Full wwPDB X-ray Structure Validation Report ⓘ

Oct 25, 2023 – 09:39 PM EDT

PDB ID	:	3H9A
Title	:	Crystal structure of BacB, an enzyme involved in Bacilysin synthesis, in triclinic form
Authors	:	Rajavel, M.; Gopal, B.
Deposited on	:	2009-04-30
Resolution	:	2.04 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:

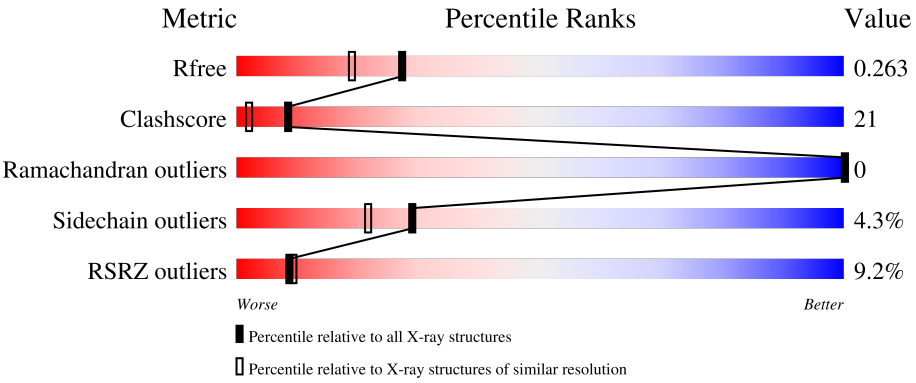
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 <=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	243	5% 71% 19% · 7%
1	B	243	11% 70% 21% · 7%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3814 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 Bacilysin biosynthesis protein bacB.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	Se	0	0	0
			1804	1142	303	344	3	12			
1	B	225	Total	C	N	O	S	Se	0	0	0
			1804	1142	303	344	3	12			

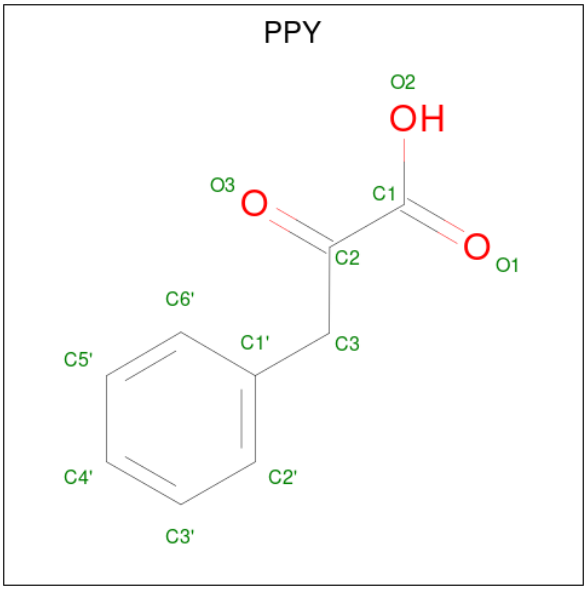
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	236	LEU	-	expression tag	UNP P39639
A	237	GLU	-	expression tag	UNP P39639
A	238	HIS	-	expression tag	UNP P39639
A	239	HIS	-	expression tag	UNP P39639
A	240	HIS	-	expression tag	UNP P39639
A	241	HIS	-	expression tag	UNP P39639
A	242	HIS	-	expression tag	UNP P39639
A	243	HIS	-	expression tag	UNP P39639
B	236	LEU	-	expression tag	UNP P39639
B	237	GLU	-	expression tag	UNP P39639
B	238	HIS	-	expression tag	UNP P39639
B	239	HIS	-	expression tag	UNP P39639
B	240	HIS	-	expression tag	UNP P39639
B	241	HIS	-	expression tag	UNP P39639
B	242	HIS	-	expression tag	UNP P39639
B	243	HIS	-	expression tag	UNP P39639

- Molecule 2 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Co	0	0
			2	2		
2	B	2	Total	Co	0	0
			2	2		

- Molecule 3 is 3-PHENYLPYRUVIC ACID (three-letter code: PPY) (formula: $C_9H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 12 9 3	0	0
3	B	1	Total C O 12 9 3	0	0

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Fe 1 1	0	0

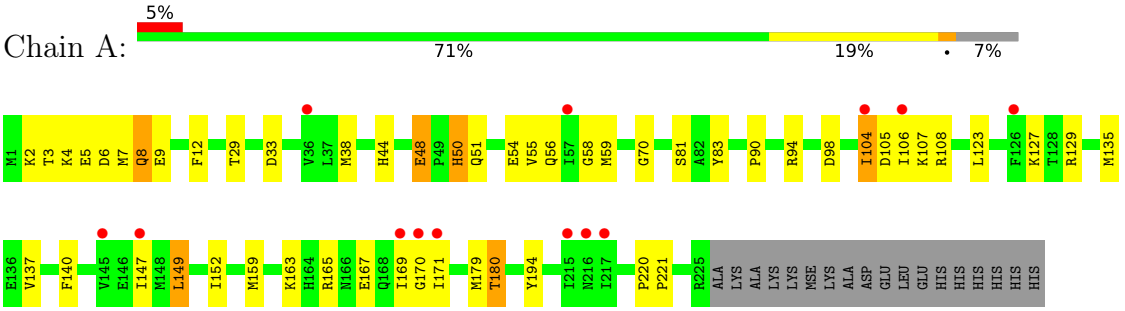
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	109	Total O 109 109	0	0
5	B	68	Total O 68 68	0	0

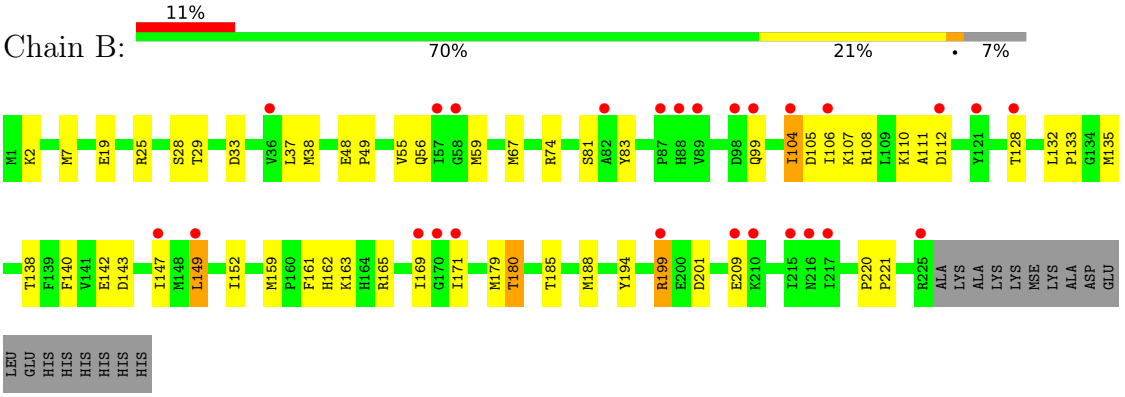
3 Residue-property plots (i)

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: Bacilysin biosynthesis protein bacB



• Molecule 1: Bacilysin biosynthesis protein bacB



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	46.25Å 47.20Å 62.91Å 89.05° 77.27° 82.28°	Depositor
Resolution (Å)	30.34 – 2.04 30.35 – 2.04	Depositor EDS
% Data completeness (in resolution range)	93.3 (30.34-2.04) 93.3 (30.35-2.04)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.82 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.206 , 0.262 0.209 , 0.263	Depositor DCC
R_{free} test set	1537 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	24.6	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 45.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	3814	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.90% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: FE, PPY, CO

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	0.33	0/1838	0.51	0/2469
1	B	0.28	0/1838	0.48	0/2469
All	All	0.31	0/3676	0.49	0/4938

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

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	1804	0	1751	72	0
1	B	1804	0	1751	73	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	12	0	7	3	0
3	B	12	0	7	3	0
4	A	1	0	0	0	0
5	A	109	0	0	14	0
5	B	68	0	0	6	0
All	All	3814	0	3516	149	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 21.

All (149) 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:179:MSE:CE	1:B:188:MSE:HE2	1.69	1.22
1:A:165:ARG:HB3	5:A:337:HOH:O	1.38	1.19
1:B:159:MSE:HE3	1:B:162:HIS:ND1	1.56	1.18
1:B:159:MSE:CE	1:B:162:HIS:CE1	2.28	1.16
1:A:3:THR:HG23	1:A:6:ASP:H	1.01	1.15
1:A:12:PHE:HE1	1:A:179:MSE:CE	1.61	1.14
1:B:111:ALA:O	1:B:112:ASP:OD2	1.69	1.11
1:B:138:THR:CG2	1:B:149:LEU:HD22	1.80	1.11
1:B:179:MSE:CE	1:B:188:MSE:CE	2.27	1.10
1:B:138:THR:HG22	1:B:149:LEU:HB2	1.15	1.07
1:B:159:MSE:HE3	1:B:162:HIS:CE1	1.90	1.07
1:B:179:MSE:HE1	1:B:188:MSE:HE2	1.35	1.05
1:A:3:THR:HG22	1:A:6:ASP:CB	1.88	1.04
1:B:159:MSE:HE2	1:B:162:HIS:CE1	1.90	1.03
1:A:179:MSE:HE1	1:A:194:TYR:CB	1.86	1.03
1:B:138:THR:HG21	1:B:149:LEU:HD22	1.31	1.03
1:A:6:ASP:HB3	5:A:334:HOH:O	1.57	1.02
1:A:179:MSE:HE1	1:A:194:TYR:HB3	1.39	1.02
1:B:138:THR:CG2	1:B:149:LEU:HB2	1.93	0.98
1:A:3:THR:HG22	1:A:6:ASP:HB2	1.45	0.97
1:A:3:THR:HG22	1:A:6:ASP:CG	1.84	0.97
1:B:59:MSE:HE1	1:B:147:ILE:HG21	1.48	0.95
1:A:167:GLU:OE2	1:A:169:ILE:HD11	1.64	0.95
1:B:179:MSE:HE2	1:B:188:MSE:HE2	1.47	0.94
1:A:12:PHE:CE1	1:A:179:MSE:CE	2.51	0.92
1:A:3:THR:CG2	1:A:6:ASP:H	1.82	0.92
1:B:38:MSE:HE3	1:B:104:ILE:HD11	1.49	0.92
1:A:3:THR:HG23	1:A:6:ASP:N	1.86	0.91
1:A:59:MSE:HE1	1:A:147:ILE:HG21	1.52	0.90
1:A:12:PHE:HE1	1:A:179:MSE:HE2	1.36	0.89
1:B:179:MSE:HE1	1:B:188:MSE:CE	2.01	0.86
1:B:138:THR:HG21	1:B:149:LEU:CD2	2.04	0.86
1:B:179:MSE:HE2	1:B:188:MSE:SE	2.27	0.84
1:B:179:MSE:HE2	1:B:188:MSE:CE	2.05	0.84
1:A:12:PHE:CE1	1:A:179:MSE:HE3	2.12	0.83
1:B:179:MSE:CE	1:B:188:MSE:SE	2.75	0.83
1:A:38:MSE:HE3	1:A:104:ILE:HD11	1.59	0.83
1:B:38:MSE:HE3	1:B:104:ILE:CD1	2.08	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:246:PPY:H2'	3:B:246:PPY:O3	1.78	0.82
1:A:33:ASP:HB2	5:A:339:HOH:O	1.78	0.82
1:A:3:THR:CG2	1:A:6:ASP:HB2	2.10	0.82
1:A:12:PHE:HE1	1:A:179:MSE:HE3	1.43	0.81
1:B:138:THR:CG2	1:B:149:LEU:CD2	2.58	0.80
1:A:105:ASP:C	1:A:106:ILE:HD12	2.01	0.80
1:A:5:GLU:H	1:A:5:GLU:CD	1.84	0.78
3:A:246:PPY:H2'	3:A:246:PPY:O3	1.85	0.76
1:A:2:LYS:HG2	5:A:334:HOH:O	1.86	0.76
1:B:110:LYS:HE2	5:B:311:HOH:O	1.86	0.74
1:B:165:ARG:HG3	1:B:199:ARG:NH1	2.03	0.74
1:A:106:ILE:HD12	1:A:106:ILE:N	2.03	0.72
1:A:33:ASP:CB	5:A:339:HOH:O	2.36	0.72
1:B:38:MSE:CE	1:B:104:ILE:CD1	2.67	0.72
1:A:179:MSE:HE1	1:A:194:TYR:CG	2.24	0.72
1:B:159:MSE:HE3	1:B:162:HIS:HD1	1.54	0.72
1:A:159:MSE:HE3	3:A:246:PPY:H31	1.71	0.71
1:A:3:THR:CG2	1:A:6:ASP:CB	2.63	0.71
1:A:12:PHE:CE1	1:A:179:MSE:HE2	2.24	0.69
1:A:127:LYS:HE2	1:A:129:ARG:HB3	1.76	0.68
1:A:50:HIS:HE1	5:A:335:HOH:O	1.79	0.66
1:B:48:GLU:HG3	1:B:49:PRO:HD2	1.77	0.66
1:B:138:THR:HG22	1:B:149:LEU:CB	2.08	0.65
1:B:67:MSE:HE2	1:B:83:TYR:CZ	2.32	0.64
1:A:8:GLN:NE2	1:A:9:GLU:HG3	2.13	0.63
1:B:138:THR:HG23	1:B:149:LEU:HD22	1.76	0.63
1:B:179:MSE:HE3	1:B:188:MSE:SE	2.47	0.63
1:A:38:MSE:HE3	1:A:104:ILE:CD1	2.28	0.63
1:A:29:THR:HG23	5:A:279:HOH:O	1.99	0.62
1:A:8:GLN:HE22	1:A:9:GLU:HG3	1.65	0.62
1:B:179:MSE:HE1	1:B:194:TYR:CD2	2.34	0.61
1:A:4:LYS:HE3	5:A:301:HOH:O	1.99	0.61
1:B:19:GLU:HG2	1:B:25:ARG:HG2	1.82	0.61
1:A:179:MSE:CE	1:A:194:TYR:CB	2.74	0.60
1:B:179:MSE:HE1	1:B:194:TYR:HD2	1.66	0.60
1:A:106:ILE:N	1:A:106:ILE:CD1	2.63	0.60
1:A:38:MSE:CE	1:A:104:ILE:CD1	2.79	0.60
1:B:165:ARG:HG3	1:B:199:ARG:HH11	1.64	0.60
1:A:50:HIS:CE1	5:A:335:HOH:O	2.52	0.60
1:B:163:LYS:HB3	1:B:201:ASP:OD2	2.00	0.60
1:B:220:PRO:HB2	1:B:221:PRO:HD2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:246:PPY:O3	3:B:246:PPY:C2'	2.50	0.59
1:A:2:LYS:CE	1:A:7:MSE:HE3	2.34	0.58
1:B:105:ASP:C	1:B:106:ILE:HD12	2.24	0.58
1:B:111:ALA:O	1:B:112:ASP:CG	2.41	0.57
1:B:33:ASP:HB2	1:B:110:LYS:HZ3	1.70	0.57
1:A:135:MSE:HE2	1:A:152:ILE:HG12	1.85	0.56
1:B:59:MSE:CE	1:B:147:ILE:HG21	2.30	0.56
1:A:4:LYS:HG3	5:A:336:HOH:O	2.05	0.56
1:A:3:THR:OG1	1:A:5:GLU:OE2	2.23	0.55
1:A:33:ASP:N	5:A:339:HOH:O	2.22	0.55
1:B:33:ASP:HB2	1:B:110:LYS:NZ	2.22	0.55
1:B:25:ARG:HD2	5:B:275:HOH:O	2.08	0.54
1:B:25:ARG:CD	5:B:275:HOH:O	2.56	0.53
1:B:38:MSE:CE	1:B:104:ILE:HD13	2.37	0.53
1:B:135:MSE:HE2	1:B:152:ILE:HG23	1.90	0.52
1:A:8:GLN:HE22	1:A:9:GLU:CG	2.22	0.52
1:B:180:THR:HG22	5:B:249:HOH:O	2.08	0.52
1:B:161:PHE:CZ	1:B:201:ASP:HB3	2.44	0.52
1:A:3:THR:CG2	1:A:6:ASP:CG	2.71	0.51
1:A:3:THR:CG2	1:A:6:ASP:N	2.60	0.51
1:B:104:ILE:HD11	1:B:171:ILE:HD13	1.92	0.51
1:A:180:THR:HG22	5:A:251:HOH:O	2.11	0.51
1:A:220:PRO:HB2	1:A:221:PRO:HD2	1.92	0.51
1:A:5:GLU:CD	1:A:5:GLU:N	2.60	0.50
1:A:48:GLU:HA	1:A:48:GLU:OE2	2.11	0.49
1:A:167:GLU:OE2	1:A:169:ILE:CD1	2.49	0.49
1:B:29:THR:HG21	5:B:282:HOH:O	2.12	0.49
1:A:51:GLN:HE21	1:A:90:PRO:HG3	1.78	0.49
1:B:220:PRO:HB2	1:B:221:PRO:CD	2.42	0.48
1:B:59:MSE:HE1	1:B:147:ILE:CG2	2.32	0.48
1:B:2:LYS:HD3	1:B:7:MSE:HE3	1.94	0.48
1:B:55:VAL:HG22	1:B:56:GLN:N	2.28	0.48
3:A:246:PPY:O3	3:A:246:PPY:C2'	2.55	0.48
1:B:38:MSE:HE2	1:B:104:ILE:HD13	1.94	0.47
1:B:38:MSE:HE2	1:B:104:ILE:CD1	2.45	0.47
1:A:44:HIS:NE2	1:A:98:ASP:OD2	2.48	0.46
1:A:140:PHE:HB2	1:A:147:ILE:HB	1.96	0.46
1:B:161:PHE:CE1	1:B:201:ASP:HB3	2.51	0.46
1:B:48:GLU:HG3	1:B:49:PRO:CD	2.46	0.46
1:B:138:THR:CG2	1:B:149:LEU:CB	2.80	0.46
1:B:152:ILE:HD11	3:B:246:PPY:H5'	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:LEU:HD22	1:A:123:LEU:N	2.32	0.44
1:B:132:LEU:O	1:B:133:PRO:C	2.54	0.44
1:B:106:ILE:HD12	1:B:106:ILE:N	2.33	0.44
1:A:2:LYS:HE2	1:A:7:MSE:HE3	2.00	0.44
1:A:8:GLN:CD	1:A:8:GLN:C	2.76	0.44
1:A:94:ARG:NH2	5:A:343:HOH:O	2.36	0.44
1:B:107:LYS:HE3	1:B:107:LYS:HB2	1.65	0.43
1:A:137:VAL:HA	1:A:149:LEU:O	2.19	0.43
1:A:38:MSE:HE2	1:A:104:ILE:CD1	2.47	0.43
1:B:128:THR:O	1:B:128:THR:HG23	2.18	0.43
1:A:104:ILE:HD11	1:A:171:ILE:HD13	2.01	0.42
1:B:28:SER:HB2	1:B:37:LEU:CD2	2.49	0.42
1:A:48:GLU:OE2	1:A:48:GLU:CA	2.67	0.42
1:A:170:GLY:HA3	1:A:194:TYR:CE2	2.54	0.42
1:B:140:PHE:HB2	1:B:147:ILE:HB	2.02	0.42
1:A:70:GLY:O	5:A:322:HOH:O	2.22	0.42
1:A:56:GLN:HA	1:A:106:ILE:O	2.19	0.42
1:B:180:THR:HB	1:B:185:THR:OG1	2.20	0.42
1:B:142:GLU:HG3	1:B:143:ASP:N	2.34	0.42
1:A:55:VAL:HG12	1:A:108:ARG:O	2.19	0.42
1:B:179:MSE:CE	1:B:194:TYR:CD2	3.03	0.42
1:B:165:ARG:CG	1:B:199:ARG:HH11	2.33	0.41
1:A:58:GLY:HA3	1:A:83:TYR:CE2	2.54	0.41
1:A:179:MSE:HE1	1:A:194:TYR:CD2	2.56	0.41
1:B:56:GLN:HA	1:B:106:ILE:O	2.21	0.41
1:A:54:GLU:HB3	1:A:107:LYS:HG3	2.03	0.41
1:B:55:VAL:HG12	1:B:108:ARG:O	2.21	0.41
1:A:179:MSE:CE	1:A:194:TYR:HB3	2.28	0.41
1:B:25:ARG:HD3	5:B:275:HOH:O	2.19	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	223/243 (92%)	217 (97%)	6 (3%)	0	100	100
1	B	223/243 (92%)	217 (97%)	6 (3%)	0	100	100
All	All	446/486 (92%)	434 (97%)	12 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	199/201 (99%)	191 (96%)	8 (4%)	31	24
1	B	199/201 (99%)	190 (96%)	9 (4%)	27	20
All	All	398/402 (99%)	381 (96%)	17 (4%)	29	22

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

Mol	Chain	Res	Type
1	A	8	GLN
1	A	48	GLU
1	A	50	HIS
1	A	81	SER
1	A	104	ILE
1	A	149	LEU
1	A	163	LYS
1	A	180	THR
1	B	74	ARG
1	B	81	SER
1	B	99	GLN
1	B	104	ILE
1	B	149	LEU
1	B	169	ILE
1	B	180	THR
1	B	199	ARG
1	B	209	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	51	GLN
1	B	8	GLN
1	B	99	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 monosaccharides in this entry.

5.6 Ligand geometry

Of 7 ligands modelled in this entry, 5 are monoatomic - leaving 2 for Mogul analysis.

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	$\#\|Z\| > 2$
3	PPY	A	246	2	12,12,12	1.63	1 (8%)	15,15,15	1.14	2 (13%)
3	PPY	B	246	2	12,12,12	1.69	1 (8%)	15,15,15	0.85	0

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
3	PPY	A	246	2	-	2/8/8/8	0/1/1/1
3	PPY	B	246	2	-	2/8/8/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	246	PPY	C2-C1	-5.32	1.46	1.53
3	A	246	PPY	C2-C1	-5.21	1.46	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	246	PPY	O1-C1-C2	-3.28	117.33	121.72
3	A	246	PPY	O2-C1-C2	2.03	119.52	113.97

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	246	PPY	C2'-C1'-C3-C2
3	A	246	PPY	C2'-C1'-C3-C2
3	A	246	PPY	C6'-C1'-C3-C2
3	B	246	PPY	C6'-C1'-C3-C2

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	246	PPY	3	0
3	B	246	PPY	3	0

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, 95^{th} 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	213/243 (87%)	0.33	13 (6%) 21 22	6, 11, 18, 40	0
1	B	213/243 (87%)	0.62	26 (12%) 4 3	11, 17, 25, 37	0
All	All	426/486 (87%)	0.48	39 (9%) 9 9	6, 15, 24, 40	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	217	ILE	5.8
1	B	217	ILE	5.2
1	A	215	ILE	4.9
1	B	215	ILE	4.9
1	B	104	ILE	4.1
1	B	147	ILE	3.8
1	A	57	ILE	3.7
1	A	147	ILE	3.7
1	B	106	ILE	3.3
1	A	106	ILE	3.3
1	B	57	ILE	3.3
1	A	145	VAL	3.2
1	B	99	GLN	3.2
1	B	58	GLY	2.9
1	B	210	LYS	2.8
1	B	87	PRO	2.7
1	B	98	ASP	2.7
1	B	128	THR	2.7
1	B	112	ASP	2.7
1	B	149	LEU	2.7
1	B	209	GLU	2.7
1	A	36	VAL	2.6
1	B	88	HIS	2.5
1	A	104	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	169	ILE	2.5
1	A	126	PHE	2.5
1	B	36	VAL	2.5
1	B	171	ILE	2.4
1	A	171	ILE	2.3
1	A	216	ASN	2.3
1	B	216	ASN	2.3
1	B	121	TYR	2.2
1	B	199	ARG	2.2
1	B	170	GLY	2.2
1	B	169	ILE	2.1
1	B	82	ALA	2.1
1	A	170	GLY	2.0
1	B	89	VAL	2.0
1	B	225	ARG	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)

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, 95^{th} 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(Å^2)	Q<0.9
2	CO	B	244	1/1	0.91	0.08	60,60,60,60	0
3	PPY	A	246	12/12	0.95	0.16	15,19,23,23	0
3	PPY	B	246	12/12	0.97	0.13	21,23,27,27	0
4	FE	A	247	1/1	0.98	0.06	30,30,30,30	0
2	CO	B	245	1/1	0.99	0.07	23,23,23,23	0
2	CO	A	244	1/1	0.99	0.06	39,39,39,39	0
2	CO	A	245	1/1	1.00	0.07	15,15,15,15	0

6.5 Other polymers (i)

There are no such residues in this entry.