



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 06:55 pm GMT

PDB ID : 2J6X  
Title : THE CRYSTAL STRUCTURE OF LACTATE OXIDASE  
Authors : Leiros, I.; Wang, E.; Rasmussen, T.; Oksanen, E.; Repo, H.; Petersen, S.B.;  
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Deposited on : 2006-10-05  
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 : trunk28620  
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) : trunk28620

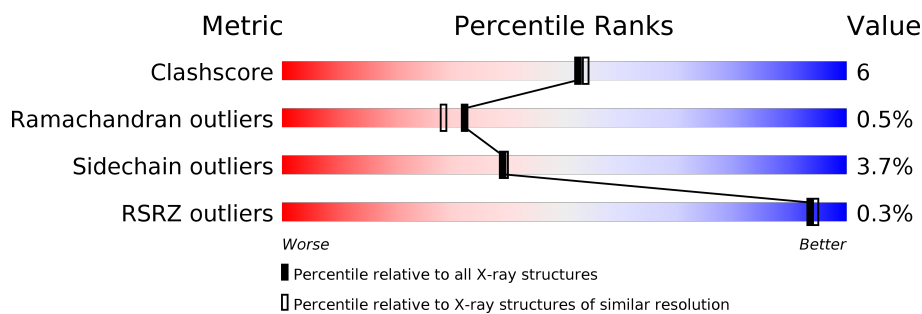
# 1 Overall quality at a glance

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))
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	374	
1	B	374	
1	C	374	
1	D	374	
1	E	374	
1	F	374	
1	G	374	

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Mol	Chain	Length	Quality of chain
1	H	374	<div><div><div>%</div><div><div></div></div><div>83%</div><div>9%</div><div>• 7%</div></div></div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23895 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 LACTATE OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	370	Total	C	N	O	S	0	0	0
			2852	1802	493	550	7			
1	B	367	Total	C	N	O	S	0	0	0
			2827	1787	490	543	7			
1	C	368	Total	C	N	O	S	0	0	0
			2836	1792	491	546	7			
1	D	353	Total	C	N	O	S	0	0	0
			2717	1721	468	522	6			
1	E	367	Total	C	N	O	S	0	0	0
			2827	1787	490	543	7			
1	F	357	Total	C	N	O	S	0	0	0
			2748	1739	474	529	6			
1	G	367	Total	C	N	O	S	0	0	0
			2827	1787	490	543	7			
1	H	348	Total	C	N	O	S	0	0	0
			2682	1697	462	517	6			

There are 16 discrepancies between the modelled and reference sequences:

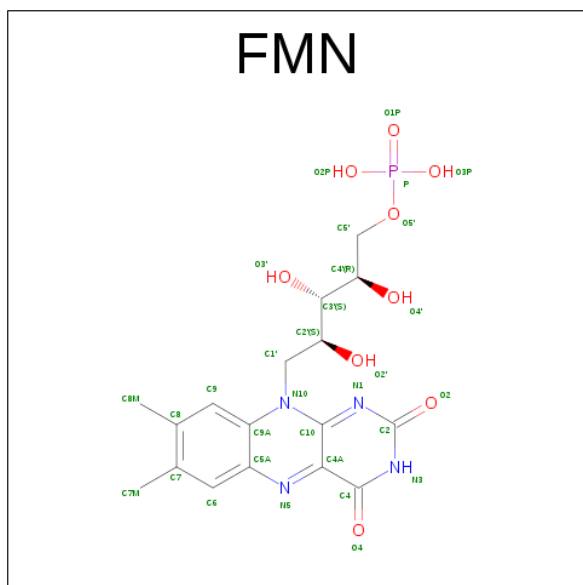
Chain	Residue	Modelled	Actual	Comment	Reference
A	232	ALA	GLY	CONFLICT	UNP Q44467
A	255	ALA	ARG	CONFLICT	UNP Q44467
B	232	ALA	GLY	CONFLICT	UNP Q44467
B	255	ALA	ARG	CONFLICT	UNP Q44467
C	232	ALA	GLY	CONFLICT	UNP Q44467
C	255	ALA	ARG	CONFLICT	UNP Q44467
D	232	ALA	GLY	CONFLICT	UNP Q44467
D	255	ALA	ARG	CONFLICT	UNP Q44467
E	232	ALA	GLY	CONFLICT	UNP Q44467
E	255	ALA	ARG	CONFLICT	UNP Q44467
F	232	ALA	GLY	CONFLICT	UNP Q44467
F	255	ALA	ARG	CONFLICT	UNP Q44467
G	232	ALA	GLY	CONFLICT	UNP Q44467

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Chain	Residue	Modelled	Actual	Comment	Reference
G	255	ALA	ARG	CONFLICT	UNP Q44467
H	232	ALA	GLY	CONFLICT	UNP Q44467
H	255	ALA	ARG	CONFLICT	UNP Q44467

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C<sub>17</sub>H<sub>21</sub>N<sub>4</sub>O<sub>9</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	C	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	D	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	E	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	F	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	G	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	H	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total 1	Zn 1	0	0

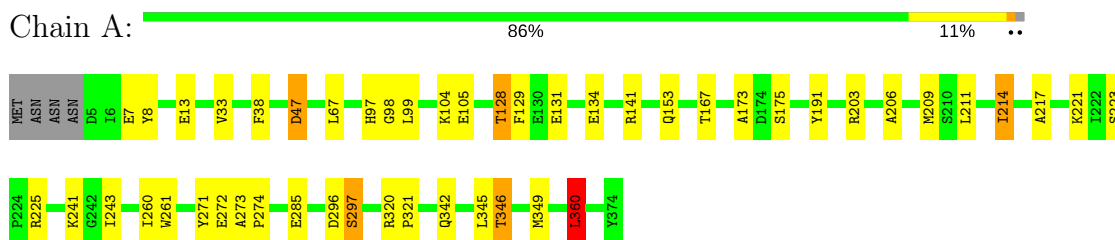
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	199	Total 199	O 199	0	0
4	B	185	Total 185	O 185	0	0
4	C	203	Total 203	O 203	0	0
4	D	145	Total 145	O 145	0	0
4	E	132	Total 132	O 132	0	0
4	F	148	Total 148	O 148	0	0
4	G	158	Total 158	O 158	0	0
4	H	160	Total 160	O 160	0	0

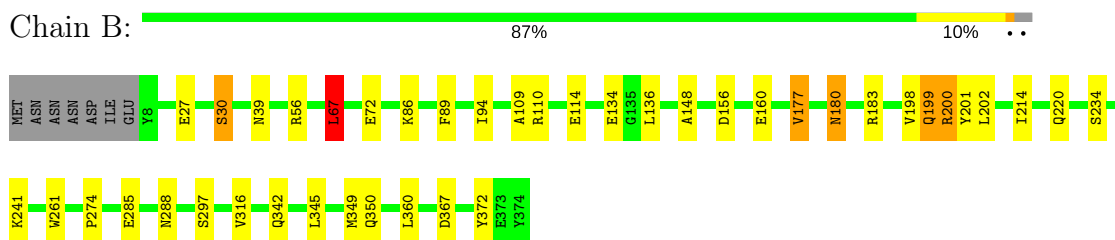
### 3 Residue-property plots

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.

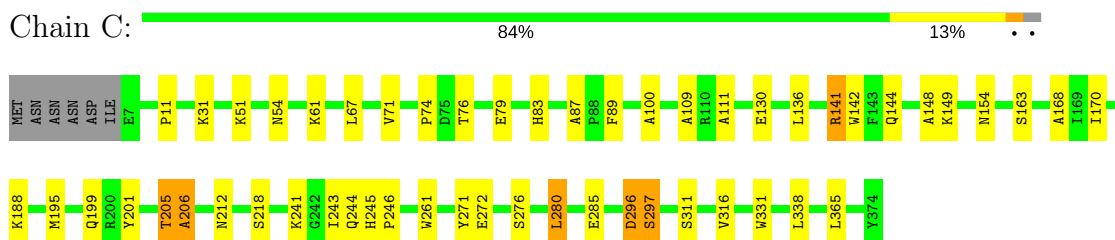
#### • Molecule 1: LACTATE OXIDASE



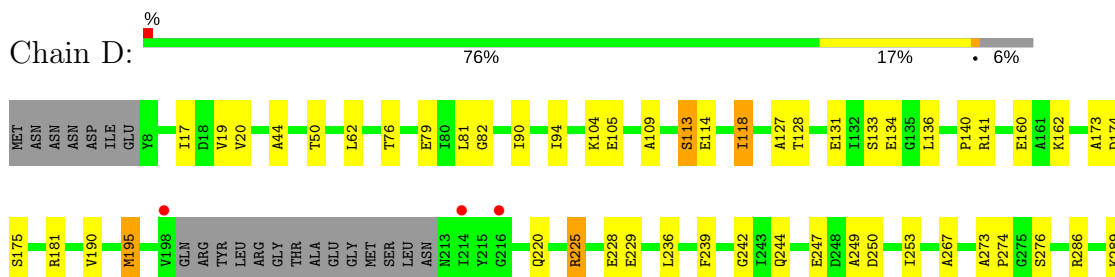
#### • Molecule 1: LACTATE OXIDASE

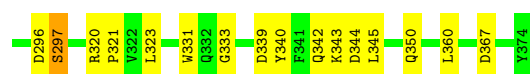


#### • Molecule 1: LACTATE OXIDASE

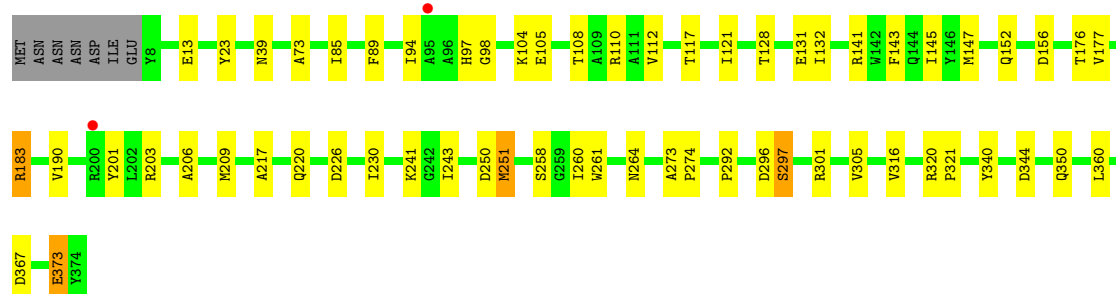
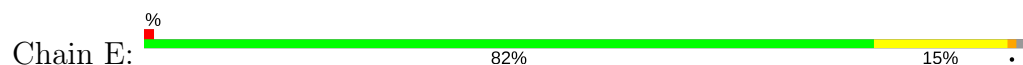


#### • Molecule 1: LACTATE OXIDASE

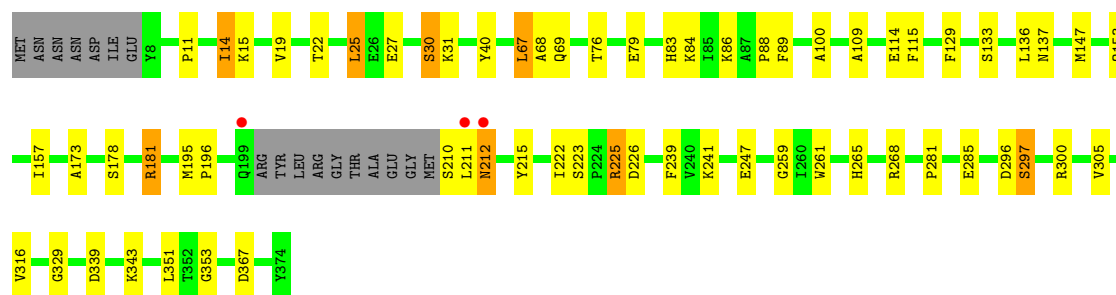
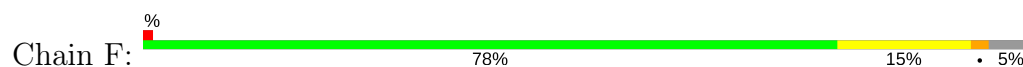




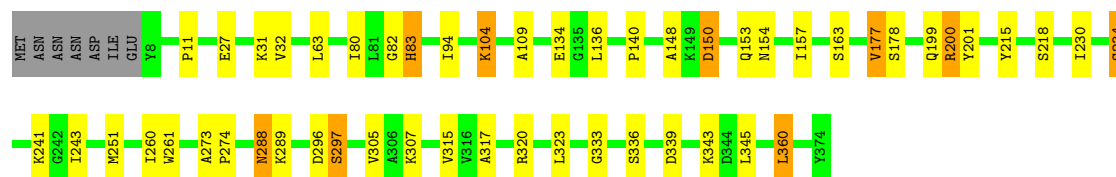
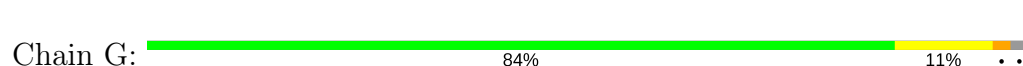
• Molecule 1: LACTATE OXIDASE



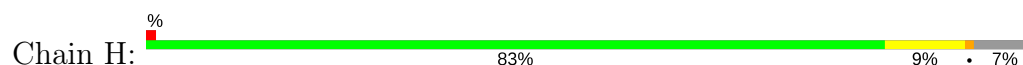
• Molecule 1: LACTATE OXIDASE



• Molecule 1: LACTATE OXIDASE



• Molecule 1: LACTATE OXIDASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	117.34Å 134.74Å 185.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.10 46.54 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (12.00-2.10) 100.0 (46.54-2.10)	Depositor EDS
$R_{merge}$	0.02	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.37 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.200 , 0.257 0.203 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	12.8	Xtriage
Anisotropy	0.847	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 36.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	23895	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	13.0	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 30.27 % 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 1.3449e-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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FMN, ZN

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	0.57	0/2916	0.65	1/3952 (0.0%)
1	B	0.52	0/2891	0.67	3/3918 (0.1%)
1	C	0.57	0/2900	0.65	0/3930
1	D	0.52	0/2779	0.63	0/3767
1	E	0.50	0/2891	0.61	0/3918
1	F	0.53	0/2810	0.63	1/3809 (0.0%)
1	G	0.52	0/2891	0.65	2/3918 (0.1%)
1	H	0.55	0/2743	0.65	1/3717 (0.0%)
All	All	0.54	0/22821	0.64	8/30929 (0.0%)

There are no bond length outliers.

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	320	ARG	NE-CZ-NH2	-7.17	116.71	120.30
1	G	320	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	B	199	GLN	C-N-CA	5.72	136.00	121.70
1	F	25	LEU	CA-CB-CG	5.65	128.31	115.30
1	B	199	GLN	N-CA-C	5.52	125.89	111.00

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	2852	0	2779	42	0
1	B	2827	0	2758	29	0
1	C	2836	0	2764	34	0
1	D	2717	0	2648	46	0
1	E	2827	0	2758	42	0
1	F	2748	0	2678	37	0
1	G	2827	0	2758	38	0
1	H	2682	0	2611	23	0
2	A	31	0	19	0	0
2	B	31	0	19	0	0
2	C	31	0	19	0	0
2	D	31	0	19	0	0
2	E	31	0	19	1	0
2	F	31	0	19	0	0
2	G	31	0	19	1	0
2	H	31	0	19	0	0
3	D	1	0	0	0	0
4	A	199	0	0	6	0
4	B	185	0	0	5	0
4	C	203	0	0	8	0
4	D	145	0	0	9	0
4	E	132	0	0	7	0
4	F	148	0	0	6	0
4	G	158	0	0	6	0
4	H	160	0	0	7	0
All	All	23895	0	21906	278	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.

The worst 5 of 278 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:F:40:TYR:O	1:F:268:ARG:NH2	1.82	1.10
1:C:141:ARG:HG2	1:C:141:ARG:HH11	1.24	0.99
1:G:82:GLY:HA3	1:G:83:HIS:HB2	1.51	0.93
1:C:71:VAL:HG13	4:C:2190:HOH:O	1.71	0.90
1:B:180:ASN:H	1:B:180:ASN:HD22	1.19	0.90

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	368/374 (98%)	359 (98%)	8 (2%)	1 (0%)	44	44
1	B	365/374 (98%)	352 (96%)	11 (3%)	2 (0%)	32	28
1	C	366/374 (98%)	352 (96%)	12 (3%)	2 (0%)	32	28
1	D	349/374 (93%)	338 (97%)	10 (3%)	1 (0%)	44	44
1	E	365/374 (98%)	351 (96%)	13 (4%)	1 (0%)	44	44
1	F	353/374 (94%)	339 (96%)	10 (3%)	4 (1%)	17	11
1	G	365/374 (98%)	350 (96%)	13 (4%)	2 (0%)	32	28
1	H	344/374 (92%)	335 (97%)	8 (2%)	1 (0%)	44	44
All	All	2875/2992 (96%)	2776 (97%)	85 (3%)	14 (0%)	32	28

5 of 14 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	297	SER
1	B	200	ARG
1	B	297	SER
1	C	297	SER
1	D	297	SER

### 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	293/297 (99%)	284 (97%)	9 (3%)	45	48

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	290/297 (98%)	280 (97%)	10 (3%)	42	43
1	C	291/297 (98%)	283 (97%)	8 (3%)	50	54
1	D	279/297 (94%)	263 (94%)	16 (6%)	24	21
1	E	290/297 (98%)	280 (97%)	10 (3%)	42	43
1	F	283/297 (95%)	272 (96%)	11 (4%)	37	37
1	G	290/297 (98%)	278 (96%)	12 (4%)	35	35
1	H	276/297 (93%)	267 (97%)	9 (3%)	43	45
All	All	2292/2376 (96%)	2207 (96%)	85 (4%)	39	39

5 of 85 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	220	GLN
1	E	183	ARG
1	H	80	ILE
1	D	225	ARG
1	E	13	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	244	GLN
1	D	244	GLN
1	G	244	GLN
1	C	69	GLN
1	C	154	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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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	FMN	A	1375	-	31,33,33	1.58	4 (12%)	38,50,50	1.44	3 (7%)
2	FMN	B	1375	-	31,33,33	1.33	5 (16%)	38,50,50	1.60	5 (13%)
2	FMN	C	1375	-	31,33,33	1.62	5 (16%)	38,50,50	1.91	7 (18%)
2	FMN	D	1375	-	31,33,33	1.43	5 (16%)	38,50,50	1.62	6 (15%)
2	FMN	E	1375	-	31,33,33	1.38	4 (12%)	38,50,50	1.61	6 (15%)
2	FMN	F	1375	-	31,33,33	1.46	5 (16%)	38,50,50	1.60	5 (13%)
2	FMN	G	1375	-	31,33,33	1.52	5 (16%)	38,50,50	1.72	4 (10%)
2	FMN	H	1375	-	31,33,33	1.41	4 (12%)	38,50,50	1.49	4 (10%)

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	FMN	A	1375	-	-	0/16/18/18	0/3/3/3
2	FMN	B	1375	-	-	0/16/18/18	0/3/3/3
2	FMN	C	1375	-	-	0/16/18/18	0/3/3/3
2	FMN	D	1375	-	-	0/16/18/18	0/3/3/3
2	FMN	E	1375	-	-	0/16/18/18	0/3/3/3
2	FMN	F	1375	-	-	0/16/18/18	0/3/3/3
2	FMN	G	1375	-	-	0/16/18/18	0/3/3/3
2	FMN	H	1375	-	-	0/16/18/18	0/3/3/3

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1375	FMN	C5A-N5	2.17	1.38	1.35
2	G	1375	FMN	C5A-N5	2.18	1.38	1.35
2	G	1375	FMN	C1'-N10	2.21	1.50	1.48
2	B	1375	FMN	C5A-N5	2.26	1.38	1.35
2	D	1375	FMN	C5A-N5	2.26	1.38	1.35

The worst 5 of 40 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1375	FMN	C4A-C4-N3	-3.03	119.16	123.48
2	G	1375	FMN	C4A-C4-N3	-2.94	119.30	123.48
2	C	1375	FMN	C4A-C4-N3	-2.94	119.30	123.48
2	F	1375	FMN	C4A-C4-N3	-2.59	119.80	123.48
2	H	1375	FMN	C4A-C4-N3	-2.43	120.03	123.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	1375	FMN	1	0
2	G	1375	FMN	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 [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	370/374 (98%)	-0.62	0 100 100	2, 9, 20, 27	0
1	B	367/374 (98%)	-0.55	0 100 100	3, 11, 21, 25	0
1	C	368/374 (98%)	-0.59	0 100 100	2, 8, 19, 28	0
1	D	353/374 (94%)	-0.32	3 (0%) 86 88	3, 14, 34, 49	0
1	E	367/374 (98%)	-0.29	2 (0%) 90 92	5, 17, 31, 40	0
1	F	357/374 (95%)	-0.37	3 (0%) 86 88	3, 12, 24, 43	0
1	G	367/374 (98%)	-0.47	0 100 100	4, 13, 25, 29	0
1	H	348/374 (93%)	-0.48	2 (0%) 89 91	3, 11, 26, 35	0
All	All	2897/2992 (96%)	-0.46	10 (0%) 93 94	2, 12, 26, 49	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	198	VAL	3.8
1	H	197	ILE	2.9
1	D	216	GLY	2.8
1	F	199	GLN	2.4
1	F	211	LEU	2.3

### 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

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	FMN	G	1375	31/31	0.97	0.10	0.30	5,9,11,11	0
2	FMN	E	1375	31/31	0.97	0.10	-0.29	9,11,13,13	0
2	FMN	F	1375	31/31	0.97	0.08	-0.30	3,7,8,9	0
2	FMN	H	1375	31/31	0.97	0.08	-0.33	5,8,10,13	0
2	FMN	B	1375	31/31	0.98	0.08	-0.38	5,7,9,9	0
2	FMN	C	1375	31/31	0.97	0.08	-0.41	2,5,7,8	0
2	FMN	A	1375	31/31	0.98	0.07	-0.45	2,6,8,9	0
2	FMN	D	1375	31/31	0.98	0.07	-0.81	5,8,11,12	0
3	ZN	D	1376	1/1	0.99	0.08	-1.62	26,26,26,26	0

## 6.5 Other polymers

There are no such residues in this entry.