



Full wwPDB X-ray Structure Validation Report

Apr 19, 2017 – 10:46 AM EDT

PDB ID : 5JK5
Title : Phenylalanine hydroxylase from dictyostelium - BH2 complex
Authors : Zhuang, N.; Lee, K.H.
Deposited on : 2016-04-26
Resolution : 2.07 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the  symbol.

The following versions of software and data (see [references](#) ) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029077
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029077

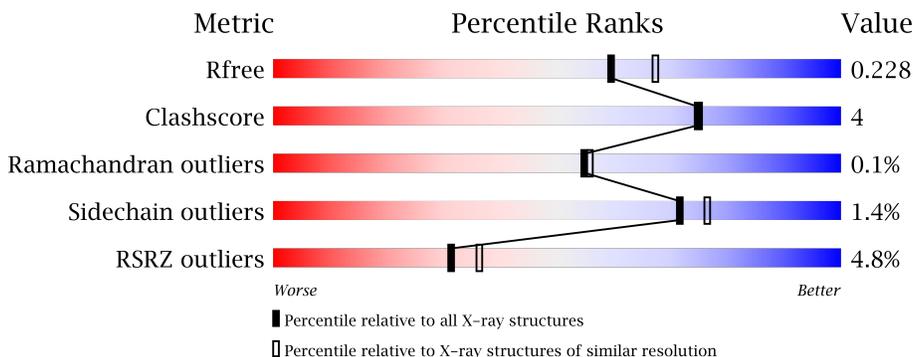
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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}	100719	4955 (2.10-2.06)
Clashscore	112137	5537 (2.10-2.06)
Ramachandran outliers	110173	5483 (2.10-2.06)
Sidechain outliers	110143	5484 (2.10-2.06)
RSRZ outliers	101464	4991 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	438	 2% 85% 6% 9%
1	B	438	 6% 79% 10% 11%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PIN	A	502	-	-	-	X
3	PIN	B	502	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6862 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 Phenylalanine-4-hydroxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	400	3245	2075	536	623	11	0	0	0
1	B	391	3174	2034	525	605	10	0	0	0

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP Q54XS1
A	-21	GLY	-	expression tag	UNP Q54XS1
A	-20	SER	-	expression tag	UNP Q54XS1
A	-19	SER	-	expression tag	UNP Q54XS1
A	-18	HIS	-	expression tag	UNP Q54XS1
A	-17	HIS	-	expression tag	UNP Q54XS1
A	-16	HIS	-	expression tag	UNP Q54XS1
A	-15	HIS	-	expression tag	UNP Q54XS1
A	-14	HIS	-	expression tag	UNP Q54XS1
A	-13	HIS	-	expression tag	UNP Q54XS1
A	-12	SER	-	expression tag	UNP Q54XS1
A	-11	SER	-	expression tag	UNP Q54XS1
A	-10	GLY	-	expression tag	UNP Q54XS1
A	-9	LEU	-	expression tag	UNP Q54XS1
A	-8	VAL	-	expression tag	UNP Q54XS1
A	-7	PRO	-	expression tag	UNP Q54XS1
A	-6	ARG	-	expression tag	UNP Q54XS1
A	-5	GLY	-	expression tag	UNP Q54XS1
A	-4	SER	-	expression tag	UNP Q54XS1
A	-3	HIS	-	expression tag	UNP Q54XS1
A	-2	MET	-	expression tag	UNP Q54XS1
A	-1	ALA	-	expression tag	UNP Q54XS1
A	0	SER	-	expression tag	UNP Q54XS1
B	-22	MET	-	expression tag	UNP Q54XS1
B	-21	GLY	-	expression tag	UNP Q54XS1

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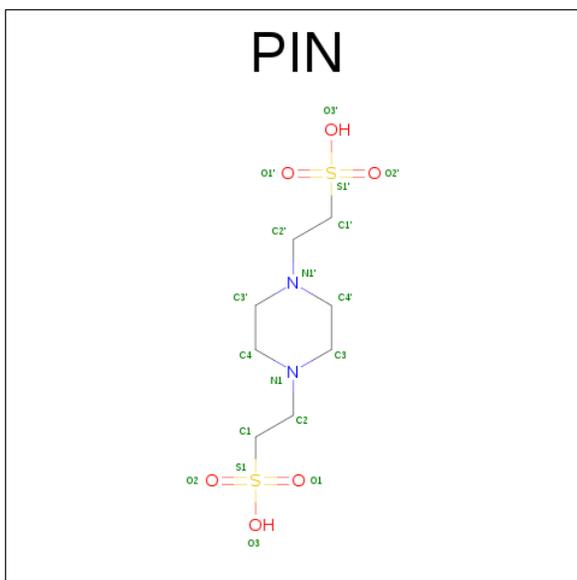
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	SER	-	expression tag	UNP Q54XS1
B	-19	SER	-	expression tag	UNP Q54XS1
B	-18	HIS	-	expression tag	UNP Q54XS1
B	-17	HIS	-	expression tag	UNP Q54XS1
B	-16	HIS	-	expression tag	UNP Q54XS1
B	-15	HIS	-	expression tag	UNP Q54XS1
B	-14	HIS	-	expression tag	UNP Q54XS1
B	-13	HIS	-	expression tag	UNP Q54XS1
B	-12	SER	-	expression tag	UNP Q54XS1
B	-11	SER	-	expression tag	UNP Q54XS1
B	-10	GLY	-	expression tag	UNP Q54XS1
B	-9	LEU	-	expression tag	UNP Q54XS1
B	-8	VAL	-	expression tag	UNP Q54XS1
B	-7	PRO	-	expression tag	UNP Q54XS1
B	-6	ARG	-	expression tag	UNP Q54XS1
B	-5	GLY	-	expression tag	UNP Q54XS1
B	-4	SER	-	expression tag	UNP Q54XS1
B	-3	HIS	-	expression tag	UNP Q54XS1
B	-2	MET	-	expression tag	UNP Q54XS1
B	-1	ALA	-	expression tag	UNP Q54XS1
B	0	SER	-	expression tag	UNP Q54XS1

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

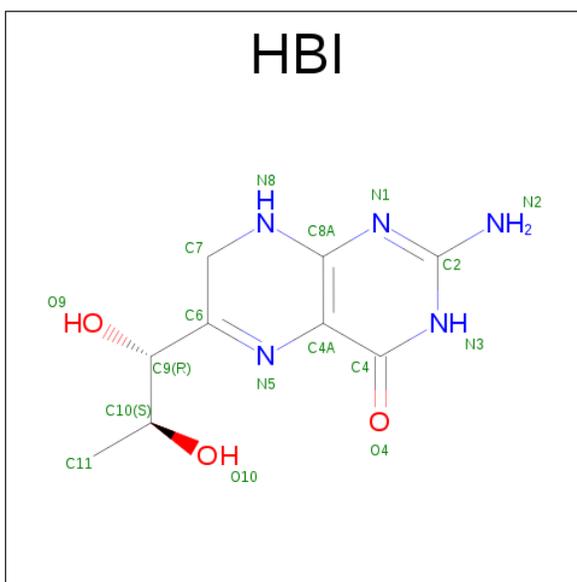
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Fe 1 1	0	0
2	A	1	Total Fe 1 1	0	0

- Molecule 3 is PIPERAZINE-N,N'-BIS(2-ETHANESULFONIC ACID) (three-letter code: PIN) (formula: C₈H₁₈N₂O₆S₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			S
3	A	1	18	8	2	6	2	0	0
3	B	1	18	8	2	6	2	0	0

- Molecule 4 is 7,8-DIHYDROBIOPTERIN (three-letter code: HBI) (formula: $C_9H_{13}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	17	9	5	3	0	0
4	B	1	17	9	5	3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	208	Total 208	O 208	0	0
5	B	163	Total 163	O 163	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.97Å 85.33Å 74.89Å 90.00° 110.23° 90.00°	Depositor
Resolution (Å)	37.45 – 2.07 37.45 – 2.07	Depositor EDS
% Data completeness (in resolution range)	99.1 (37.45-2.07) 99.1 (37.45-2.07)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.06Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.165 , 0.228 0.165 , 0.228	Depositor DCC
R_{free} test set	2443 reflections (4.83%)	DCC
Wilson B-factor (Å ²)	30.9	Xtrriage
Anisotropy	0.132	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6862	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HBI, FE, PIN

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	0.39	0/3331	0.51	2/4516 (0.0%)
1	B	0.37	0/3258	0.50	0/4414
All	All	0.38	0/6589	0.50	2/8930 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	VAL	C-N-CA	5.31	134.98	121.70
1	A	68	LEU	CA-CB-CG	5.31	127.51	115.30

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	3245	0	3135	20	0
1	B	3174	0	3071	26	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	18	0	18	2	0
3	B	18	0	18	2	0
4	A	17	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	17	0	13	0	0
5	A	208	0	0	6	0
5	B	163	0	0	4	1
All	All	6862	0	6268	47	1

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 4.

All (47) 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:123:VAL:HA	1:A:124:LEU:HB2	1.52	0.89
1:B:249:ARG:NH2	1:B:289:ASP:OD1	2.25	0.67
1:B:218:ILE:HD13	1:B:275:LEU:HD22	1.78	0.65
1:B:111:ARG:O	5:B:601:HOH:O	2.13	0.65
1:B:335:ILE:HG22	1:B:342:MET:HG2	1.80	0.62
1:B:213:PRO:HB2	1:B:218:ILE:HD11	1.82	0.62
1:B:137:ALA:O	1:B:143:ARG:NH2	2.32	0.61
1:B:181:ARG:HG3	1:B:339:THR:HG21	1.87	0.54
1:A:43:LYS:O	5:A:601:HOH:O	2.19	0.53
1:A:213:PRO:HB2	1:A:218:ILE:HD11	1.91	0.52
1:A:123:VAL:HG11	1:A:234:GLN:HB2	1.92	0.51
1:A:367:ILE:HG13	1:A:368:THR:HG23	1.94	0.50
1:B:143:ARG:HG2	1:B:147:ARG:NH2	2.28	0.49
1:B:194:HIS:CD2	1:B:336:LEU:HD13	2.47	0.49
1:B:49:ILE:HD13	1:B:52:ILE:HD11	1.95	0.48
1:A:47:ILE:HG12	5:A:722:HOH:O	2.15	0.47
1:B:29:SER:HB3	1:B:32:ILE:HG13	1.95	0.47
1:A:259:HIS:CG	1:A:260:PRO:HD2	2.50	0.46
1:B:259:HIS:CG	1:B:260:PRO:HD2	2.50	0.46
1:B:270:ASP:O	1:B:274:GLU:HG2	2.15	0.46
1:A:405:TYR:CZ	1:B:404:PRO:HB3	2.50	0.46
1:B:73:PRO:HG2	1:B:74:THR:HG23	1.97	0.46
1:B:219:SER:O	1:B:223:GLN:HG3	2.14	0.46
3:A:502:PIN:H1'2	3:A:502:PIN:H4'2	1.65	0.46
1:B:367:ILE:HG13	1:B:368:THR:HG23	1.97	0.46
1:A:30:ASP:OD2	5:A:602:HOH:O	2.21	0.46
1:A:304:GLU:HG3	5:A:777:HOH:O	2.16	0.45
1:A:58:LYS:NZ	3:B:502:PIN:O1'	2.49	0.44
1:B:224:GLU:OE1	5:B:602:HOH:O	2.21	0.44
1:A:346:LEU:HB2	5:A:705:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:ARG:HG2	1:B:147:ARG:HH21	1.82	0.44
1:A:321:LEU:HD13	1:A:328:ILE:HG21	1.99	0.43
1:B:23:PHE:HE1	1:B:35:LEU:HD21	1.83	0.43
1:A:123:VAL:HA	1:A:124:LEU:CB	2.37	0.42
1:B:23:PHE:CE1	1:B:35:LEU:HD21	2.54	0.42
1:A:229:ARG:HG2	5:A:670:HOH:O	2.19	0.42
1:A:123:VAL:CA	1:A:124:LEU:HB2	2.36	0.42
1:A:289:ASP:OD2	3:A:502:PIN:H41	2.20	0.41
3:B:502:PIN:H12	5:B:713:HOH:O	2.20	0.41
1:B:392:ASP:OD1	5:B:603:HOH:O	2.22	0.41
1:A:228:TRP:CD1	1:A:288:ALA:HB1	2.56	0.41
1:A:335:ILE:O	1:A:342:MET:HG2	2.19	0.41
1:B:233:VAL:HB	1:B:252:HIS:HB3	2.03	0.41
1:B:212:ILE:HA	1:B:213:PRO:HD3	1.90	0.41
1:A:123:VAL:HG11	1:A:234:GLN:CB	2.50	0.40
1:B:259:HIS:CE1	1:B:261:SER:HG	2.34	0.40
1:B:30:ASP:N	1:B:30:ASP:OD1	2.31	0.40

All (1) 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)
5:B:674:HOH:O	5:B:754:HOH:O[2_756]	2.03	0.17

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	398/438 (91%)	391 (98%)	6 (2%)	1 (0%)	44 43
1	B	387/438 (88%)	376 (97%)	11 (3%)	0	100 100
All	All	785/876 (90%)	767 (98%)	17 (2%)	1 (0%)	55 56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	LEU

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	358/390 (92%)	354 (99%)	4 (1%)	78	82
1	B	348/390 (89%)	342 (98%)	6 (2%)	66	70
All	All	706/780 (90%)	696 (99%)	10 (1%)	71	76

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	124	LEU
1	A	166	ASP
1	A	348	ASP
1	B	30	ASP
1	B	41	ILE
1	B	103	GLN
1	B	208	SER
1	B	327	THR
1	B	413	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	PIN	A	502	-	18,18,18	2.03	6 (33%)	22,26,26	1.60	6 (27%)
4	HBI	A	503	-	13,18,18	2.17	3 (23%)	10,26,26	3.05	5 (50%)
3	PIN	B	502	-	18,18,18	2.00	6 (33%)	22,26,26	1.79	6 (27%)
4	HBI	B	503	-	13,18,18	2.17	3 (23%)	10,26,26	3.11	5 (50%)

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	PIN	A	502	-	-	0/12/22/22	0/1/1/1
4	HBI	A	503	-	-	0/4/17/17	0/2/2/2
3	PIN	B	502	-	-	0/12/22/22	0/1/1/1
4	HBI	B	503	-	-	0/4/17/17	0/2/2/2

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	PIN	O2'-S1'	2.66	1.52	1.45
3	A	502	PIN	O2-S1	2.68	1.52	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	503	HBI	C2-N3	2.76	1.40	1.35
3	A	502	PIN	O1'-S1'	2.78	1.53	1.45
3	B	502	PIN	O1-S1	2.83	1.53	1.45
3	B	502	PIN	O2-S1	2.87	1.53	1.45
3	B	502	PIN	O1'-S1'	2.92	1.53	1.45
4	A	503	HBI	C2-N3	2.97	1.40	1.35
3	A	502	PIN	O1-S1	2.98	1.53	1.45
3	A	502	PIN	O2'-S1'	2.98	1.53	1.45
3	A	502	PIN	C1'-S1'	3.90	1.83	1.77
3	B	502	PIN	C1'-S1'	3.97	1.83	1.77
4	A	503	HBI	C4-N3	4.12	1.40	1.33
3	B	502	PIN	C1-S1	4.15	1.83	1.77
3	A	502	PIN	C1-S1	4.32	1.83	1.77
4	B	503	HBI	C4-N3	4.49	1.41	1.33
4	B	503	HBI	C6-N5	5.25	1.35	1.28
4	A	503	HBI	C6-N5	5.34	1.35	1.28

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	503	HBI	C4A-C4-N3	-3.82	118.04	123.48
4	B	503	HBI	C4A-C4-N3	-3.68	118.25	123.48
4	B	503	HBI	N3-C2-N1	-2.84	120.85	125.45
4	A	503	HBI	N3-C2-N1	-2.52	121.37	125.45
3	B	502	PIN	O3'-S1'-C1'	2.08	108.61	106.06
3	A	502	PIN	O1-S1-C1	2.12	108.61	106.79
4	B	503	HBI	C4-N3-C2	2.17	119.18	116.06
4	A	503	HBI	C4-N3-C2	2.27	119.32	116.06
3	A	502	PIN	O2-S1-C1	2.44	108.89	106.79
3	A	502	PIN	O3'-S1'-C1'	2.57	109.21	106.06
4	A	503	HBI	C2-N1-C8A	2.57	120.29	114.51
3	B	502	PIN	O3-S1-C1	2.62	109.28	106.06
4	B	503	HBI	C2-N1-C8A	2.87	120.98	114.51
3	A	502	PIN	O1'-S1'-C1'	2.92	109.30	106.79
3	B	502	PIN	O2-S1-C1	3.24	109.57	106.79
3	A	502	PIN	O3-S1-C1	3.31	110.13	106.06
3	A	502	PIN	O2'-S1'-C1'	3.35	109.67	106.79
3	B	502	PIN	O1-S1-C1	3.52	109.82	106.79
3	B	502	PIN	O2'-S1'-C1'	3.62	109.90	106.79
3	B	502	PIN	O1'-S1'-C1'	3.64	109.92	106.79
4	A	503	HBI	C4-C4A-C8A	7.49	119.52	114.52
4	B	503	HBI	C4-C4A-C8A	7.56	119.57	114.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	PIN	2	0
3	B	502	PIN	2	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, 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	400/438 (91%)	-0.15	10 (2%) 58 62	20, 35, 65, 87	0
1	B	391/438 (89%)	0.21	28 (7%) 16 20	18, 40, 77, 101	0
All	All	791/876 (90%)	0.03	38 (4%) 31 36	18, 37, 73, 101	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	138	SER	4.5
1	B	264	LEU	4.5
1	B	137	ALA	4.3
1	B	47	ILE	3.9
1	A	414	ASN	3.9
1	B	262	VAL	3.8
1	A	124	LEU	3.7
1	B	142	TYR	3.6
1	B	143	ARG	3.6
1	B	265	TYR	3.5
1	B	261	SER	3.4
1	B	134	HIS	3.3
1	A	128	SER	3.2
1	A	76	ASN	3.2
1	B	267	PRO	3.1
1	B	141	VAL	3.0
1	B	25	ILE	2.9
1	B	45	HIS	2.9
1	B	30	ASP	2.9
1	B	123	VAL	2.8
1	B	263	PRO	2.8
1	A	78	LYS	2.7
1	B	59	THR	2.7
1	B	28	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	259	HIS	2.6
1	A	126	MET	2.6
1	B	78	LYS	2.5
1	A	77	ASN	2.5
1	B	135	PRO	2.5
1	B	136	GLY	2.4
1	A	326	ASP	2.4
1	A	101	SER	2.3
1	A	395	LYS	2.2
1	B	139	ASP	2.2
1	B	80	VAL	2.2
1	B	42	ILE	2.1
1	B	140	PRO	2.1
1	B	73	PRO	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	PIN	B	502	18/18	0.89	0.21	2.26	67,76,119,119	3
3	PIN	A	502	18/18	0.92	0.17	2.17	55,63,69,70	1
4	HBI	B	503	17/17	0.93	0.10	-0.42	33,39,52,55	0
4	HBI	A	503	17/17	0.93	0.11	-0.56	27,36,45,52	0
2	FE	A	501	1/1	0.99	0.13	-0.77	28,28,28,28	0
2	FE	B	501	1/1	0.99	0.10	-0.78	28,28,28,28	0

6.5 Other polymers

There are no such residues in this entry.