



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2016 – 08:21 AM EST

PDB ID : 5KF6
Title : Structure of proline utilization A from Sinorhizobium meliloti complexed with L-tetrahydrofuroic acid and NAD⁺ in space group P21
Authors : Tanner, J.J.
Deposited on : 2016-06-12
Resolution : 1.70 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028320
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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-20028320

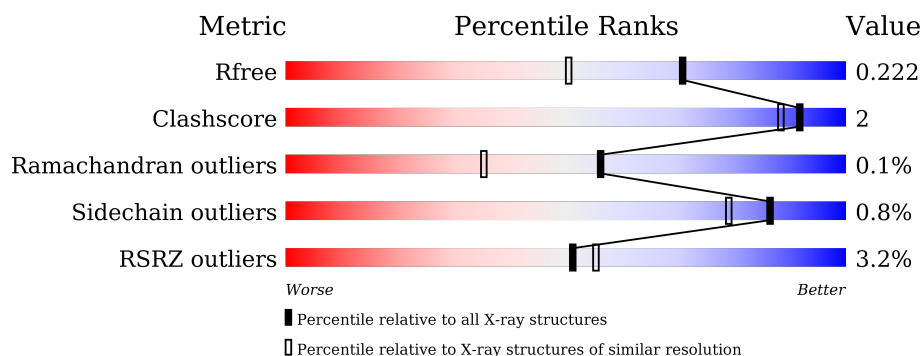
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	91344	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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	1235	<div> <div>2%</div> <div>93%</div> <div>5%</div> </div>
1	B	1235	<div> <div>4%</div> <div>91%</div> <div>6%</div> </div>

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
6	PGE	A	2005	-	-	-	X
6	PGE	A	2006	-	-	-	X
6	PGE	A	2007	-	-	-	X
6	PGE	B	2005	-	-	-	X
6	PGE	B	2009	-	-	-	X
7	PG4	B	2010	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 19406 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.

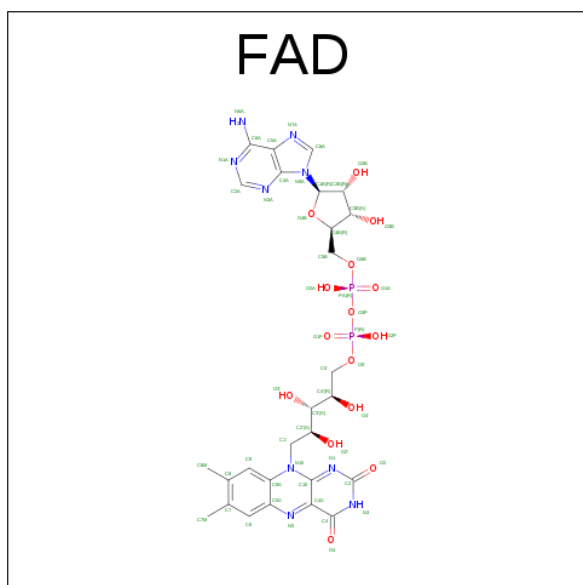
- Molecule 1 is a protein called Bifunctional protein PutA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1207	Total	C	N	O	S	0	3	0
			8948	5632	1605	1677	34			
1	B	1200	Total	C	N	O	S	0	8	0
			8879	5595	1582	1670	32			

There are 4 discrepancies between the modelled and reference sequences:

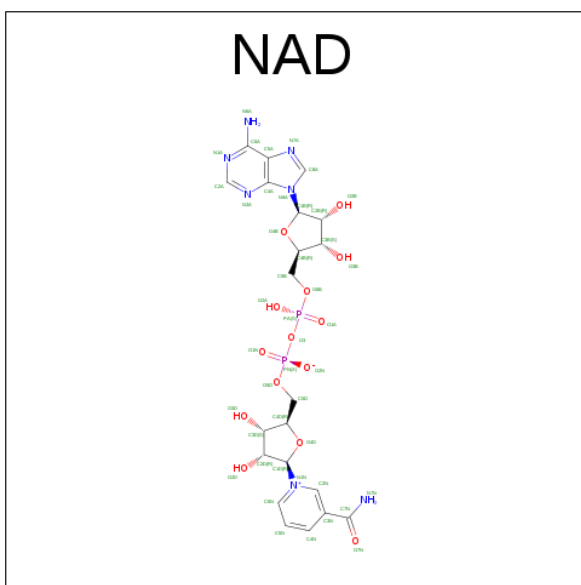
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP F7X6I3
A	0	MET	-	expression tag	UNP F7X6I3
B	-1	SER	-	expression tag	UNP F7X6I3
B	0	MET	-	expression tag	UNP F7X6I3

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



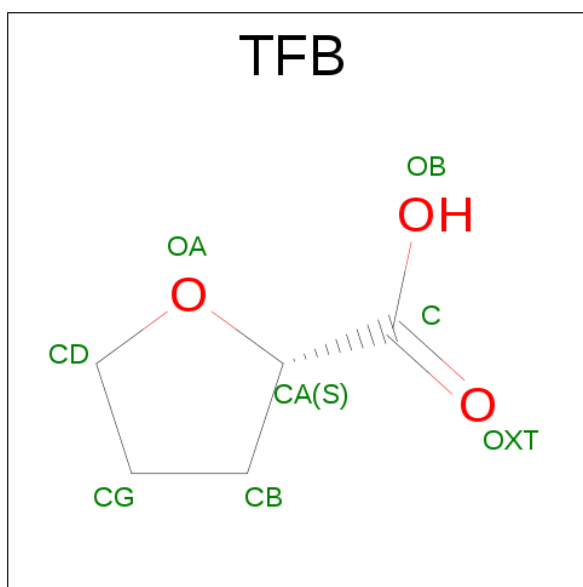
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is TETRAHYDROFURAN-2-CARBOXYLIC ACID (three-letter code: TFB) (formula: $C_5H_8O_3$).

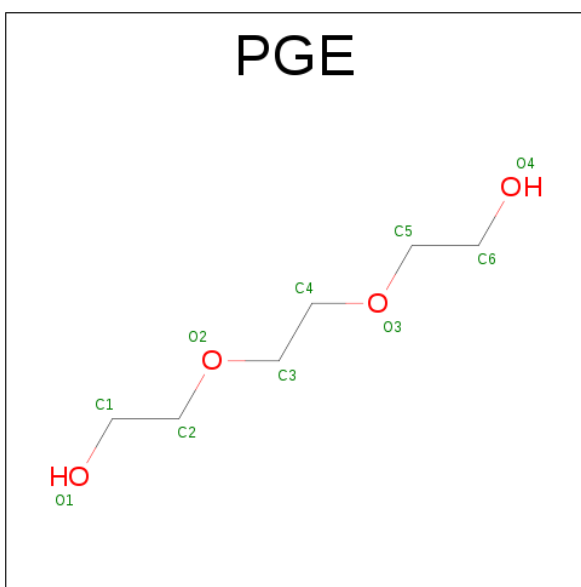


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	5	3		
4	B	1	Total	C	O	0	0
			8	5	3		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

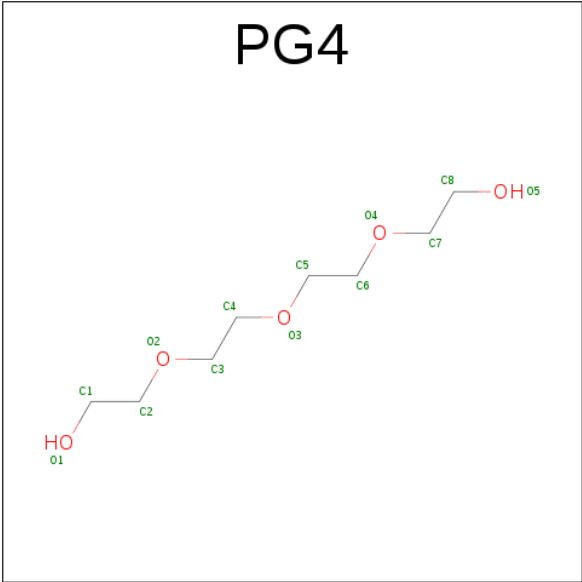
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		
5	A	1	Total	Mg	0	0
			1	1		

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			13	8	5		

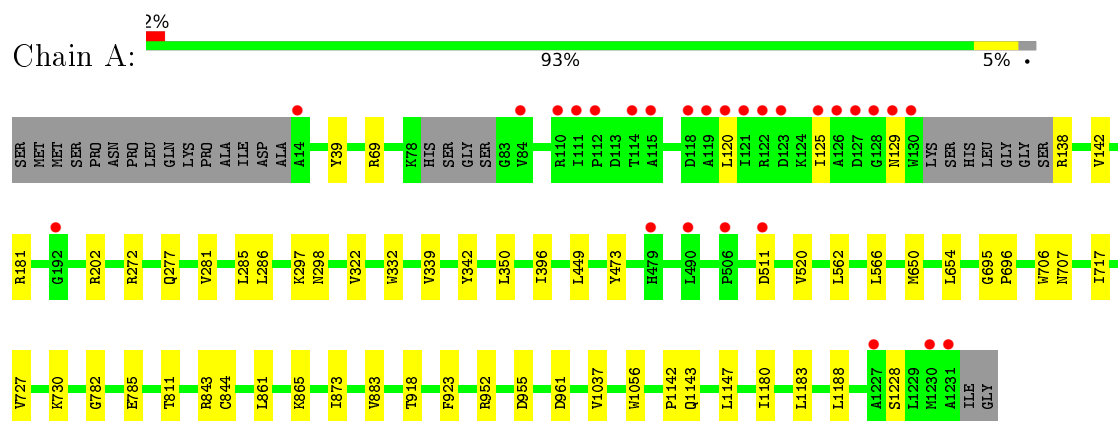
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	626	Total	O	0	0
			626	626		
8	B	660	Total	O	0	0
			660	660		

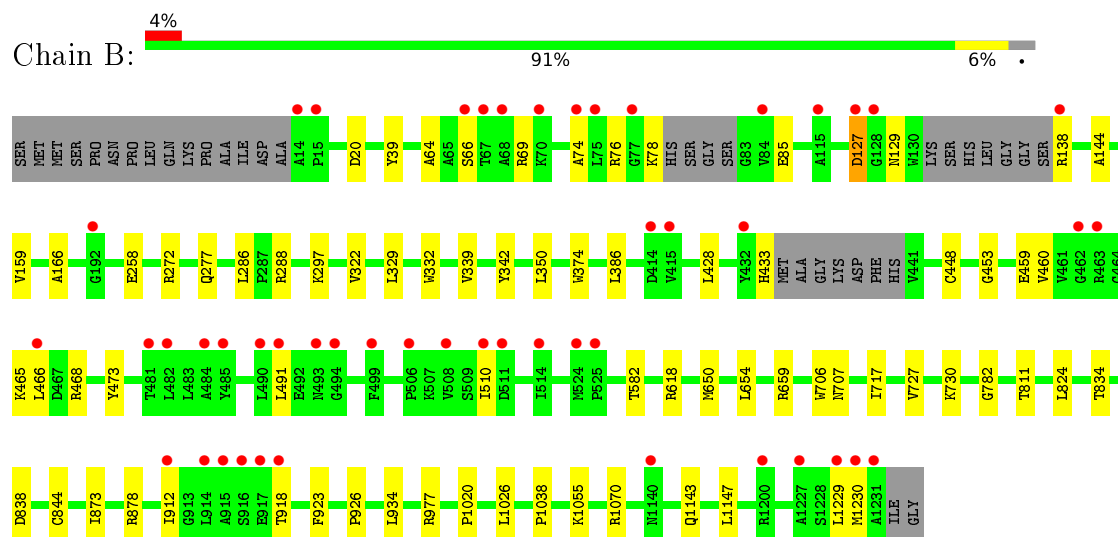
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Bifunctional protein PutA



• Molecule 1: Bifunctional protein PutA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	101.40 Å 102.33 Å 125.91 Å 90.00° 106.53° 90.00°	Depositor
Resolution (Å)	51.38 – 1.70 51.38 – 1.70	Depositor EDS
% Data completeness (in resolution range)	97.8 (51.38-1.70) 97.9 (51.38-1.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 1.70 Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.190 , 0.222 0.190 , 0.222	Depositor DCC
R_{free} test set	13377 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	15.8	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.1	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	19406	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.62% 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: MG, PGE, NAD, TFB, PG4, FAD

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	0.36	0/9113	0.55	0/12399
1	B	0.36	0/9061	0.56	0/12337
All	All	0.36	0/18174	0.55	0/24736

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 [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	8948	0	8978	34	0
1	B	8879	0	8907	42	0
2	A	53	0	31	2	0
2	B	53	0	31	2	0
3	A	44	0	26	1	0
3	B	44	0	26	2	0
4	A	8	0	7	0	0
4	B	8	0	7	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	24	0	32	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	44	0	60	2	0
7	B	13	0	18	0	0
8	A	626	0	0	3	0
8	B	660	0	0	5	0
All	All	19406	0	18123	75	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 2.

All (75) 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:582:THR:HA	6:B:2005:PGE:H5	1.76	0.67
1:A:286:LEU:HD21	1:A:322:VAL:HG11	1.76	0.67
1:A:138:ARG:HE	1:A:142:VAL:HG21	1.61	0.66
1:B:844:CYS:SG	3:B:2002:NAD:C4N	2.83	0.66
1:B:20:ASP:OD2	1:B:878:ARG:NH1	2.29	0.65
1:B:473:TYR:HB2	2:B:2001:FAD:HM72	1.80	0.63
1:A:961:ASP:OD2	1:B:1055:LYS:NZ	2.30	0.62
1:A:861:LEU:HD22	1:A:865:LYS:HE3	1.81	0.62
1:A:844:CYS:SG	3:A:2002:NAD:C4N	2.87	0.62
1:A:873:ILE:HG13	1:A:883:VAL:HB	1.82	0.61
1:B:297:LYS:HG3	1:B:332:TRP:HB2	1.81	0.60
1:B:288[B]:ARG:NE	8:B:2104:HOH:O	2.35	0.59
1:B:339:VAL:HG21	1:B:350:LEU:HD21	1.85	0.59
1:A:473:TYR:HB2	2:A:2001:FAD:HM72	1.86	0.58
1:B:286:LEU:HD21	1:B:322:VAL:HG11	1.85	0.58
1:A:339:VAL:HG21	1:A:350:LEU:HD21	1.85	0.57
1:B:650:MET:O	1:B:654:LEU:HG	2.04	0.57
1:B:297:LYS:HD2	1:B:329:LEU:HA	1.86	0.56
1:A:298:ASN:ND2	8:A:2104:HOH:O	2.35	0.55
1:A:69:ARG:NH2	1:A:511:ASP:OD1	2.40	0.54
1:A:449:LEU:HD23	1:A:473:TYR:HB3	1.90	0.53
1:B:459:GLU:OE1	1:B:465:LYS:NZ	2.42	0.53
1:B:1143:GLN:O	1:B:1147:LEU:HG	2.11	0.51
1:A:281:VAL:HG13	1:A:285:LEU:HD23	1.93	0.51
1:B:69:ARG:HG3	1:B:510:ILE:HG21	1.92	0.50
1:B:127:ASP:N	1:B:127:ASP:OD1	2.45	0.49
1:A:1180:ILE:HG23	1:A:1188:LEU:HD12	1.94	0.49
1:B:448:CYS:HB2	1:B:453:GLY:HA3	1.94	0.48
1:A:650:MET:O	1:A:654:LEU:HG	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:952:ARG:NH2	1:A:955:ASP:OD2	2.47	0.48
1:A:706:TRP:CE3	1:A:707:ASN:HA	2.49	0.48
1:B:64:ALA:HA	1:B:433:HIS:CD2	2.49	0.47
1:B:144:ALA:HB3	1:B:159:VAL:HG11	1.97	0.47
1:A:785:GLU:H	1:A:785:GLU:CD	2.17	0.47
1:B:374:TRP:HZ3	1:B:1229:LEU:HB3	1.80	0.47
1:A:843:ARG:HH12	6:A:2006:PGE:H12	1.80	0.46
1:A:202:ARG:NH2	8:A:2113:HOH:O	2.47	0.46
1:A:782:GLY:O	1:A:811:THR:HA	2.16	0.46
1:B:466:LEU:O	1:B:468:ARG:NH1	2.49	0.46
1:A:1056:TRP:CD1	1:A:1142:PRO:HD3	2.51	0.46
1:B:844:CYS:SG	3:B:2002:NAD:C3N	3.04	0.46
1:A:1183:LEU:O	8:A:2101:HOH:O	2.20	0.45
1:A:449:LEU:HG	2:A:2001:FAD:HM82	1.98	0.45
1:B:918:THR:HB	1:B:923:PHE:CD1	2.52	0.44
1:A:1147:LEU:HD22	1:B:1147:LEU:HD22	2.00	0.44
1:A:272:ARG:HB3	1:A:277:GLN:HG3	1.99	0.44
1:A:120:LEU:O	1:A:125:ILE:HG12	2.17	0.44
1:B:428:LEU:HD11	1:B:460:VAL:HG21	2.00	0.43
1:A:918:THR:HB	1:A:923:PHE:CD1	2.54	0.43
1:B:824:LEU:HB2	1:B:977:ARG:HH21	1.83	0.43
1:A:396:ILE:HD11	1:A:520:VAL:HB	2.01	0.43
6:B:2008:PGE:H32	6:B:2008:PGE:H52	1.60	0.42
1:B:1026:LEU:HD23	1:B:1038:PRO:HG2	2.00	0.42
1:A:1037:VAL:HG11	1:B:166:ALA:HB1	2.01	0.42
1:A:717:ILE:HG12	1:A:727:VAL:HG11	2.02	0.42
1:B:74:ALA:O	1:B:78:LYS:HE3	2.19	0.42
1:B:706:TRP:CE3	1:B:707:ASN:HA	2.55	0.42
1:B:873:ILE:HD11	1:B:912:ILE:HD11	2.01	0.42
2:B:2001:FAD:H1'1	2:B:2001:FAD:H9	1.82	0.41
1:B:491:LEU:HD12	1:B:1230:MET:SD	2.60	0.41
1:B:717:ILE:HG12	1:B:727:VAL:HG11	2.03	0.41
1:B:834:THR:O	1:B:838:ASP:HB3	2.20	0.41
1:B:659:ARG:HD3	8:B:2387:HOH:O	2.19	0.41
1:B:272:ARG:HB3	1:B:277:GLN:HG3	2.02	0.41
1:B:460:VAL:HG22	1:B:466:LEU:HD12	2.02	0.41
1:B:618:ARG:NH1	8:B:2106:HOH:O	2.40	0.41
1:B:1070:ARG:HD3	8:B:2384:HOH:O	2.20	0.41
1:B:782:GLY:O	1:B:811:THR:HA	2.21	0.41
1:A:562:LEU:O	1:A:566:LEU:HG	2.21	0.41
1:A:1143:GLN:O	1:A:1147:LEU:HG	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:912:ILE:HD11	1:B:926:PRO:HD2	2.03	0.41
1:B:386:LEU:HA	1:B:386:LEU:HD23	1.92	0.40
1:A:695:GLY:HA2	1:A:696:PRO:HD3	1.95	0.40
1:A:297:LYS:HA	1:A:332:TRP:CD1	2.56	0.40
1:B:1020:PRO:HA	8:B:2164:HOH:O	2.21	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	1204/1235 (98%)	1178 (98%)	25 (2%)	1 (0%)	56	35
1	B	1202/1235 (97%)	1172 (98%)	29 (2%)	1 (0%)	56	35
All	All	2406/2470 (97%)	2350 (98%)	54 (2%)	2 (0%)	56	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	129	ASN
1	B	129	ASN

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	897/951 (94%)	892 (99%)	5 (1%)	90	85
1	B	892/951 (94%)	882 (99%)	10 (1%)	80	69
All	All	1789/1902 (94%)	1774 (99%)	15 (1%)	86	79

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	TYR
1	A	181	ARG
1	A	342	TYR
1	A	730	LYS
1	A	1228	SER
1	B	39	TYR
1	B	66	SER
1	B	76	ARG
1	B	85	GLU
1	B	127	ASP
1	B	138	ARG
1	B	258	GLU
1	B	342	TYR
1	B	730	LYS
1	B	934	LEU

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

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 17 ligands modelled in this entry, 2 are monoatomic - leaving 15 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$
2	FAD	A	2001	-	52,58,58	1.70	14 (26%)	52,89,89	2.26	8 (15%)
3	NAD	A	2002	5	42,48,48	1.69	10 (23%)	46,73,73	2.26	6 (13%)
4	TFB	A	2003	-	5,8,8	2.40	1 (20%)	2,10,10	0.48	0
6	PGE	A	2005	-	9,9,9	0.62	0	8,8,8	0.78	0
6	PGE	A	2006	-	6,6,9	0.58	0	5,5,8	0.89	0
6	PGE	A	2007	-	6,6,9	0.55	0	5,5,8	0.93	0
2	FAD	B	2001	-	52,58,58	1.71	13 (25%)	52,89,89	2.39	11 (21%)
3	NAD	B	2002	5	42,48,48	1.78	9 (21%)	46,73,73	2.08	3 (6%)
4	TFB	B	2003	-	5,8,8	2.31	1 (20%)	2,10,10	0.51	0
6	PGE	B	2005	-	9,9,9	0.69	0	8,8,8	0.77	0
6	PGE	B	2006	-	6,6,9	0.54	0	5,5,8	0.70	0
6	PGE	B	2007	-	9,9,9	0.67	0	8,8,8	0.71	0
6	PGE	B	2008	-	9,9,9	0.68	0	8,8,8	0.86	0
6	PGE	B	2009	-	6,6,9	0.51	0	5,5,8	0.63	0
7	PG4	B	2010	-	12,12,12	0.56	0	11,11,11	0.89	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
2	FAD	A	2001	-	-	0/30/50/50	0/6/6/6
3	NAD	A	2002	5	-	0/22/62/62	0/5/5/5
4	TFB	A	2003	-	-	0/0/11/11	0/1/1/1
6	PGE	A	2005	-	-	0/7/7/7	0/0/0/0
6	PGE	A	2006	-	-	0/4/4/7	0/0/0/0
6	PGE	A	2007	-	-	0/4/4/7	0/0/0/0
2	FAD	B	2001	-	-	0/30/50/50	0/6/6/6
3	NAD	B	2002	5	-	0/22/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	TFB	B	2003	-	-	0/0/11/11	0/1/1/1
6	PGE	B	2005	-	-	0/7/7/7	0/0/0/0
6	PGE	B	2006	-	-	0/4/4/7	0/0/0/0
6	PGE	B	2007	-	-	0/7/7/7	0/0/0/0
6	PGE	B	2008	-	-	0/7/7/7	0/0/0/0
6	PGE	B	2009	-	-	0/4/4/7	0/0/0/0
7	PG4	B	2010	-	-	0/10/10/10	0/0/0/0

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2003	TFB	CB-CA	-4.92	1.45	1.53
4	B	2003	TFB	CB-CA	-4.67	1.46	1.53
3	B	2002	NAD	C2B-C1B	-3.11	1.48	1.53
3	A	2002	NAD	C2B-C1B	-2.88	1.49	1.53
2	A	2001	FAD	C2B-C1B	-2.87	1.49	1.53
2	B	2001	FAD	C2B-C1B	-2.66	1.49	1.53
2	B	2001	FAD	PA-O2A	-2.60	1.44	1.55
3	A	2002	NAD	PA-O2A	-2.44	1.44	1.55
2	A	2001	FAD	O2'-C2'	-2.44	1.37	1.43
3	B	2002	NAD	O3D-C3D	-2.43	1.37	1.43
2	A	2001	FAD	PA-O2A	-2.42	1.44	1.55
3	B	2002	NAD	C2D-C1D	-2.39	1.49	1.53
2	A	2001	FAD	PA-O5B	-2.31	1.49	1.59
2	B	2001	FAD	O2'-C2'	-2.29	1.38	1.43
3	A	2002	NAD	PA-O5B	-2.26	1.49	1.59
3	B	2002	NAD	PA-O2A	-2.18	1.45	1.55
2	B	2001	FAD	O2B-C2B	-2.14	1.37	1.43
3	A	2002	NAD	O4D-C4D	-2.07	1.40	1.45
3	B	2002	NAD	PA-O5B	-2.06	1.50	1.59
2	A	2001	FAD	O4B-C4B	-2.01	1.40	1.45
3	A	2002	NAD	C2D-C1D	-2.01	1.50	1.53
2	B	2001	FAD	C2-N1	2.01	1.42	1.38
3	A	2002	NAD	C6N-N1N	2.01	1.41	1.35
3	A	2002	NAD	C2A-N3A	2.12	1.35	1.32
2	A	2001	FAD	C2A-N3A	2.25	1.36	1.32
2	A	2001	FAD	C2-N1	2.26	1.42	1.38
2	A	2001	FAD	C4-N3	2.30	1.37	1.33
3	A	2002	NAD	C6A-N6A	2.35	1.43	1.34
3	B	2002	NAD	C2A-N3A	2.36	1.36	1.32
2	B	2001	FAD	C6A-N6A	2.39	1.43	1.34
2	B	2001	FAD	C5X-N5	2.41	1.39	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2001	FAD	C6A-N6A	2.43	1.44	1.34
2	B	2001	FAD	C2A-N3A	2.51	1.36	1.32
3	B	2002	NAD	C6A-N6A	2.51	1.44	1.34
2	B	2001	FAD	C4-N3	2.54	1.37	1.33
2	A	2001	FAD	C9A-N10	3.20	1.43	1.38
2	B	2001	FAD	C9A-N10	3.28	1.43	1.38
2	A	2001	FAD	C4-C4X	3.38	1.48	1.41
2	A	2001	FAD	C5X-N5	3.46	1.40	1.35
2	B	2001	FAD	C4X-N5	3.71	1.39	1.33
2	A	2001	FAD	C4X-N5	3.91	1.39	1.33
2	B	2001	FAD	C4-C4X	4.02	1.49	1.41
2	B	2001	FAD	O4-C4	4.31	1.35	1.24
2	A	2001	FAD	O4-C4	4.33	1.35	1.24
3	A	2002	NAD	C7N-N7N	4.68	1.42	1.33
3	A	2002	NAD	C2N-N1N	4.91	1.42	1.35
3	B	2002	NAD	C7N-N7N	5.14	1.43	1.33
3	B	2002	NAD	C2N-N1N	5.31	1.42	1.35

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2002	NAD	N3A-C2A-N1A	-12.21	119.28	128.87
3	B	2002	NAD	N3A-C2A-N1A	-12.01	119.43	128.87
2	B	2001	FAD	N3A-C2A-N1A	-11.09	120.16	128.87
2	A	2001	FAD	N3A-C2A-N1A	-10.43	120.68	128.87
2	A	2001	FAD	C4-C4X-C10	-6.31	115.90	119.94
3	A	2002	NAD	C4B-O4B-C1B	-4.02	105.38	109.64
2	B	2001	FAD	C4-C4X-C10	-3.92	117.43	119.94
2	A	2001	FAD	N3-C2-N1	-3.86	121.19	127.69
3	A	2002	NAD	C1B-N9A-C4A	-3.86	122.50	126.81
2	B	2001	FAD	C4X-C4-N3	-3.77	118.59	123.52
3	B	2002	NAD	C4B-O4B-C1B	-3.40	106.04	109.64
2	B	2001	FAD	N3-C2-N1	-3.17	122.36	127.69
3	A	2002	NAD	C4D-O4D-C1D	-2.79	106.69	109.64
2	B	2001	FAD	O2P-P-O3P	-2.65	93.89	105.27
2	A	2001	FAD	C1B-N9A-C4A	-2.37	124.16	126.81
2	B	2001	FAD	O2P-P-O5'	-2.34	97.09	108.24
2	A	2001	FAD	C4B-O4B-C1B	-2.33	107.17	109.64
3	B	2002	NAD	C4D-O4D-C1D	-2.28	107.23	109.64
2	B	2001	FAD	O2A-PA-O5B	-2.28	97.39	108.24
3	A	2002	NAD	O5B-PA-O1A	2.01	117.43	109.21
2	B	2001	FAD	C4-C4X-N5	2.35	121.55	118.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2001	FAD	C2A-N1A-C6A	2.43	123.10	118.77
3	A	2002	NAD	C3N-C7N-N7N	2.51	120.66	117.82
2	A	2001	FAD	C5X-C9A-N10	2.66	119.57	117.58
2	B	2001	FAD	C5X-C9A-N10	3.35	120.09	117.58
2	A	2001	FAD	C4-C4X-N5	3.69	123.19	118.70
2	A	2001	FAD	C4-N3-C2	5.41	119.67	115.16
2	B	2001	FAD	C4-N3-C2	7.64	121.53	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001	FAD	2	0
3	A	2002	NAD	1	0
6	A	2006	PGE	1	0
2	B	2001	FAD	2	0
3	B	2002	NAD	2	0
6	B	2005	PGE	1	0
6	B	2008	PGE	1	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	1207/1235 (97%)	-0.18	27 (2%) 65 70	10, 18, 39, 67	0
1	B	1200/1235 (97%)	-0.06	49 (4%) 41 44	9, 18, 43, 63	0
All	All	2407/2470 (97%)	-0.12	76 (3%) 51 55	9, 18, 41, 67	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1231	ALA	6.8
1	A	129	ASN	5.9
1	B	485	TYR	5.8
1	A	115	ALA	5.3
1	B	506	PRO	4.4
1	B	462	GLY	4.1
1	A	111	ILE	4.0
1	A	1231	ALA	4.0
1	B	1230	MET	3.9
1	A	119	ALA	3.7
1	B	1227	ALA	3.7
1	A	14	ALA	3.6
1	A	121	ILE	3.5
1	B	463	ARG	3.5
1	A	114	THR	3.5
1	A	1230	MET	3.4
1	A	506	PRO	3.4
1	B	915	ALA	3.4
1	B	14	ALA	3.2
1	A	130	TRP	3.2
1	B	482	LEU	3.1
1	B	493	ASN	3.1
1	B	484	ALA	3.1
1	B	514	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	125	ILE	3.0
1	A	122	ARG	3.0
1	B	508	VAL	3.0
1	A	112	PRO	2.9
1	B	481	THR	2.9
1	A	127	ASP	2.9
1	B	115	ALA	2.8
1	B	415	VAL	2.8
1	B	499	PHE	2.8
1	A	490	LEU	2.7
1	B	912	ILE	2.7
1	A	126	ALA	2.7
1	B	511	ASP	2.7
1	A	1227	ALA	2.6
1	A	192	GLY	2.6
1	B	1229	LEU	2.6
1	B	1140	ASN	2.6
1	A	84	VAL	2.6
1	A	118	ASP	2.6
1	B	128	GLY	2.5
1	B	490	LEU	2.5
1	B	510	ILE	2.5
1	B	491	LEU	2.5
1	A	123	ASP	2.4
1	A	128	GLY	2.4
1	B	75	LEU	2.4
1	B	918	THR	2.4
1	A	511	ASP	2.3
1	B	917	GLU	2.3
1	B	916	SER	2.3
1	B	414	ASP	2.3
1	B	138	ARG	2.2
1	B	84	VAL	2.2
1	B	914	LEU	2.2
1	B	524	MET	2.2
1	B	192	GLY	2.2
1	B	77	GLY	2.2
1	A	479	HIS	2.1
1	B	74	ALA	2.1
1	B	432	TYR	2.1
1	B	15	PRO	2.1
1	B	67	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	110	ARG	2.1
1	B	525	PRO	2.1
1	B	127	ASP	2.1
1	B	66	SER	2.1
1	B	70	LYS	2.0
1	B	1200	ARG	2.0
1	B	68	ALA	2.0
1	A	120	LEU	2.0
1	B	466	LEU	2.0
1	B	494	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 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
7	PG4	B	2010	13/13	0.72	0.20	10.44	33,42,52,54	0
6	PGE	A	2007	7/10	0.90	0.15	6.45	20,33,40,42	0
6	PGE	B	2009	7/10	0.92	0.14	3.71	15,31,39,44	0
6	PGE	A	2006	7/10	0.85	0.16	3.57	31,44,46,49	0
6	PGE	A	2005	10/10	0.89	0.11	3.51	23,36,47,51	0
6	PGE	B	2005	10/10	0.91	0.11	2.06	31,41,45,50	0
6	PGE	B	2006	7/10	0.89	0.11	1.39	23,34,41,45	0
6	PGE	B	2007	10/10	0.92	0.09	0.60	24,34,41,43	0
6	PGE	B	2008	10/10	0.84	0.16	0.50	30,43,45,46	0
4	TFB	B	2003	8/8	0.89	0.14	0.36	24,25,28,34	8
3	NAD	B	2002	44/44	0.97	0.09	0.01	9,13,17,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	FAD	A	2001	53/53	0.97	0.08	-0.15	9,13,17,19	0
2	FAD	B	2001	53/53	0.94	0.09	-0.19	12,19,23,25	0
3	NAD	A	2002	44/44	0.97	0.09	-0.27	13,16,19,26	0
4	TFB	A	2003	8/8	0.97	0.06	-2.14	17,18,20,20	0
5	MG	B	2004	1/1	0.97	0.04	-	23,23,23,23	0
5	MG	A	2004	1/1	0.98	0.04	-	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.