



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:27 AM GMT

PDB ID : 3LVM
Title : Crystal Structure of E.coli IscS
Authors : Shi, R.; Proteau, A.; Matte, A.; Cygler, M.; Montreal-Kingston Bacterial Structural Genomics Initiative (BSGI)
Deposited on : 2010-02-22
Resolution : 2.05 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

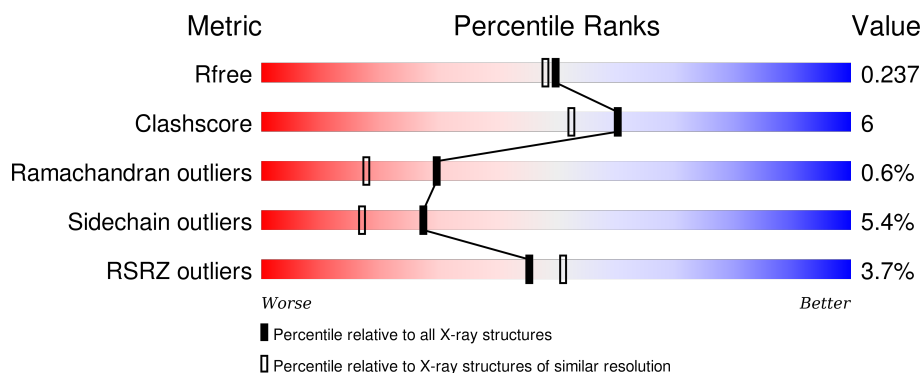
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (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 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	423	
1	B	423	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6495 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 Cysteine desulfurase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	0	0
			3070	1925	548	580	17			
1	B	394	Total	C	N	O	S	0	2	0
			3085	1940	551	577	17			

There are 38 discrepancies between the modelled and reference sequences:

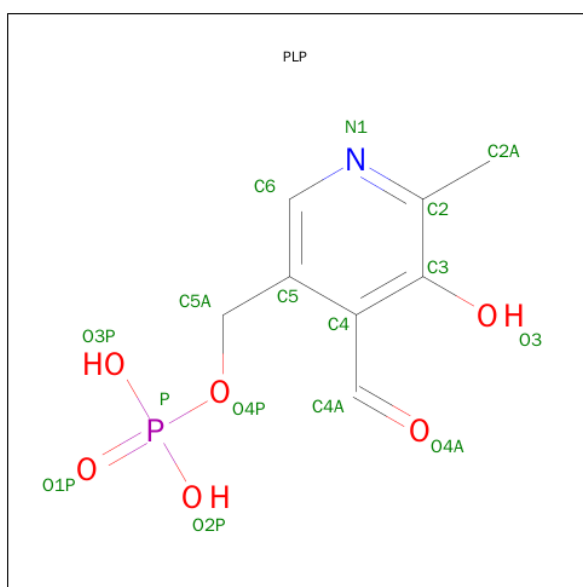
Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	EXPRESSION TAG	UNP P0A6B9
A	-17	GLY	-	EXPRESSION TAG	UNP P0A6B9
A	-16	SER	-	EXPRESSION TAG	UNP P0A6B9
A	-15	SER	-	EXPRESSION TAG	UNP P0A6B9
A	-14	HIS	-	EXPRESSION TAG	UNP P0A6B9
A	-13	HIS	-	EXPRESSION TAG	UNP P0A6B9
A	-12	HIS	-	EXPRESSION TAG	UNP P0A6B9
A	-11	HIS	-	EXPRESSION TAG	UNP P0A6B9
A	-10	HIS	-	EXPRESSION TAG	UNP P0A6B9
A	-9	HIS	-	EXPRESSION TAG	UNP P0A6B9
A	-8	GLY	-	EXPRESSION TAG	UNP P0A6B9
A	-7	SER	-	EXPRESSION TAG	UNP P0A6B9
A	-6	MET	-	EXPRESSION TAG	UNP P0A6B9
A	-5	TYR	-	EXPRESSION TAG	UNP P0A6B9
A	-4	GLY	-	EXPRESSION TAG	UNP P0A6B9
A	-3	VAL	-	EXPRESSION TAG	UNP P0A6B9
A	-2	TYR	-	EXPRESSION TAG	UNP P0A6B9
A	-1	ARG	-	EXPRESSION TAG	UNP P0A6B9
A	0	ALA	-	EXPRESSION TAG	UNP P0A6B9
B	-18	MET	-	EXPRESSION TAG	UNP P0A6B9
B	-17	GLY	-	EXPRESSION TAG	UNP P0A6B9
B	-16	SER	-	EXPRESSION TAG	UNP P0A6B9
B	-15	SER	-	EXPRESSION TAG	UNP P0A6B9
B	-14	HIS	-	EXPRESSION TAG	UNP P0A6B9
B	-13	HIS	-	EXPRESSION TAG	UNP P0A6B9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	HIS	-	EXPRESSION TAG	UNP P0A6B9
B	-11	HIS	-	EXPRESSION TAG	UNP P0A6B9
B	-10	HIS	-	EXPRESSION TAG	UNP P0A6B9
B	-9	HIS	-	EXPRESSION TAG	UNP P0A6B9
B	-8	GLY	-	EXPRESSION TAG	UNP P0A6B9
B	-7	SER	-	EXPRESSION TAG	UNP P0A6B9
B	-6	MET	-	EXPRESSION TAG	UNP P0A6B9
B	-5	TYR	-	EXPRESSION TAG	UNP P0A6B9
B	-4	GLY	-	EXPRESSION TAG	UNP P0A6B9
B	-3	VAL	-	EXPRESSION TAG	UNP P0A6B9
B	-2	TYR	-	EXPRESSION TAG	UNP P0A6B9
B	-1	ARG	-	EXPRESSION TAG	UNP P0A6B9
B	0	ALA	-	EXPRESSION TAG	UNP P0A6B9

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	152	Total	O	0	0
			152	152		

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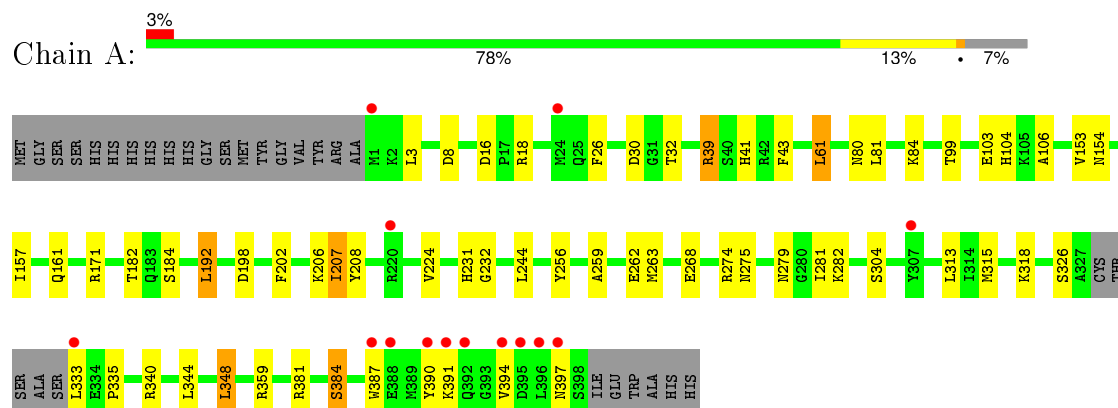
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	158	Total	O	0	0
			158	158		

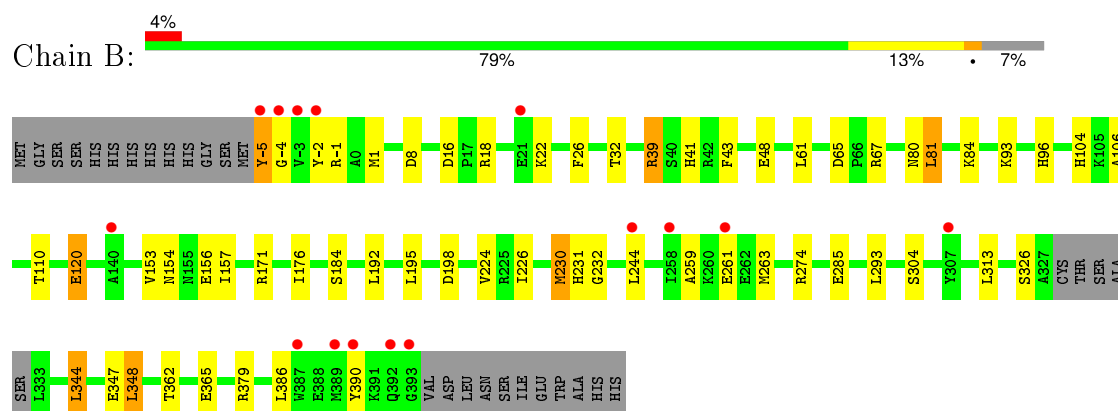
3 Residue-property plots [i](#)

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: Cysteine desulfurase



• Molecule 1: Cysteine desulfurase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.78 Å 99.20 Å 118.07 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.05 49.60 – 2.05	Depositor EDS
% Data completeness (in resolution range)	93.5 (50.00-2.05) 93.4 (49.60-2.05)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.05 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.198 , 0.238 0.199 , 0.237	Depositor DCC
R_{free} test set	2681 reflections (5.41%)	DCC
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.709	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	1 of 52319 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6495	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.47 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.0313e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: PLP

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	0.60	0/3121	0.70	2/4207 (0.0%)
1	B	0.60	0/3141	0.68	1/4235 (0.0%)
All	All	0.60	0/6262	0.69	3/8442 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	344	LEU	CA-CB-CG	6.08	129.28	115.30
1	A	192	LEU	CA-CB-CG	5.11	127.04	115.30
1	A	344	LEU	CA-CB-CG	5.03	126.86	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	3070	0	3071	39	0
1	B	3085	0	3083	36	0
2	A	15	0	7	0	0
2	B	15	0	7	0	0
3	A	152	0	0	2	0
3	B	158	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6495	0	6168	69	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 6.

All (69) 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:39:ARG:HG3	1:A:39:ARG:HH11	0.93	1.06
1:A:39:ARG:CG	1:A:39:ARG:HH11	1.74	1.01
1:A:39:ARG:NH1	1:A:39:ARG:HG3	1.73	0.95
1:A:80:ASN:HD21	1:B:231:HIS:H	1.18	0.89
1:A:231:HIS:H	1:B:80:ASN:HD21	1.18	0.88
1:A:259:ALA:O	1:A:263:MET:HB2	1.79	0.82
1:A:104:HIS:HD2	1:A:106:ALA:H	1.30	0.80
1:B:104:HIS:HD2	1:B:106:ALA:H	1.31	0.78
1:B:259:ALA:O	1:B:263:MET:HB2	1.85	0.76
1:B:198:ASP:HB3	1:B:224[A]:VAL:HG21	1.68	0.76
1:A:304:SER:HB3	1:A:348:LEU:HD22	1.68	0.75
1:A:39:ARG:NH1	1:A:39:ARG:CG	2.40	0.71
1:B:39:ARG:NH1	1:B:48:GLU:OE2	2.24	0.71
1:B:154:ASN:HD22	1:B:157:ILE:H	1.39	0.71
1:A:274:ARG:HH21	1:A:275:ASN:HD21	1.39	0.70
1:B:104:HIS:CD2	1:B:106:ALA:H	2.09	0.69
1:B:347:GLU:HG2	3:B:462:HOH:O	1.92	0.67
1:A:198:ASP:HB3	1:A:224:VAL:HG21	1.75	0.67
1:B:80:ASN:HD22	1:B:84:LYS:HZ2	1.43	0.66
1:B:274:ARG:NH2	1:B:293:LEU:O	2.30	0.65
1:B:-5:TYR:HD1	1:B:-4:GLY:N	1.95	0.64
1:B:16:ASP:OD1	1:B:18:ARG:HD3	1.97	0.64
1:A:326:SER:HB3	1:A:333:LEU:HD11	1.81	0.63
1:A:104:HIS:CD2	1:A:106:ALA:H	2.16	0.62
1:A:154:ASN:HD22	1:A:157:ILE:H	1.47	0.62
1:A:207:ILE:HD11	1:A:256:TYR:CE1	2.36	0.61
1:B:304:SER:HB3	1:B:348:LEU:HD22	1.84	0.60
1:B:362:THR:OG1	1:B:365:GLU:HG3	2.02	0.59
1:A:335:PRO:O	1:A:340:ARG:NH2	2.36	0.58
1:A:274:ARG:HH21	1:A:275:ASN:ND2	2.02	0.56
1:A:80:ASN:HD22	1:A:84:LYS:HZ2	1.52	0.56
1:B:-2:TYR:HA	1:B:1:MET:HE1	1.89	0.55
1:B:81:LEU:HD22	1:B:226:ILE:HD12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:HIS:HD2	1:B:120:GLU:HG3	1.74	0.53
1:A:231:HIS:H	1:B:80:ASN:ND2	1.97	0.52
1:A:8:ASP:OD1	1:B:41:HIS:HE1	1.93	0.52
1:A:16:ASP:OD1	1:A:18:ARG:HD3	2.10	0.51
1:B:80:ASN:HD22	1:B:84:LYS:NZ	2.09	0.51
1:A:61:LEU:HD12	1:A:256:TYR:CD2	2.47	0.50
1:A:397:ASN:HB2	3:A:496:HOH:O	2.11	0.50
1:A:80:ASN:ND2	1:B:231:HIS:H	1.97	0.50
1:A:279:ASN:HA	1:A:282:LYS:HD3	1.92	0.50
1:A:381:ARG:HB3	1:A:387:TRP:CG	2.48	0.49
1:A:41:HIS:HE1	1:B:8:ASP:OD1	1.96	0.48
1:B:41:HIS:CD2	1:B:43:PHE:HB2	2.50	0.47
1:A:80:ASN:HD22	1:A:84:LYS:NZ	2.12	0.46
1:B:154:ASN:ND2	1:B:156:GLU:H	2.13	0.46
1:B:176:ILE:HD12	1:B:224[B]:VAL:HG23	1.97	0.46
1:B:386:LEU:O	1:B:390:TYR:N	2.49	0.45
1:A:41:HIS:CD2	1:A:43:PHE:HB2	2.51	0.45
1:B:93:LYS:HB2	1:B:93:LYS:HE3	1.62	0.45
1:A:207:ILE:HD11	1:A:256:TYR:HE1	1.78	0.45
1:B:26:PHE:CD2	1:B:32:THR:HB	2.52	0.44
1:B:65:ASP:OD1	1:B:67:ARG:HB2	2.16	0.44
1:B:230:MET:CE	3:B:414:HOH:O	2.65	0.44
1:A:153:VAL:HG23	1:A:184:SER:HB3	1.99	0.43
1:B:-1:ARG:HE	1:B:-1:ARG:HB3	1.70	0.43
1:A:18:ARG:NH2	1:A:262:GLU:OE2	2.48	0.43
1:B:153:VAL:HG23	1:B:184:SER:HB3	2.00	0.43
1:B:84:LYS:HE3	1:B:110:THR:HG23	2.01	0.42
1:A:182:THR:HG22	1:A:206:LYS:HG3	2.01	0.42
1:A:99:THR:OG1	1:A:103:GLU:HG3	2.19	0.42
1:A:381:ARG:HB3	1:A:387:TRP:CD1	2.55	0.42
1:B:-5:TYR:CD1	1:B:-4:GLY:N	2.83	0.41
1:A:154:ASN:ND2	1:A:157:ILE:H	2.15	0.41
1:A:161:GLN:HG2	3:A:408:HOH:O	2.21	0.41
1:A:208:TYR:CZ	1:A:359:ARG:HG3	2.56	0.40
1:A:26:PHE:CD2	1:A:32:THR:HB	2.56	0.40
1:B:285:GLU:HB2	3:B:530:HOH:O	2.21	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	389/423 (92%)	371 (95%)	14 (4%)	4 (1%)	19	8
1	B	392/423 (93%)	381 (97%)	10 (3%)	1 (0%)	46	36
All	All	781/846 (92%)	752 (96%)	24 (3%)	5 (1%)	30	18

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	390	TYR
1	A	391	LYS
1	A	232	GLY
1	A	384	SER
1	B	232	GLY

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	326/352 (93%)	308 (94%)	18 (6%)	27	16
1	B	324/352 (92%)	307 (95%)	17 (5%)	29	18
All	All	650/704 (92%)	615 (95%)	35 (5%)	27	17

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU

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Mol	Chain	Res	Type
1	A	30	ASP
1	A	39	ARG
1	A	61	LEU
1	A	81	LEU
1	A	171	ARG
1	A	192	LEU
1	A	202	PHE
1	A	207	ILE
1	A	244	LEU
1	A	268	GLU
1	A	281	ILE
1	A	313	LEU
1	A	315	MET
1	A	318	LYS
1	A	348	LEU
1	A	384	SER
1	A	394	VAL
1	B	-5	TYR
1	B	22	LYS
1	B	39	ARG
1	B	61	LEU
1	B	81	LEU
1	B	120	GLU
1	B	171	ARG
1	B	192	LEU
1	B	195	LEU
1	B	230	MET
1	B	244	LEU
1	B	261	GLU
1	B	313	LEU
1	B	326	SER
1	B	344	LEU
1	B	348	LEU
1	B	379	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	HIS
1	A	57	GLN
1	A	80	ASN
1	A	104	HIS

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Mol	Chain	Res	Type
1	A	127	GLN
1	A	154	ASN
1	A	275	ASN
1	B	41	HIS
1	B	57	GLN
1	B	80	ASN
1	B	104	HIS
1	B	127	GLN
1	B	154	ASN
1	B	275	ASN

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 ⓘ

2 ligands are 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	PLP	A	764	1	15,15,16	1.27	2 (13%)	21,22,23	1.15	3 (14%)
2	PLP	B	764	1	15,15,16	1.19	2 (13%)	21,22,23	1.06	1 (4%)

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	PLP	A	764	1	-	0/6/6/8	0/1/1/1
2	PLP	B	764	1	-	0/6/6/8	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	764	PLP	C6-N1	2.30	1.39	1.34
2	A	764	PLP	C6-N1	2.56	1.39	1.34
2	B	764	PLP	C2-N1	2.76	1.39	1.34
2	A	764	PLP	C2-N1	2.86	1.40	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	764	PLP	C5-C6-N1	-2.21	120.03	123.86
2	A	764	PLP	C4A-C4-C5	2.27	123.25	120.88
2	B	764	PLP	C6-C5-C4	2.41	120.19	118.15
2	A	764	PLP	C6-C5-C4	2.85	120.56	118.15

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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	393/423 (92%)	0.29	14 (3%) 46 53	28, 37, 50, 64	0
1	B	394/423 (93%)	0.30	15 (3%) 44 50	29, 37, 49, 64	0
All	All	787/846 (93%)	0.30	29 (3%) 45 51	28, 37, 49, 64	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-5	TYR	7.8
1	A	394	VAL	5.7
1	B	-2	TYR	4.8
1	B	392	GLN	4.7
1	A	1	MET	4.1
1	A	307	TYR	3.9
1	A	387	TRP	3.6
1	A	390	TYR	3.5
1	A	397	ASN	3.5
1	B	387	TRP	3.3
1	A	391	LYS	3.2
1	B	-4	GLY	3.2
1	B	390	TYR	3.0
1	A	333	LEU	2.9
1	B	-3	VAL	2.9
1	A	392	GLN	2.7
1	B	389	MET	2.7
1	A	388	GLU	2.6
1	B	261	GLU	2.6
1	B	307	TYR	2.5
1	B	258	ILE	2.3
1	A	220	ARG	2.2
1	B	140	ALA	2.2
1	A	24	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	395	ASP	2.1
1	B	21	GLU	2.1
1	A	396	LEU	2.0
1	B	393	GLY	2.0
1	B	244	LEU	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
2	PLP	B	764	15/16	0.92	0.14	-0.00	43,46,51,52	0
2	PLP	A	764	15/16	0.93	0.14	-0.20	45,47,49,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.