:HB1	1:A:194:VAL:HG21	1.99	0.44
1:A:249:VAL:HG13	1:A:250:HIS:N	2.33	0.44
1:B:200:PHE:CE1	1:B:213:LEU:HG	2.53	0.44
1:A:213:LEU:HD23	1:B:285:LEU:HD21	1.99	0.43
1:A:20:ARG:NH2	1:A:120:LEU:O	2.51	0.43
1:A:243:GLU:OE2	3:A:403:HOH:O	2.21	0.43
1:B:136:GLU:HG2	1:B:185:ARG:HH21	1.83	0.43
1:A:196:LEU:HD11	1:A:224:HIS:HD2	1.83	0.43
1:A:55:LEU:O	1:A:58:ALA:HB3	2.19	0.43
1:B:89:LYS:HD3	3:B:1434:HOH:O	2.19	0.43
1:B:30:ARG:HB2	1:B:266:LEU:HD23	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:PRO:HA	1:B:162:MET:HE2	2.00	0.42
1:B:15:ARG:NH2	3:B:1442:HOH:O	2.35	0.42
1:B:114:GLU:H	1:B:114:GLU:HG3	1.53	0.42
1:B:37:LEU:N	1:B:72:GLY:O	2.48	0.42
1:B:159:PRO:CA	1:B:162:MET:HE2	2.49	0.42
1:A:158:ASP:HA	1:A:159:PRO:HD2	1.92	0.42
1:B:158:ASP:HA	1:B:159:PRO:HD3	1.89	0.42
1:B:55:LEU:HD12	1:B:55:LEU:HA	1.90	0.42
1:B:78:PRO:HA	1:B:162:MET:HE3	2.02	0.41
1:A:129:ASP:OD1	1:A:131:THR:OG1	2.30	0.41
1:B:136:GLU:HG3	3:B:1413:HOH:O	2.19	0.41
1:A:55:LEU:H	1:A:55:LEU:CD2	2.31	0.41
1:B:156:VAL:CG1	3:B:1356:HOH:O	2.69	0.41
1:A:112:LYS:HA	1:A:113:PRO:HD3	1.88	0.40
1:B:90:ARG:HH11	1:B:90:ARG:HD3	1.59	0.40
1:B:19:LEU:HD13	1:B:105:PRO:CB	2.52	0.40
1:B:120:LEU:HD23	1:B:124:ALA:HB3	2.04	0.40
1:B:161:THR:O	1:B:164:ALA:HB3	2.21	0.40
1:B:162:MET:HG3	1:B:163:MET:N	2.35	0.40
1:A:285:LEU:HD21	1:B:213:LEU:CD1	2.51	0.40
1:B:77:ARG:NH2	3:B:1365:HOH:O	2.53	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	239/294 (81%)	229 (96%)	9 (4%)	1 (0%)	30	22
1	B	251/294 (85%)	243 (97%)	7 (3%)	1 (0%)	30	22
All	All	490/588 (83%)	472 (96%)	16 (3%)	2 (0%)	30	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	39	PRO
1	A	130	VAL

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	193/232 (83%)	184 (95%)	9 (5%)	23	15
1	B	206/232 (89%)	202 (98%)	4 (2%)	50	47
All	All	399/464 (86%)	386 (97%)	13 (3%)	33	26

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ARG
1	A	22	ARG
1	A	35	LEU
1	A	74	GLU
1	A	110	THR
1	A	111	ARG
1	A	114	GLU
1	A	122	LEU
1	A	169	ARG
1	B	59	ARG
1	B	87	GLU
1	B	127	LEU
1	B	167	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	270	HIS
1	B	145	HIS
1	B	247	GLN

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 ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PAB	B	1301	-	10,10,10	1.71	3 (30%)	13,13,13	1.79	2 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PAB	B	1301	-	-	0/4/4/4	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1301	PAB	C2-C1	2.48	1.43	1.39
2	B	1301	PAB	O2'-C1'	-2.42	1.23	1.30
2	B	1301	PAB	C6-C1	2.36	1.43	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1301	PAB	O2'-C1'-O1'	-4.56	113.56	123.35
2	B	1301	PAB	O2'-C1'-C1	2.83	122.10	114.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	245/294 (83%)	0.82	36 (14%) 6 6	16, 34, 63, 86	0
1	B	257/294 (87%)	0.99	50 (19%) 3 3	16, 37, 70, 86	0
All	All	502/588 (85%)	0.91	86 (17%) 4 4	16, 35, 67, 86	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	55	LEU	7.3
1	B	83	VAL	5.5
1	B	85	VAL	5.0
1	A	14	VAL	4.8
1	B	54	ALA	4.8
1	A	38	THR	4.6
1	B	41	SER	4.3
1	B	55	LEU	4.3
1	A	286	TYR	4.2
1	A	37	LEU	4.1
1	B	101	SER	4.0
1	A	110	THR	3.8
1	A	35	LEU	3.7
1	B	77	ARG	3.7
1	B	38	THR	3.6
1	A	39	PRO	3.6
1	A	138	MET	3.6
1	B	286	TYR	3.5
1	B	82	PRO	3.5
1	B	88	GLU	3.4
1	B	156	VAL	3.4
1	A	53	ARG	3.3
1	B	87	GLU	3.3
1	B	36	ASN	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	141	LEU	3.2
1	A	287	GLY	3.2
1	A	56	GLU	3.2
1	A	129	ASP	3.1
1	B	39	PRO	3.1
1	A	130	VAL	3.1
1	B	102	LEU	3.1
1	A	54	ALA	3.0
1	B	84	PRO	3.0
1	B	109	ASP	2.9
1	A	58	ALA	2.9
1	B	160	ALA	2.9
1	B	78	PRO	2.8
1	A	92	LEU	2.8
1	A	75	SER	2.8
1	A	188	SER	2.8
1	B	37	LEU	2.8
1	B	56	GLU	2.8
1	A	119	ALA	2.7
1	B	93	LEU	2.7
1	B	95	VAL	2.7
1	A	57	ARG	2.6
1	B	76	THR	2.6
1	B	94	PRO	2.5
1	A	93	LEU	2.5
1	B	121	LYS	2.5
1	A	158	ASP	2.5
1	B	155	PRO	2.5
1	A	137	ARG	2.4
1	B	223	GLY	2.4
1	B	150	VAL	2.4
1	A	159	PRO	2.4
1	B	86	GLU	2.4
1	B	165	HIS	2.4
1	A	140	ALA	2.3
1	A	157	PRO	2.3
1	B	164	ALA	2.3
1	A	111	ARG	2.3
1	B	144	ARG	2.3
1	B	59	ARG	2.2
1	A	161	THR	2.2
1	A	136	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	159	PRO	2.2
1	A	181	ALA	2.2
1	B	154	MET	2.2
1	B	90	ARG	2.2
1	A	139	VAL	2.2
1	B	75	SER	2.2
1	B	120	LEU	2.2
1	B	122	LEU	2.2
1	A	99	VAL	2.1
1	B	161	THR	2.1
1	B	188	SER	2.1
1	A	36	ASN	2.1
1	A	166	ALA	2.1
1	B	34	VAL	2.1
1	A	183	ALA	2.1
1	B	113	PRO	2.0
1	B	92	LEU	2.0
1	B	91	ARG	2.0
1	B	99	VAL	2.0
1	B	115	VAL	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PAB	B	1301	10/10	0.94	0.08	32,33,37,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.