



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 18, 2018 – 02:48 AM EST

PDB ID : 5O5C  
Title : The crystal structure of DfoJ, the desferrioxamine biosynthetic pathway lysine decarboxylase from the fire blight disease pathogen *Erwinia amylovora*  
Authors : Salomone-Stagni, M.; Bartho, J.D.; Polsinelli, I.; Bellini, D.; Walsh, M.A.; Demitri, N.; Benini, S.  
Deposited on : 2017-06-01  
Resolution : 2.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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-20030736
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-20030736

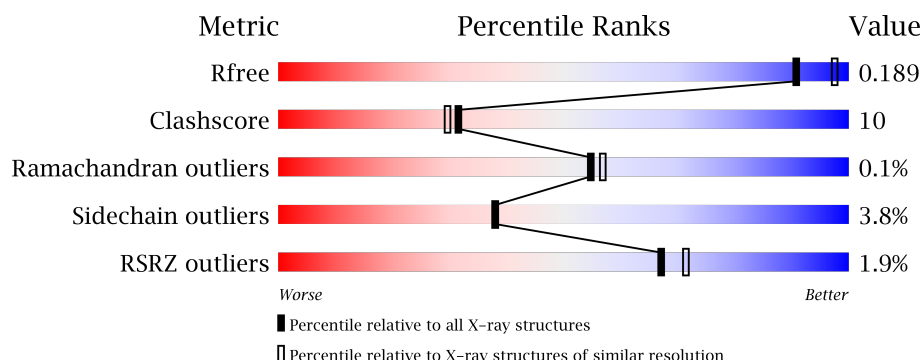
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

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  $\geq 3$ , 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	519	<div> <div>3%</div> <div>68%18%11%</div> </div>
1	B	519	<div> <div>2%</div> <div>71%17%10%</div> </div>
1	C	519	<div> <div>2%</div> <div>68%18%11%</div> </div>
1	D	519	<div> <div>2%</div> <div>72%17%10%</div> </div>
1	E	519	<div> <div>0%</div> <div>74%13%12%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	519	

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
2	PLP	B	601	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 22003 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 Putative decarboxylase involved in desferrioxamine biosynthesis.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3621	2302	641	661	17			
1	B	468	Total	C	N	O	S	0	0	0
			3645	2316	645	667	17			
1	C	462	Total	C	N	O	S	0	0	0
			3599	2289	635	658	17			
1	D	468	Total	C	N	O	S	0	0	0
			3646	2317	645	667	17			
1	E	457	Total	C	N	O	S	0	0	0
			3565	2266	631	651	17			
1	F	467	Total	C	N	O	S	0	0	0
			3637	2312	644	664	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP D4I245
A	0	ALA	-	expression tag	UNP D4I245
B	-1	GLY	-	expression tag	UNP D4I245
B	0	ALA	-	expression tag	UNP D4I245
C	-1	GLY	-	expression tag	UNP D4I245
C	0	ALA	-	expression tag	UNP D4I245
D	-1	GLY	-	expression tag	UNP D4I245
D	0	ALA	-	expression tag	UNP D4I245
E	-1	GLY	-	expression tag	UNP D4I245
E	0	ALA	-	expression tag	UNP D4I245
F	-1	GLY	-	expression tag	UNP D4I245
F	0	ALA	-	expression tag	UNP D4I245

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C<sub>8</sub>H<sub>10</sub>NO<sub>6</sub>P).



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		
2	C	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	D	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	E	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	F	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

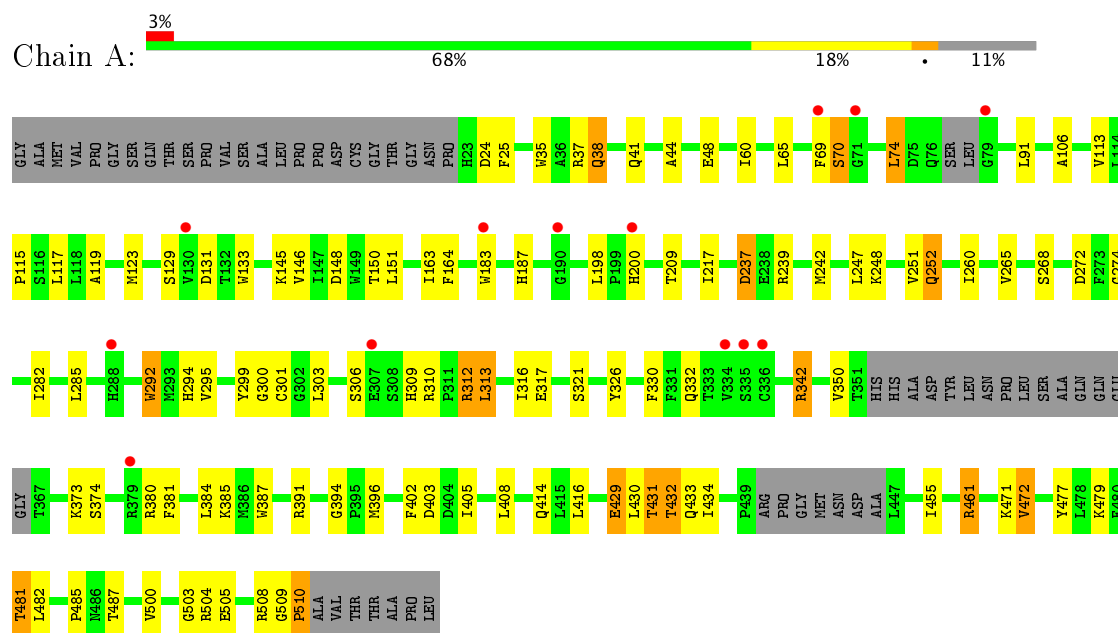
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total	O	0	0
			30	30		
3	B	31	Total	O	0	0
			31	31		
3	C	43	Total	O	0	0
			43	43		
3	D	42	Total	O	0	0
			42	42		
3	E	23	Total	O	0	0
			23	23		
3	F	31	Total	O	0	0
			31	31		

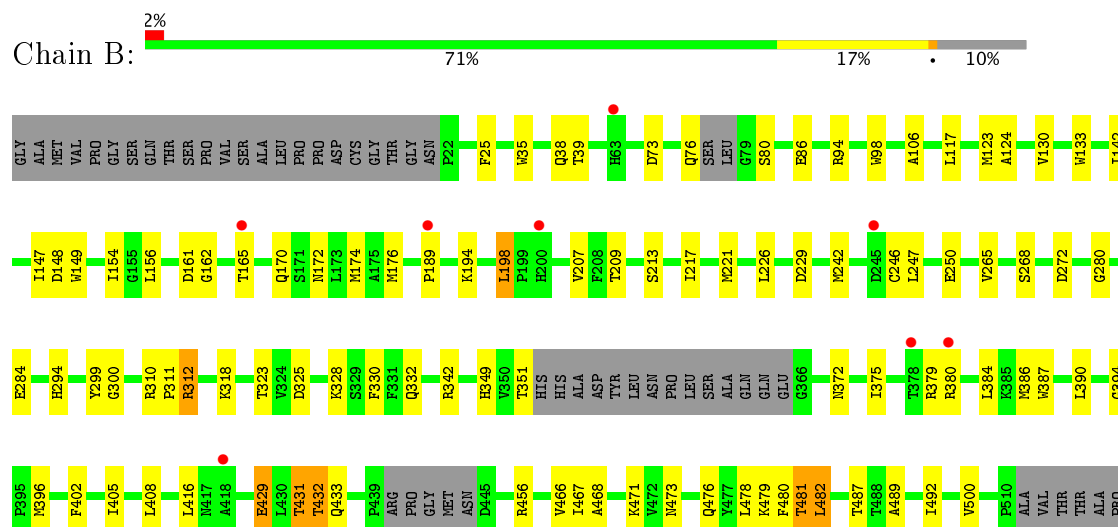
### 3 Residue-property plots [i](#)

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

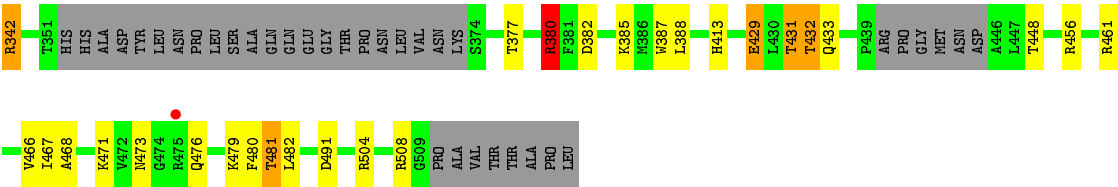
- Molecule 1: Putative decarboxylase involved in desferrioxamine biosynthesis



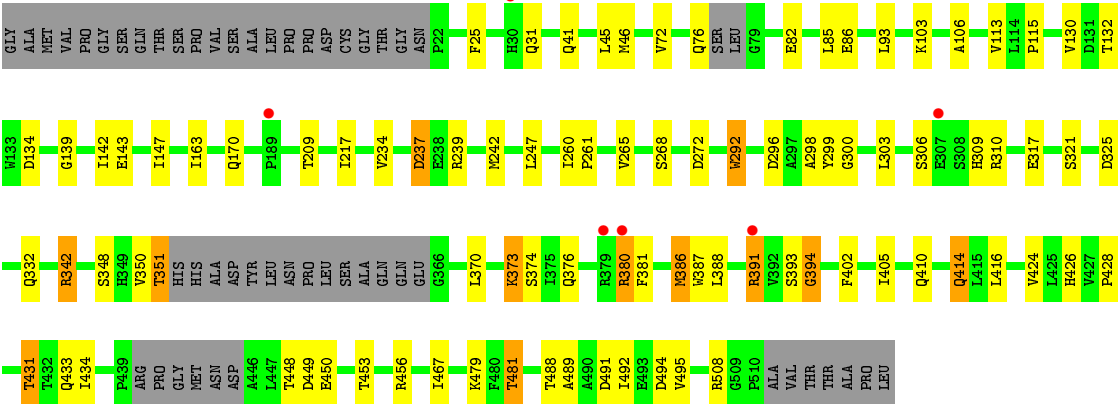
- Molecule 1: Putative decarboxylase involved in desferrioxamine biosynthesis



GLY	ALA	MET	VAL	PRO	GLY	SER	GLN	THR	SER	PRO	VAL	SER	ALA	LEU	PRO	ASP	CYS	GLY	THR	ASN	PRO														
K90	L91	K103	A106	A119	E120	M123	D131	D181	S182	W183	P189	G190	H200	V207	D237	E238	M242	C246	E250	P261	V262	A263	S268	D272	H288	M291	Y299	G300	R310	P311	R312	E317	K318	V324	S335



● Molecule 1: Putative decarboxylase involved in desferrioxamine biosynthesis





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	269.79 Å   269.79 Å   56.19 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	233.65 – 2.10 47.41 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (233.65-2.10) 99.9 (47.41-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	110.40 (at 2.10 Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.165 , 0.187 0.169 , 0.189	Depositor DCC
$R_{free}$ test set	13243 reflections (5.23%)	DCC
Wilson B-factor (Å <sup>2</sup> )	29.4	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 45.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l 0.000 for h,-h-k,-l 0.036 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	22003	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	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 23.63 % 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 4.4894e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: 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	0.79	2/3704 (0.1%)	0.94	11/5031 (0.2%)
1	B	0.82	2/3729 (0.1%)	0.95	8/5065 (0.2%)
1	C	0.79	0/3682	1.00	17/5001 (0.3%)
1	D	0.75	0/3730	0.91	5/5066 (0.1%)
1	E	0.78	1/3646 (0.0%)	0.94	9/4950 (0.2%)
1	F	0.81	1/3721 (0.0%)	0.93	4/5054 (0.1%)
All	All	0.79	6/22212 (0.0%)	0.95	54/30167 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
1	F	0	1
All	All	0	4

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	292	TRP	CB-CG	-9.63	1.32	1.50
1	E	237	ASP	CB-CG	-5.86	1.39	1.51
1	A	429	GLU	CG-CD	-5.33	1.44	1.51
1	B	133	TRP	CE3-CZ3	-5.33	1.29	1.38
1	F	394	GLY	N-CA	5.07	1.53	1.46

The worst 5 of 54 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	380	ARG	NE-CZ-NH2	-15.33	112.64	120.30
1	C	380	ARG	NE-CZ-NH1	11.11	125.86	120.30
1	E	380	ARG	NE-CZ-NH2	-10.93	114.84	120.30
1	E	380	ARG	NE-CZ-NH1	10.70	125.65	120.30
1	A	380	ARG	NE-CZ-NH2	-10.07	115.26	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	394	GLY	Peptide
1	A	509	GLY	Peptide
1	B	394	GLY	Peptide
1	F	394	GLY	Peptide

## 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	3621	0	3592	82	2
1	B	3645	0	3615	83	1
1	C	3599	0	3568	109	0
1	D	3646	0	3617	71	1
1	E	3565	0	3531	41	3
1	F	3637	0	3608	74	3
2	A	15	0	6	0	0
2	B	15	0	7	5	0
2	C	15	0	6	1	0
2	D	15	0	6	4	0
2	E	15	0	7	0	0
2	F	15	0	6	1	0
3	A	30	0	0	0	0
3	B	31	0	0	3	0
3	C	43	0	0	2	0
3	D	42	0	0	1	0
3	E	23	0	0	0	0
3	F	31	0	0	1	0
All	All	22003	0	21569	442	5

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 10.

The worst 5 of 442 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:D:471:LYS:NZ	1:D:476:GLN:OE1	1.74	1.18
1:B:328:LYS:NZ	2:B:601:PLP:C4A	2.07	1.17
1:A:292:TRP:HH2	1:A:321:SER:OG	1.28	1.13
1:A:74:LEU:HD21	1:B:390:LEU:HD13	1.25	1.12
1:D:328:LYS:NZ	2:D:601:PLP:C4A	2.13	1.11

All (5) 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)
1:A:70:SER:OG	1:B:76:GLN:OE1[1_556]	1.66	0.54
1:E:335:SER:OG	1:F:380:ARG:NH1[1_556]	1.73	0.47
1:D:475:ARG:NH2	1:F:508:ARG:O[1_556]	1.80	0.40
1:E:84:ALA:O	1:F:391:ARG:NH1[1_556]	1.86	0.34
1:A:508:ARG:O	1:E:473:ASN:ND2[2_664]	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	456/519 (88%)	439 (96%)	17 (4%)	0	100	100
1	B	460/519 (89%)	436 (95%)	22 (5%)	2 (0%)	38	35
1	C	454/519 (88%)	428 (94%)	25 (6%)	1 (0%)	51	52
1	D	460/519 (89%)	435 (95%)	25 (5%)	0	100	100
1	E	449/519 (86%)	423 (94%)	26 (6%)	0	100	100
1	F	459/519 (88%)	432 (94%)	26 (6%)	1 (0%)	51	52

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	2738/3114 (88%)	2593 (95%)	141 (5%)	4 (0%)	55 57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	24	ASP
1	F	237	ASP
1	B	80	SER
1	B	189	PRO

### 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	390/432 (90%)	369 (95%)	21 (5%)	26 23
1	B	392/432 (91%)	382 (97%)	10 (3%)	51 55
1	C	387/432 (90%)	369 (95%)	18 (5%)	30 28
1	D	392/432 (91%)	380 (97%)	12 (3%)	45 48
1	E	382/432 (88%)	370 (97%)	12 (3%)	45 48
1	F	391/432 (90%)	376 (96%)	15 (4%)	38 38
All	All	2334/2592 (90%)	2246 (96%)	88 (4%)	38 38

5 of 88 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	373	LYS
1	D	25	PHE
1	F	380	ARG
1	C	380	ARG
1	C	430	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	170	GLN

### 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](#)

6 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	601	1	15,15,16	3.44	3 (20%)	20,22,23	1.41	3 (15%)
2	PLP	B	601	-	15,15,16	3.11	3 (20%)	20,22,23	1.61	4 (20%)
2	PLP	C	601	1	15,15,16	3.78	3 (20%)	20,22,23	1.12	1 (5%)
2	PLP	D	601	-	15,15,16	3.06	3 (20%)	20,22,23	1.56	4 (20%)
2	PLP	E	601	1	15,15,16	3.49	3 (20%)	20,22,23	1.28	1 (5%)
2	PLP	F	601	1	15,15,16	3.44	4 (26%)	20,22,23	1.51	5 (25%)

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	601	1	-	0/6/6/8	0/1/1/1
2	PLP	B	601	-	-	0/6/6/8	0/1/1/1
2	PLP	C	601	1	-	0/6/6/8	0/1/1/1
2	PLP	D	601	-	-	0/6/6/8	0/1/1/1
2	PLP	E	601	1	-	0/6/6/8	0/1/1/1
2	PLP	F	601	1	-	0/6/6/8	0/1/1/1

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	601	PLP	C4A-C4	-2.64	1.46	1.51
2	E	601	PLP	C3-C4	3.94	1.48	1.40
2	C	601	PLP	C3-C4	3.97	1.48	1.40
2	B	601	PLP	C3-C4	4.02	1.48	1.40
2	F	601	PLP	C3-C4	4.05	1.48	1.40

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	PLP	O4P-C5A-C5	-2.76	103.76	109.32
2	F	601	PLP	C4A-C4-C5	-2.59	118.24	120.86
2	B	601	PLP	C4-C3-C2	-2.43	116.06	120.06
2	D	601	PLP	C4A-C4-C5	-2.42	118.42	120.86
2	A	601	PLP	C2A-C2-C3	-2.38	118.13	120.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	PLP	5	0
2	C	601	PLP	1	0
2	D	601	PLP	4	0
2	F	601	PLP	1	0

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	464/519 (89%)	-0.12	13 (2%) 53 60	19, 31, 49, 76	0
1	B	468/519 (90%)	-0.18	8 (1%) 70 75	16, 28, 45, 64	0
1	C	462/519 (89%)	-0.17	10 (2%) 62 67	16, 28, 45, 62	0
1	D	468/519 (90%)	-0.15	8 (1%) 70 75	17, 30, 49, 63	0
1	E	457/519 (88%)	-0.22	7 (1%) 74 77	15, 28, 46, 72	0
1	F	467/519 (89%)	-0.27	6 (1%) 77 81	16, 28, 43, 67	0
All	All	2786/3114 (89%)	-0.18	52 (1%) 67 71	15, 29, 47, 76	0

The worst 5 of 52 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	71	GLY	6.8
1	F	391	ARG	5.0
1	E	71	GLY	4.7
1	D	200	HIS	4.2
1	C	480	PHE	4.1

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	PLP	B	601	15/16	0.98	0.22	2.36	35,37,40,41	0
2	PLP	E	601	15/16	0.98	0.20	1.33	35,41,46,49	0
2	PLP	D	601	15/16	0.96	0.21	1.32	39,42,47,48	0
2	PLP	A	601	15/16	0.99	0.20	0.96	35,40,42,42	0
2	PLP	C	601	15/16	0.98	0.20	0.94	29,39,42,44	0
2	PLP	F	601	15/16	0.97	0.18	0.80	36,37,41,42	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.