



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 12:35 am GMT

PDB ID : 2YR4
Title : Crystal structure of L-phenylalanine oxiase from Psuedomonas sp. P-501
Authors : Ida, K.; Kurabayashi, M.; Suguro, M.; Suzuki, H.
Deposited on : 2007-04-02
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 : 1.7.2 (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) : recalc28949

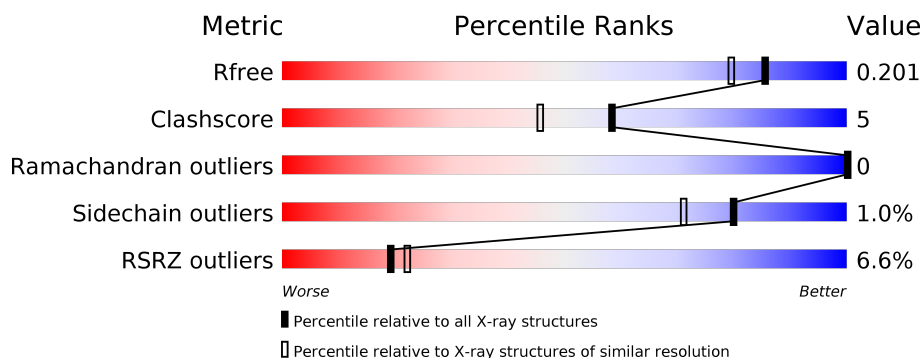
1 Overall quality at a glance

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}	100719	3453 (1.70-1.70)
Clashscore	112137	3876 (1.70-1.70)
Ramachandran outliers	110173	3815 (1.70-1.70)
Sidechain outliers	110143	3815 (1.70-1.70)
RSRZ outliers	101464	3491 (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	721	<div> <div>6%</div> <div> <div></div> <div>84%</div> <div>9%</div> <div>7%</div> </div> </div>
1	B	721	<div> <div>7%</div> <div> <div></div> <div>86%</div> <div>7%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11927 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.

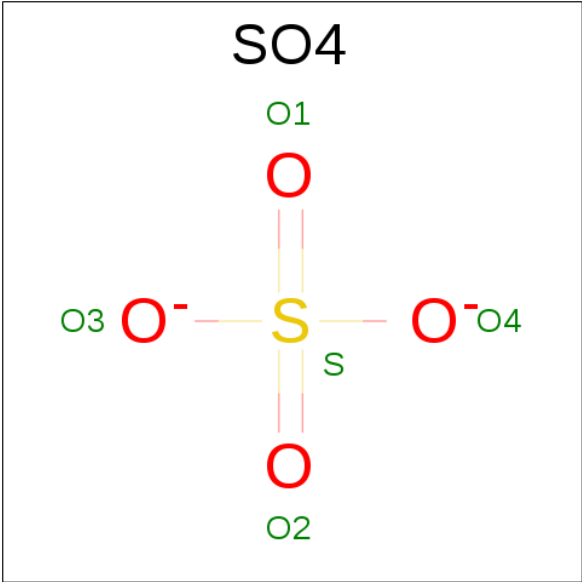
- Molecule 1 is a protein called Pro-enzyme of L-phenylalanine oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	674	Total	C	N	O	S	0	0	0
			5145	3284	892	958	11			
1	B	673	Total	C	N	O	S	0	0	0
			5145	3287	894	953	11			

There are 16 discrepancies between the modelled and reference sequences:

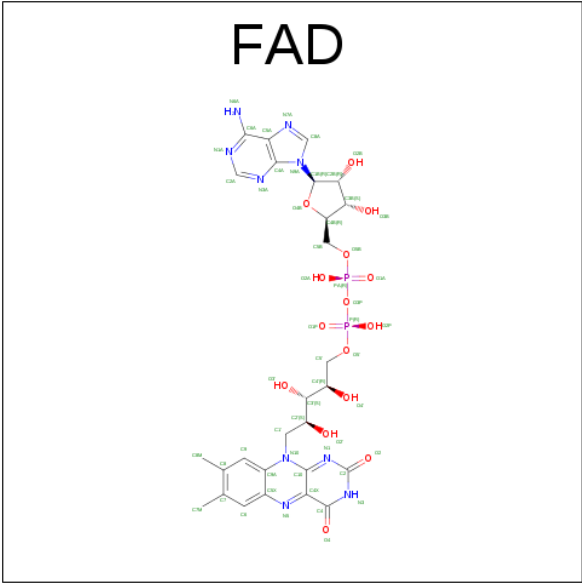
Chain	Residue	Modelled	Actual	Comment	Reference
A	714	LEU	-	EXPRESSION TAG	UNP Q5W9R9
A	715	GLU	-	EXPRESSION TAG	UNP Q5W9R9
A	716	HIS	-	EXPRESSION TAG	UNP Q5W9R9
A	717	HIS	-	EXPRESSION TAG	UNP Q5W9R9
A	718	HIS	-	EXPRESSION TAG	UNP Q5W9R9
A	719	HIS	-	EXPRESSION TAG	UNP Q5W9R9
A	720	HIS	-	EXPRESSION TAG	UNP Q5W9R9
A	721	HIS	-	EXPRESSION TAG	UNP Q5W9R9
B	714	LEU	-	EXPRESSION TAG	UNP Q5W9R9
B	715	GLU	-	EXPRESSION TAG	UNP Q5W9R9
B	716	HIS	-	EXPRESSION TAG	UNP Q5W9R9
B	717	HIS	-	EXPRESSION TAG	UNP Q5W9R9
B	718	HIS	-	EXPRESSION TAG	UNP Q5W9R9
B	719	HIS	-	EXPRESSION TAG	UNP Q5W9R9
B	720	HIS	-	EXPRESSION TAG	UNP Q5W9R9
B	721	HIS	-	EXPRESSION TAG	UNP Q5W9R9

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

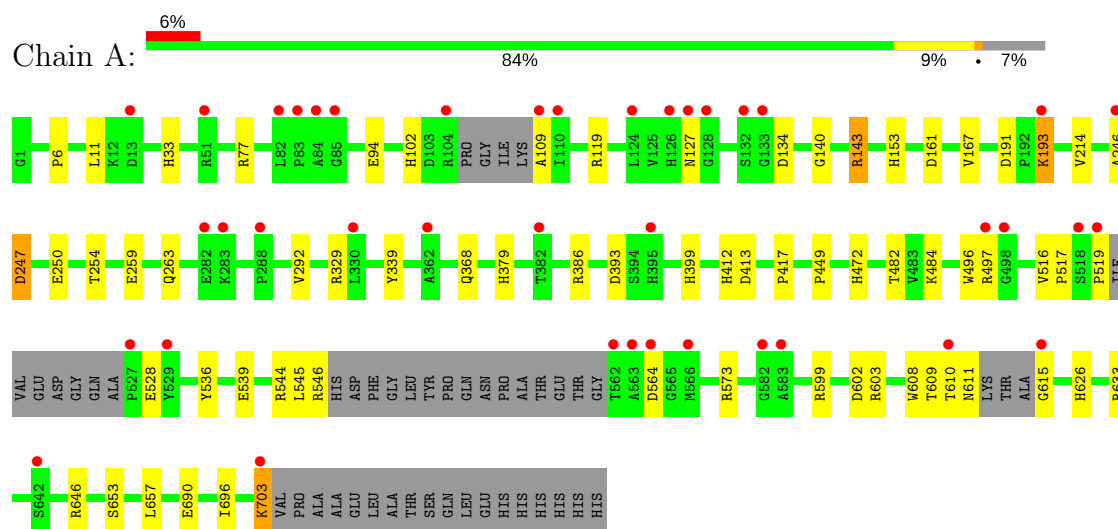
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	755	Total 755	O 755	0	0
4	B	771	Total 771	O 771	0	0

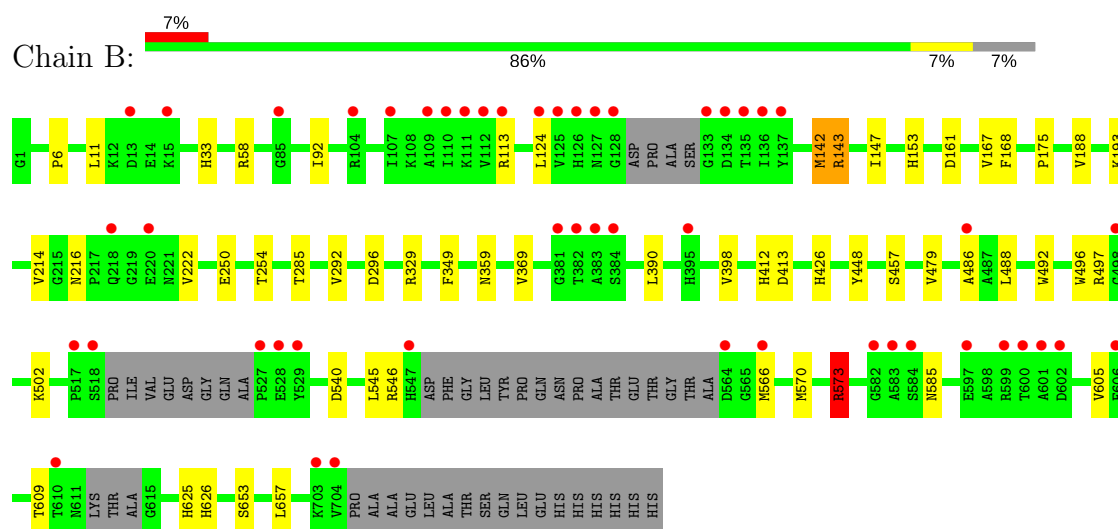
3 Residue-property plots

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: Pro-enzyme of L-phenylalanine oxidase



- Molecule 1: Pro-enzyme of L-phenylalanine oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	141.58Å 145.41Å 82.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.77 – 1.70 50.80 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.77-1.70) 99.3 (50.80-1.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.96 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.166 , 0.195 0.175 , 0.201	Depositor DCC
R_{free} test set	9271 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	19.0	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 62.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.010 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11927	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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: SO4, 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	0.80	0/5273	0.81	4/7189 (0.1%)
1	B	0.81	0/5273	0.81	2/7186 (0.0%)
All	All	0.80	0/10546	0.81	6/14375 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	386	ARG	NE-CZ-NH1	8.18	124.39	120.30
1	B	573	ARG	NE-CZ-NH2	-6.62	116.99	120.30
1	A	329	ARG	NE-CZ-NH1	6.20	123.40	120.30
1	A	161	ASP	CB-CG-OD1	5.86	123.57	118.30
1	A	633	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	B	161	ASP	CB-CG-OD1	5.01	122.81	118.30

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	5145	0	5060	62	0
1	B	5145	0	5070	55	0
2	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	53	0	31	2	0
3	B	53	0	31	2	0
4	A	755	0	0	23	0
4	B	771	0	0	16	0
All	All	11927	0	10192	113	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 5.

All (113) 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:127:ASN:HD21	1:A:599:ARG:NH2	1.44	1.14
1:A:127:ASN:HD21	1:A:599:ARG:CZ	1.87	0.86
1:B:479:VAL:HG21	1:B:570:MET:SD	2.19	0.83
1:B:143:ARG:NH1	3:B:801:FAD:O4	2.14	0.81
1:A:519:PRO:HA	4:A:1513:HOH:O	1.81	0.79
1:A:379:HIS:HE1	1:A:646:ARG:H	1.30	0.79
1:A:127:ASN:ND2	1:A:599:ARG:NH2	2.29	0.73
1:A:545:LEU:HD21	1:A:573:ARG:HG3	1.73	0.71
1:A:143:ARG:NH1	3:A:801:FAD:O4	2.24	0.70
1:A:496:TRP:CD1	1:A:497:ARG:HG2	2.25	0.70
1:A:472:HIS:HD2	1:B:540:ASP:OD2	1.75	0.69
1:A:134:ASP:O	4:A:1503:HOH:O	2.11	0.69
1:B:11:LEU:HD11	1:B:167:VAL:HG11	1.74	0.68
1:A:11:LEU:HD11	1:A:167:VAL:HG11	1.75	0.67
1:A:696:ILE:HD11	4:A:1259:HOH:O	1.96	0.65
1:B:412:HIS:HD2	1:B:653:SER:OG	1.80	0.64
1:A:412:HIS:HD2	1:A:653:SER:OG	1.80	0.64
1:B:58:ARG:HD2	4:B:2514:HOH:O	1.96	0.64
1:B:329:ARG:HH22	1:B:625:HIS:HD2	1.42	0.64
1:A:127:ASN:ND2	1:A:599:ARG:CZ	2.60	0.62
1:B:216:ASN:HB3	4:B:2664:HOH:O	2.00	0.61
1:A:393:ASP:OD2	1:A:399:HIS:HE1	1.83	0.61
1:B:113:ARG:C	1:B:124:LEU:HD11	2.21	0.61
1:A:472:HIS:HE1	1:A:539:GLU:OE2	1.84	0.60
1:A:690:GLU:HG3	4:A:1245:HOH:O	2.00	0.60
1:A:545:LEU:CD2	1:A:573:ARG:HG3	2.32	0.58
1:B:142:MET:CE	1:B:168:PHE:CZ	2.86	0.58
1:A:6:PRO:HB3	1:A:214:VAL:HG21	1.84	0.58
1:A:484:LYS:NZ	1:A:528:GLU:OE2	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:482:THR:HG21	1:A:599:ARG:HH21	1.69	0.58
1:B:488:LEU:HD13	1:B:502:LYS:HE2	1.86	0.57
1:B:329:ARG:HH22	1:B:625:HIS:CD2	2.21	0.57
1:A:413:ASP:HB2	4:A:1489:HOH:O	2.05	0.57
1:B:113:ARG:O	1:B:124:LEU:HD11	2.05	0.56
1:B:175:PRO:HA	1:B:188:VAL:HG22	1.88	0.54
1:B:216:ASN:HD21	1:B:222:VAL:CG1	2.21	0.54
1:A:246:ALA:O	1:A:247:ASP:HB2	2.08	0.54
1:B:216:ASN:ND2	1:B:222:VAL:HG13	2.23	0.53
1:B:142:MET:CE	1:B:168:PHE:CE1	2.92	0.53
1:B:585:ASN:HA	4:B:2246:HOH:O	2.07	0.53
3:B:801:FAD:O2'	3:B:801:FAD:H9	2.09	0.52
1:A:250:GLU:OE2	1:A:254:THR:HG21	2.10	0.52
1:A:379:HIS:CE1	1:A:646:ARG:H	2.19	0.52
1:A:77:ARG:HD3	4:A:962:HOH:O	2.09	0.52
1:B:413:ASP:HB2	4:B:2716:HOH:O	2.09	0.51
1:B:566:MET:HG2	1:B:605:VAL:HG11	1.91	0.51
1:B:359:ASN:HB2	4:B:2717:HOH:O	2.09	0.51
1:A:153:HIS:HE1	4:A:956:HOH:O	1.94	0.51
1:A:573:ARG:HD3	4:A:1485:HOH:O	2.11	0.51
1:B:33:HIS:CE1	1:B:657:LEU:HD11	2.47	0.51
1:A:94:GLU:O	1:A:102:HIS:HE1	1.94	0.50
1:B:250:GLU:OE2	1:B:254:THR:HG21	2.11	0.50
1:A:546:ARG:C	4:A:1258:HOH:O	2.50	0.50
1:A:626:HIS:HE1	4:A:1031:HOH:O	1.94	0.50
1:B:142:MET:HE3	1:B:168:PHE:CZ	2.46	0.50
1:A:615:GLY:C	4:A:1444:HOH:O	2.50	0.49
1:B:124:LEU:HD22	4:B:2727:HOH:O	2.10	0.49
1:B:142:MET:HE3	1:B:168:PHE:CE1	2.46	0.49
1:B:626:HIS:HE1	4:B:2103:HOH:O	1.95	0.49
1:B:486:ALA:HB3	4:B:2731:HOH:O	2.12	0.49
1:B:496:TRP:CD1	1:B:497:ARG:HG2	2.47	0.49
1:A:703:LYS:HE3	1:B:292:VAL:HG13	1.93	0.49
4:A:1198:HOH:O	1:B:626:HIS:HD2	1.96	0.48
1:A:646:ARG:CZ	4:A:1488:HOH:O	2.60	0.48
1:B:6:PRO:HB3	1:B:214:VAL:HG21	1.96	0.48
1:A:33:HIS:CE1	1:A:657:LEU:HD11	2.49	0.48
1:A:482:THR:HG21	1:A:599:ARG:NH2	2.28	0.47
1:B:285:THR:HG22	1:B:285:THR:O	2.14	0.47
1:B:193:LYS:NZ	4:B:2772:HOH:O	2.46	0.47
1:A:626:HIS:HD2	1:B:296:ASP:OD2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:609:THR:C	4:A:1445:HOH:O	2.52	0.47
1:A:519:PRO:CG	4:A:1508:HOH:O	2.62	0.47
1:A:615:GLY:N	4:A:1204:HOH:O	2.48	0.46
1:A:102:HIS:HD2	4:A:1093:HOH:O	1.98	0.46
1:A:610:THR:O	1:A:610:THR:HG22	2.15	0.46
1:A:167:VAL:HG22	1:A:339:TYR:CE1	2.50	0.46
1:B:573:ARG:NH2	4:B:2159:HOH:O	2.49	0.45
1:A:259:GLU:O	1:A:263:GLN:HG3	2.16	0.45
1:A:412:HIS:CD2	1:A:653:SER:OG	2.66	0.45
1:B:153:HIS:HE1	4:B:2087:HOH:O	1.99	0.45
1:B:492:TRP:HE3	4:B:2415:HOH:O	2.00	0.44
1:A:609:THR:O	1:B:609:THR:O	2.35	0.44
1:A:246:ALA:O	1:A:247:ASP:CB	2.66	0.44
3:A:801:FAD:H9	3:A:801:FAD:O2'	2.18	0.44
1:B:546:ARG:NH1	4:B:2666:HOH:O	2.42	0.44
1:B:570:MET:HB3	1:B:570:MET:HE3	1.88	0.43
1:A:564:ASP:OD1	1:A:603:ARG:NH1	2.52	0.43
1:B:216:ASN:ND2	1:B:222:VAL:CG1	2.81	0.43
1:A:191:ASP:OD1	1:A:193:LYS:HD3	2.18	0.43
1:A:544:ARG:NH2	4:A:855:HOH:O	2.50	0.43
1:B:412:HIS:HE1	4:B:2177:HOH:O	2.01	0.43
1:B:147:ILE:HD12	4:B:2356:HOH:O	2.19	0.43
1:A:608:TRP:O	1:B:546:ARG:NH2	2.53	0.42
1:A:127:ASN:ND2	1:A:602:ASP:OD2	2.52	0.42
1:A:127:ASN:O	1:A:127:ASN:OD1	2.37	0.42
1:A:292:VAL:HG21	1:B:457:SER:HB2	2.01	0.42
1:B:545:LEU:HD22	1:B:573:ARG:CZ	2.50	0.42
1:B:113:ARG:HB3	1:B:124:LEU:HD21	2.02	0.42
1:A:119:ARG:O	1:A:140:GLY:HA3	2.20	0.42
1:A:413:ASP:HB2	4:A:850:HOH:O	2.19	0.42
1:B:426:HIS:HB2	1:B:448:TYR:CE1	2.54	0.42
1:A:109:ALA:HB1	4:A:1550:HOH:O	2.20	0.41
1:A:516:VAL:HG13	1:A:517:PRO:HD2	2.01	0.41
1:A:703:LYS:CE	4:A:1411:HOH:O	2.68	0.41
1:B:390:LEU:HD11	1:B:398:VAL:CG2	2.51	0.41
1:A:703:LYS:HE3	4:A:1411:HOH:O	2.19	0.41
1:B:175:PRO:CA	1:B:188:VAL:HG22	2.50	0.41
1:B:92:ILE:CG2	1:B:369:VAL:HG22	2.51	0.41
1:A:417:PRO:HB2	4:A:1085:HOH:O	2.21	0.41
1:B:142:MET:CE	1:B:143:ARG:HD2	2.51	0.41
1:A:379:HIS:HD2	1:A:449:PRO:O	2.04	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:MET:HE2	1:B:143:ARG:HB3	2.02	0.41
1:B:413:ASP:HB2	4:B:2187:HOH:O	2.20	0.41

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	664/721 (92%)	648 (98%)	16 (2%)	0	100	100
1	B	663/721 (92%)	649 (98%)	14 (2%)	0	100	100
All	All	1327/1442 (92%)	1297 (98%)	30 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	529/566 (94%)	522 (99%)	7 (1%)	73	60
1	B	529/566 (94%)	525 (99%)	4 (1%)	85	78
All	All	1058/1132 (94%)	1047 (99%)	11 (1%)	80	71

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	143	ARG
1	A	193	LYS
1	A	247	ASP
1	A	368	GLN
1	A	536	TYR
1	A	611	ASN
1	A	703	LYS
1	B	142	MET
1	B	143	ARG
1	B	349	PHE
1	B	573	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	HIS
1	A	127	ASN
1	A	153	HIS
1	A	170	ASN
1	A	263	GLN
1	A	379	HIS
1	A	399	HIS
1	A	412	HIS
1	A	472	HIS
1	A	626	HIS
1	B	153	HIS
1	B	216	ASN
1	B	346	ASN
1	B	412	HIS
1	B	431	ASN
1	B	442	HIS
1	B	446	GLN
1	B	625	HIS
1	B	626	HIS

5.3.3 RNA ⓘ

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](#)

3 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	$\# Z > 2$
3	FAD	A	801	-	51,58,58	1.77	11 (21%)	54,89,89	1.93	11 (20%)
2	SO4	B	2001	-	4,4,4	0.24	0	6,6,6	0.29	0
3	FAD	B	801	-	51,58,58	1.66	12 (23%)	54,89,89	2.10	15 (27%)

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
3	FAD	A	801	-	-	0/28/50/50	0/6/6/6
2	SO4	B	2001	-	-	0/0/0/0	0/0/0/0
3	FAD	B	801	-	-	0/28/50/50	0/6/6/6

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	801	FAD	C4X-C10	-4.19	1.33	1.41
3	A	801	FAD	C4X-C10	-3.46	1.34	1.41
3	A	801	FAD	C9A-C5X	-3.20	1.36	1.42
3	B	801	FAD	C2-N1	-3.10	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	801	FAD	C4-C4X	-2.81	1.36	1.41
3	B	801	FAD	C9A-C5X	-2.79	1.36	1.42
3	A	801	FAD	C2-N3	-2.10	1.34	1.38
3	B	801	FAD	C2B-C1B	-2.09	1.50	1.53
3	A	801	FAD	C8M-C8	2.11	1.55	1.51
3	B	801	FAD	C10-N1	2.18	1.36	1.33
3	B	801	FAD	C4-N3	2.28	1.37	1.33
3	B	801	FAD	C9A-N10	2.37	1.41	1.38
3	B	801	FAD	C5X-N5	2.45	1.39	1.35
3	A	801	FAD	C9A-N10	2.57	1.42	1.38
3	A	801	FAD	C4-N3	2.64	1.37	1.33
3	B	801	FAD	C4X-N5	2.92	1.37	1.33
3	A	801	FAD	C5X-N5	2.97	1.39	1.35
3	B	801	FAD	C2A-N3A	3.56	1.38	1.32
3	A	801	FAD	C10-N1	3.67	1.38	1.33
3	B	801	FAD	C1'-N10	4.12	1.52	1.48
3	A	801	FAD	C2A-N3A	4.24	1.39	1.32
3	A	801	FAD	C4X-N5	4.71	1.40	1.33
3	A	801	FAD	C1'-N10	4.99	1.53	1.48

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801	FAD	C4-C4X-C10	-6.98	114.31	119.96
3	B	801	FAD	C1'-N10-C10	-6.17	112.17	118.50
3	A	801	FAD	C4-C4X-C10	-5.44	115.56	119.96
3	B	801	FAD	C4X-C10-N10	-4.87	117.14	120.52
3	A	801	FAD	C1'-N10-C10	-4.75	113.63	118.50
3	B	801	FAD	C9A-C5X-N5	-3.76	116.64	122.24
3	A	801	FAD	C4X-C10-N10	-3.14	118.34	120.52
3	A	801	FAD	C9A-C5X-N5	-2.91	117.91	122.24
3	B	801	FAD	C7M-C7-C8	-2.80	114.83	120.72
3	A	801	FAD	C7M-C7-C8	-2.71	115.03	120.72
3	B	801	FAD	O2'-C2'-C1'	-2.39	104.26	109.79
3	B	801	FAD	C4B-O4B-C1B	-2.15	107.48	109.77
3	B	801	FAD	C8M-C8-C9	-2.14	114.97	120.34
3	A	801	FAD	O2'-C2'-C1'	-2.05	105.04	109.79
3	A	801	FAD	C5X-C9A-N10	2.12	119.23	117.66
3	B	801	FAD	C5A-C6A-N6A	2.20	124.96	120.47
3	B	801	FAD	C4-C4X-N5	2.51	121.43	118.68
3	B	801	FAD	C4-N3-C2	2.83	117.63	115.16
3	B	801	FAD	C6-C5X-N5	2.98	122.46	118.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801	FAD	C10-C4X-N5	2.98	124.02	120.59
3	B	801	FAD	C4X-N5-C5X	3.07	120.00	116.76
3	A	801	FAD	C6-C5X-N5	3.25	122.78	118.97
3	A	801	FAD	C4X-N5-C5X	3.44	120.39	116.76
3	A	801	FAD	C4-C4X-N5	4.22	123.31	118.68
3	B	801	FAD	C5X-C9A-N10	4.38	120.91	117.66
3	A	801	FAD	C4-N3-C2	5.15	119.66	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	FAD	2	0
3	B	801	FAD	2	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	674/721 (93%)	0.25	40 (5%) 23 26	12, 23, 38, 53	0
1	B	673/721 (93%)	0.15	49 (7%) 16 19	11, 21, 41, 62	0
All	All	1347/1442 (93%)	0.20	89 (6%) 19 22	11, 22, 39, 62	0

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	133	GLY	10.0
1	B	128	GLY	8.6
1	A	562	THR	7.6
1	B	110	ILE	6.9
1	B	124	LEU	6.9
1	B	127	ASN	6.9
1	B	126	HIS	6.8
1	B	134	ASP	6.7
1	A	13	ASP	6.2
1	B	125	VAL	6.0
1	B	527	PRO	5.7
1	A	519	PRO	5.7
1	A	104	ARG	5.6
1	B	583	ALA	5.6
1	B	601	ALA	5.5
1	A	246	ALA	5.1
1	A	583	ALA	5.0
1	B	518	SER	5.0
1	A	563	ALA	4.6
1	B	529	TYR	4.6
1	B	703	LYS	4.2
1	A	518	SER	4.2
1	A	610	THR	4.1
1	A	582	GLY	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	126	HIS	4.1
1	B	104	ARG	4.0
1	A	109	ALA	4.0
1	B	13	ASP	3.9
1	B	136	ILE	3.9
1	B	564	ASP	3.8
1	B	486	ALA	3.8
1	B	111	LYS	3.7
1	B	704	VAL	3.7
1	B	600	THR	3.6
1	A	127	ASN	3.6
1	B	395	HIS	3.5
1	A	51	ARG	3.5
1	A	128	GLY	3.3
1	A	498	GLY	3.3
1	A	193	LYS	3.3
1	B	135	THR	3.2
1	A	497	ARG	3.1
1	B	602	ASP	3.1
1	B	606	PHE	3.1
1	A	83	PRO	3.1
1	B	382	THR	3.0
1	A	283	LYS	3.0
1	A	382	THR	3.0
1	B	384	SER	2.9
1	B	109	ALA	2.9
1	B	220	GLU	2.9
1	B	597	GLU	2.8
1	A	288	PRO	2.7
1	B	381	GLY	2.6
1	A	564	ASP	2.6
1	A	132	SER	2.6
1	A	362	ALA	2.6
1	A	615	GLY	2.6
1	B	498	GLY	2.6
1	B	15	LYS	2.6
1	A	84	ALA	2.5
1	A	133	GLY	2.5
1	A	82	LEU	2.5
1	A	703	LYS	2.5
1	B	547	HIS	2.5
1	A	527	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	582	GLY	2.4
1	A	642	SER	2.3
1	B	137	TYR	2.3
1	A	110	ILE	2.3
1	B	610	THR	2.3
1	B	113	ARG	2.3
1	A	395	HIS	2.3
1	B	584	SER	2.2
1	A	282	GLU	2.2
1	B	517	PRO	2.2
1	A	566	MET	2.2
1	B	383	ALA	2.2
1	B	85	GLY	2.2
1	B	112	VAL	2.1
1	A	330	LEU	2.1
1	A	85	GLY	2.1
1	B	218	GLN	2.1
1	B	566	MET	2.1
1	B	528	GLU	2.1
1	A	124	LEU	2.1
1	B	107	ILE	2.0
1	B	599	ARG	2.0
1	A	529	TYR	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(\AA^2)	Q<0.9
3	FAD	A	801	53/53	0.93	0.12	0.75	12,16,22,29	0
3	FAD	B	801	53/53	0.94	0.11	0.61	11,15,22,29	0
2	SO4	B	2001	5/5	0.93	0.24	0.24	59,60,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.