



wwPDB X-ray Structure Validation Summary Report

Aug 17, 2022 – 05:39 PM EDT

PDB ID : 3X43
Title : Crystal structure of O-ureido-L-serine synthase
Authors : Matoba, Y.; Uda, N.; Oda, K.; Sugiyama, M.
Deposited on : 2015-03-13
Resolution : 2.25 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

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 : 2.29
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) : 2.29

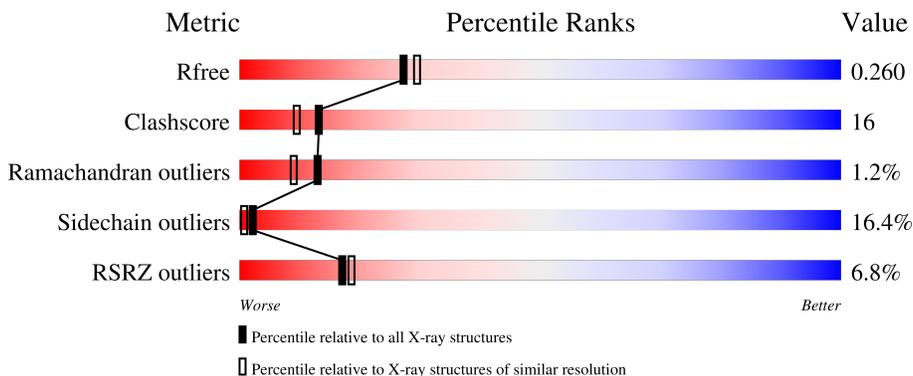
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 $\leq 5\%$. 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	332	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 62%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 27%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div>
1	B	332	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 60%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 27%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div>
1	C	332	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 62%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 26%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 7%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div>
1	D	332	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 64%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 25%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div>
1	E	332	<div style="display: flex; align-items: center;"> <div style="width: 12%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 58%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 30%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div>

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Mol	Chain	Length	Quality of chain
1	F	332	<p>% 64% 23% 8% 5%</p>
1	G	332	<p>7% 64% 26% 6% 5%</p>
1	H	332	<p>9% 60% 28% 7% 5%</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20024 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 O-ureido-L-serine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	316	2381	1502	419	452	8	0	0	0
1	B	319	2404	1516	422	458	8	0	0	0
1	C	317	2389	1508	420	453	8	0	0	0
1	D	316	2381	1502	419	452	8	0	0	0
1	E	317	2389	1508	420	453	8	0	0	0
1	F	316	2381	1502	419	452	8	0	0	0
1	G	317	2389	1508	420	453	8	0	0	0
1	H	316	2381	1502	419	452	8	0	0	0

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	325	LEU	-	expression tag	UNP D2Z027
A	326	GLU	-	expression tag	UNP D2Z027
A	327	HIS	-	expression tag	UNP D2Z027
A	328	HIS	-	expression tag	UNP D2Z027
A	329	HIS	-	expression tag	UNP D2Z027
A	330	HIS	-	expression tag	UNP D2Z027
A	331	HIS	-	expression tag	UNP D2Z027
A	332	HIS	-	expression tag	UNP D2Z027
B	325	LEU	-	expression tag	UNP D2Z027
B	326	GLU	-	expression tag	UNP D2Z027
B	327	HIS	-	expression tag	UNP D2Z027
B	328	HIS	-	expression tag	UNP D2Z027
B	329	HIS	-	expression tag	UNP D2Z027

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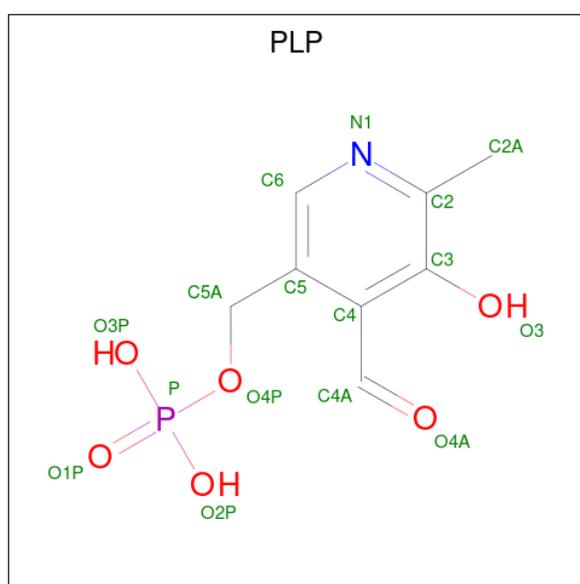
Chain	Residue	Modelled	Actual	Comment	Reference
B	330	HIS	-	expression tag	UNP D2Z027
B	331	HIS	-	expression tag	UNP D2Z027
B	332	HIS	-	expression tag	UNP D2Z027
C	325	LEU	-	expression tag	UNP D2Z027
C	326	GLU	-	expression tag	UNP D2Z027
C	327	HIS	-	expression tag	UNP D2Z027
C	328	HIS	-	expression tag	UNP D2Z027
C	329	HIS	-	expression tag	UNP D2Z027
C	330	HIS	-	expression tag	UNP D2Z027
C	331	HIS	-	expression tag	UNP D2Z027
C	332	HIS	-	expression tag	UNP D2Z027
D	325	LEU	-	expression tag	UNP D2Z027
D	326	GLU	-	expression tag	UNP D2Z027
D	327	HIS	-	expression tag	UNP D2Z027
D	328	HIS	-	expression tag	UNP D2Z027
D	329	HIS	-	expression tag	UNP D2Z027
D	330	HIS	-	expression tag	UNP D2Z027
D	331	HIS	-	expression tag	UNP D2Z027
D	332	HIS	-	expression tag	UNP D2Z027
E	325	LEU	-	expression tag	UNP D2Z027
E	326	GLU	-	expression tag	UNP D2Z027
E	327	HIS	-	expression tag	UNP D2Z027
E	328	HIS	-	expression tag	UNP D2Z027
E	329	HIS	-	expression tag	UNP D2Z027
E	330	HIS	-	expression tag	UNP D2Z027
E	331	HIS	-	expression tag	UNP D2Z027
E	332	HIS	-	expression tag	UNP D2Z027
F	325	LEU	-	expression tag	UNP D2Z027
F	326	GLU	-	expression tag	UNP D2Z027
F	327	HIS	-	expression tag	UNP D2Z027
F	328	HIS	-	expression tag	UNP D2Z027
F	329	HIS	-	expression tag	UNP D2Z027
F	330	HIS	-	expression tag	UNP D2Z027
F	331	HIS	-	expression tag	UNP D2Z027
F	332	HIS	-	expression tag	UNP D2Z027
G	325	LEU	-	expression tag	UNP D2Z027
G	326	GLU	-	expression tag	UNP D2Z027
G	327	HIS	-	expression tag	UNP D2Z027
G	328	HIS	-	expression tag	UNP D2Z027
G	329	HIS	-	expression tag	UNP D2Z027
G	330	HIS	-	expression tag	UNP D2Z027
G	331	HIS	-	expression tag	UNP D2Z027

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Chain	Residue	Modelled	Actual	Comment	Reference
G	332	HIS	-	expression tag	UNP D2Z027
H	325	LEU	-	expression tag	UNP D2Z027
H	326	GLU	-	expression tag	UNP D2Z027
H	327	HIS	-	expression tag	UNP D2Z027
H	328	HIS	-	expression tag	UNP D2Z027
H	329	HIS	-	expression tag	UNP D2Z027
H	330	HIS	-	expression tag	UNP D2Z027
H	331	HIS	-	expression tag	UNP D2Z027
H	332	HIS	-	expression tag	UNP D2Z027

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



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		
2	G	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	H	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	111	Total	O	0	0
			111	111		
3	B	102	Total	O	0	0
			102	102		
3	C	96	Total	O	0	0
			96	96		
3	D	114	Total	O	0	0
			114	114		
3	E	106	Total	O	0	0
			106	106		
3	F	98	Total	O	0	0
			98	98		
3	G	97	Total	O	0	0
			97	97		
3	H	85	Total	O	0	0
			85	85		

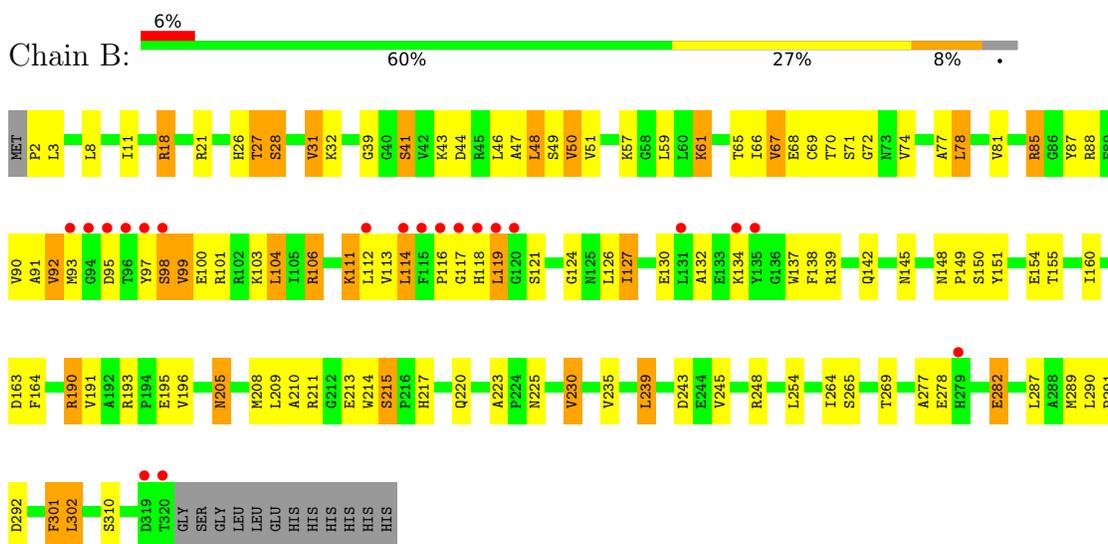
3 Residue-property plots i

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: O-ureido-L-serine synthase

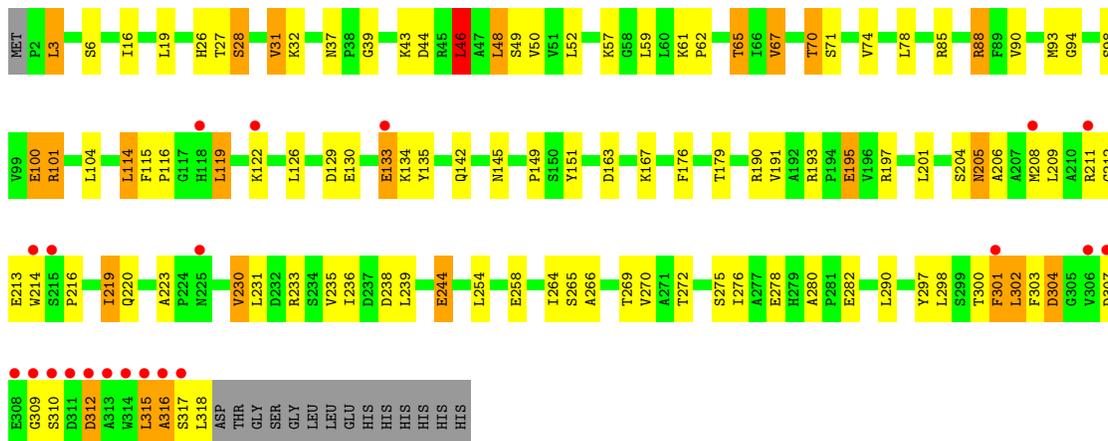


- Molecule 1: O-ureido-L-serine synthase

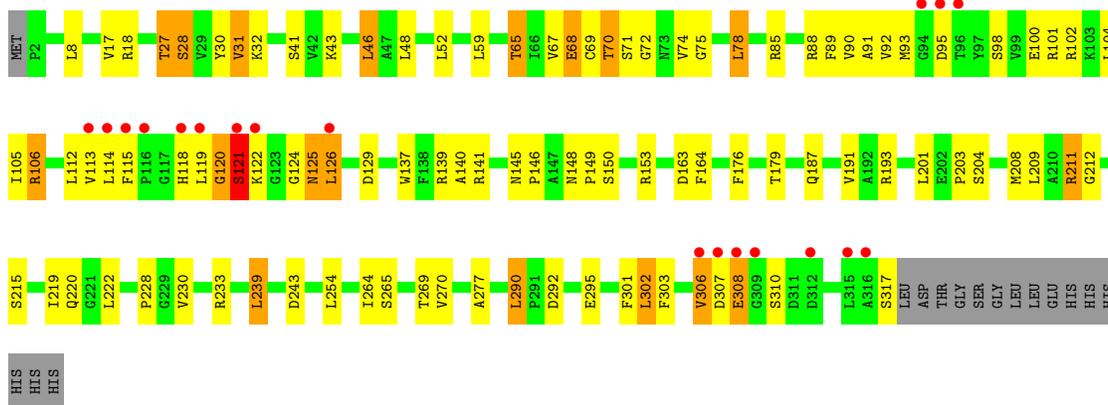


- Molecule 1: O-ureido-L-serine synthase

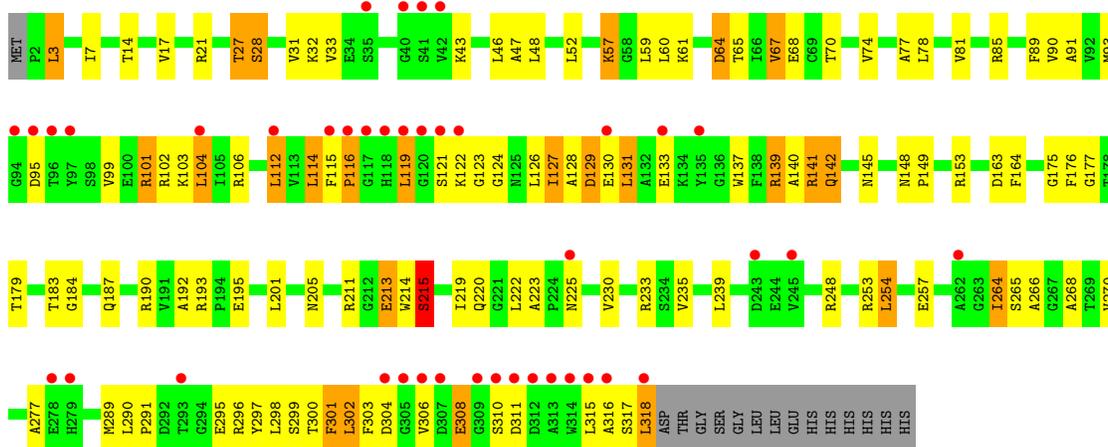




- Molecule 1: O-ureido-L-serine synthase



- Molecule 1: O-ureido-L-serine synthase



- Molecule 1: O-ureido-L-serine synthase

4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.48Å 154.13Å 118.52Å 90.00° 90.48° 90.00°	Depositor
Resolution (Å)	29.83 – 2.25 36.64 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.83-2.25) 99.9 (36.64-2.25)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 2.24Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.209 , 0.264 0.207 , 0.260	Depositor DCC
R_{free} test set	6181 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	34.1	Xtrriage
Anisotropy	0.260	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.022 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20024	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.39 % 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 1.3181e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

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	0.34	0/2427	0.60	0/3295
1	B	0.33	0/2450	0.59	0/3327
1	C	0.34	0/2435	0.61	1/3306 (0.0%)
1	D	0.37	0/2427	0.61	1/3295 (0.0%)
1	E	0.33	0/2435	0.58	2/3306 (0.1%)
1	F	0.37	0/2427	0.62	0/3295
1	G	0.33	0/2435	0.59	0/3306
1	H	0.33	0/2427	0.59	0/3295
All	All	0.34	0/19463	0.60	4/26425 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	46	LEU	CA-CB-CG	5.46	127.85	115.30
1	E	175	GLY	N-CA-C	-5.32	99.81	113.10
1	D	46	LEU	CA-CB-CG	5.26	127.41	115.30
1	E	177	GLY	N-CA-C	-5.01	100.56	113.10

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	2381	0	2374	75	0
1	B	2404	0	2396	95	0
1	C	2389	0	2385	85	0
1	D	2381	0	2374	87	0
1	E	2389	0	2385	97	0
1	F	2381	0	2374	70	0
1	G	2389	0	2385	81	0
1	H	2381	0	2374	81	0
2	A	15	0	6	0	0
2	B	15	0	6	0	0
2	C	15	0	6	0	0
2	D	15	0	6	0	0
2	E	15	0	6	0	0
2	F	15	0	6	0	0
2	G	15	0	6	0	0
2	H	15	0	6	0	0
3	A	111	0	0	2	0
3	B	102	0	0	1	0
3	C	96	0	0	6	0
3	D	114	0	0	2	0
3	E	106	0	0	0	0
3	F	98	0	0	1	0
3	G	97	0	0	1	0
3	H	85	0	0	1	0
All	All	20024	0	19095	626	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 16.

The worst 5 of 626 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:57:LYS:HB3	1:A:59:LEU:HD13	1.30	1.07
1:B:114:LEU:H	1:C:310:SER:HB2	1.21	1.06
1:D:68:GLU:HA	1:D:139:ARG:HH21	1.17	1.04
1:B:67:VAL:HG12	1:B:90:VAL:HB	1.44	0.95
1:F:26:HIS:HB2	1:F:282:GLU:HG2	1.48	0.94

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	314/332 (95%)	294 (94%)	17 (5%)	3 (1%)	15	13
1	B	317/332 (96%)	293 (92%)	19 (6%)	5 (2%)	9	5
1	C	315/332 (95%)	285 (90%)	26 (8%)	4 (1%)	12	8
1	D	314/332 (95%)	292 (93%)	20 (6%)	2 (1%)	25	25
1	E	315/332 (95%)	288 (91%)	23 (7%)	4 (1%)	12	8
1	F	314/332 (95%)	294 (94%)	15 (5%)	5 (2%)	9	5
1	G	315/332 (95%)	288 (91%)	24 (8%)	3 (1%)	15	13
1	H	314/332 (95%)	286 (91%)	24 (8%)	4 (1%)	12	8
All	All	2518/2656 (95%)	2320 (92%)	168 (7%)	30 (1%)	13	9

5 of 30 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	316	ALA
1	D	121	SER
1	E	215	SER
1	G	312	ASP
1	H	117	GLY

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	247/261 (95%)	203 (82%)	44 (18%)	2	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	250/261 (96%)	206 (82%)	44 (18%)	2	0
1	C	248/261 (95%)	210 (85%)	38 (15%)	2	1
1	D	247/261 (95%)	215 (87%)	32 (13%)	4	2
1	E	248/261 (95%)	201 (81%)	47 (19%)	1	0
1	F	247/261 (95%)	205 (83%)	42 (17%)	2	0
1	G	248/261 (95%)	214 (86%)	34 (14%)	3	2
1	H	247/261 (95%)	203 (82%)	44 (18%)	2	0
All	All	1982/2088 (95%)	1657 (84%)	325 (16%)	2	1

5 of 325 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	139	ARG
1	H	48	LEU
1	F	254	LEU
1	G	99	VAL
1	H	112	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	220	GLN
1	G	20	GLN
1	H	142	GLN
1	H	20	GLN
1	H	125	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	F	401	1	15,15,16	2.13	7 (46%)	20,22,23	1.47	5 (25%)
2	PLP	H	401	1	15,15,16	2.07	7 (46%)	20,22,23	1.50	5 (25%)
2	PLP	E	401	1	15,15,16	2.16	6 (40%)	20,22,23	1.46	5 (25%)
2	PLP	B	401	1	15,15,16	2.06	7 (46%)	20,22,23	1.46	4 (20%)
2	PLP	A	401	1	15,15,16	2.09	7 (46%)	20,22,23	1.41	4 (20%)
2	PLP	G	401	1	15,15,16	2.05	6 (40%)	20,22,23	1.48	4 (20%)
2	PLP	D	401	1	15,15,16	2.06	6 (40%)	20,22,23	1.46	5 (25%)
2	PLP	C	401	1	15,15,16	2.13	7 (46%)	20,22,23	1.50	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	F	401	1	-	0/6/6/8	0/1/1/1
2	PLP	H	401	1	-	0/6/6/8	0/1/1/1
2	PLP	E	401	1	-	0/6/6/8	0/1/1/1
2	PLP	B	401	1	-	0/6/6/8	0/1/1/1
2	PLP	A	401	1	-	0/6/6/8	0/1/1/1
2	PLP	G	401	1	-	0/6/6/8	0/1/1/1
2	PLP	D	401	1	-	0/6/6/8	0/1/1/1
2	PLP	C	401	1	-	0/6/6/8	0/1/1/1

The worst 5 of 53 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	PLP	C3-C2	-4.29	1.36	1.40
2	C	401	PLP	C3-C2	-3.83	1.37	1.40
2	G	401	PLP	C2-N1	3.79	1.41	1.33
2	A	401	PLP	C3-C2	-3.78	1.37	1.40
2	B	401	PLP	C3-C2	-3.75	1.37	1.40

The worst 5 of 37 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	401	PLP	O4P-C5A-C5	2.91	114.90	109.35
2	H	401	PLP	O4P-C5A-C5	2.90	114.89	109.35
2	C	401	PLP	O4P-C5A-C5	2.83	114.75	109.35
2	B	401	PLP	C3-C2-N1	-2.80	117.14	120.77
2	F	401	PLP	C6-N1-C2	2.80	124.36	119.17

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 [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 [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, 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	316/332 (95%)	0.08	14 (4%) 34 37	23, 44, 86, 102	0
1	B	319/332 (96%)	0.09	20 (6%) 20 22	26, 46, 102, 114	0
1	C	317/332 (95%)	0.20	21 (6%) 18 20	22, 48, 89, 110	0
1	D	316/332 (95%)	0.01	19 (6%) 21 23	22, 39, 97, 119	0
1	E	317/332 (95%)	0.49	41 (12%) 3 3	23, 52, 106, 126	0
1	F	316/332 (95%)	-0.14	3 (0%) 84 85	23, 41, 80, 93	0
1	G	317/332 (95%)	0.34	23 (7%) 15 15	27, 49, 90, 107	0
1	H	316/332 (95%)	0.26	31 (9%) 7 8	27, 47, 111, 132	0
All	All	2534/2656 (95%)	0.17	172 (6%) 17 18	22, 46, 99, 132	0

The worst 5 of 172 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	119	LEU	10.5
1	B	96	THR	8.5
1	D	119	LEU	8.3
1	H	313	ALA	7.3
1	E	118	HIS	6.6

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, 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	B-factors(Å ²)	Q<0.9
2	PLP	B	401	15/16	0.97	0.14	31,38,41,41	0
2	PLP	C	401	15/16	0.97	0.16	37,42,45,46	0
2	PLP	E	401	15/16	0.97	0.17	43,45,48,49	0
2	PLP	H	401	15/16	0.97	0.11	36,41,43,43	0
2	PLP	G	401	15/16	0.98	0.17	36,39,43,46	0
2	PLP	A	401	15/16	0.98	0.16	30,34,36,38	0
2	PLP	D	401	15/16	0.99	0.13	28,32,34,34	0
2	PLP	F	401	15/16	0.99	0.15	27,30,33,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.