



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:47 PM GMT

PDB ID : 3S5W
Title : Ornithine Hydroxylase (PvdA) from Pseudomonas aeruginosa
Authors : Olucha, J.; Lamb, A.L.
Deposited on : 2011-05-23
Resolution : 1.90 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

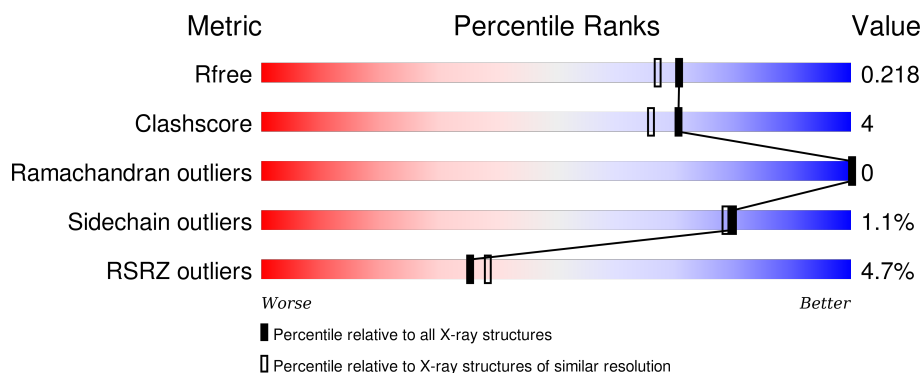
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	463	 4% 83% 5% • 11%
1	B	463	 4% 81% 8% 11%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7145 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 L-ornithine 5-monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	410	Total	C	N	O	S	0	2	0
			3267	2066	580	612	9			
1	B	414	Total	C	N	O	S	0	0	0
			3291	2080	585	617	9			

There are 40 discrepancies between the modelled and reference sequences:

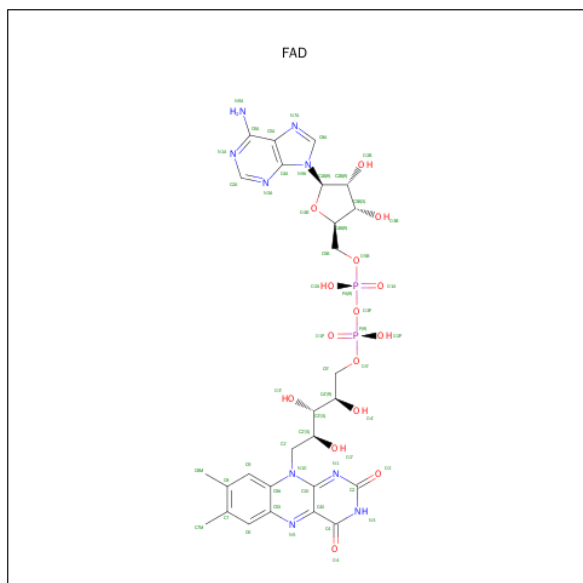
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q51548
A	-18	GLY	-	EXPRESSION TAG	UNP Q51548
A	-17	SER	-	EXPRESSION TAG	UNP Q51548
A	-16	SER	-	EXPRESSION TAG	UNP Q51548
A	-15	HIS	-	EXPRESSION TAG	UNP Q51548
A	-14	HIS	-	EXPRESSION TAG	UNP Q51548
A	-13	HIS	-	EXPRESSION TAG	UNP Q51548
A	-12	HIS	-	EXPRESSION TAG	UNP Q51548
A	-11	HIS	-	EXPRESSION TAG	UNP Q51548
A	-10	HIS	-	EXPRESSION TAG	UNP Q51548
A	-9	SER	-	EXPRESSION TAG	UNP Q51548
A	-8	SER	-	EXPRESSION TAG	UNP Q51548
A	-7	GLY	-	EXPRESSION TAG	UNP Q51548
A	-6	LEU	-	EXPRESSION TAG	UNP Q51548
A	-5	VAL	-	EXPRESSION TAG	UNP Q51548
A	-4	PRO	-	EXPRESSION TAG	UNP Q51548
A	-3	ARG	-	EXPRESSION TAG	UNP Q51548
A	-2	GLY	-	EXPRESSION TAG	UNP Q51548
A	-1	SER	-	EXPRESSION TAG	UNP Q51548
A	0	HIS	-	EXPRESSION TAG	UNP Q51548
B	-19	MET	-	EXPRESSION TAG	UNP Q51548
B	-18	GLY	-	EXPRESSION TAG	UNP Q51548
B	-17	SER	-	EXPRESSION TAG	UNP Q51548
B	-16	SER	-	EXPRESSION TAG	UNP Q51548
B	-15	HIS	-	EXPRESSION TAG	UNP Q51548

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	EXPRESSION TAG	UNP Q51548
B	-13	HIS	-	EXPRESSION TAG	UNP Q51548
B	-12	HIS	-	EXPRESSION TAG	UNP Q51548
B	-11	HIS	-	EXPRESSION TAG	UNP Q51548
B	-10	HIS	-	EXPRESSION TAG	UNP Q51548
B	-9	SER	-	EXPRESSION TAG	UNP Q51548
B	-8	SER	-	EXPRESSION TAG	UNP Q51548
B	-7	GLY	-	EXPRESSION TAG	UNP Q51548
B	-6	LEU	-	EXPRESSION TAG	UNP Q51548
B	-5	VAL	-	EXPRESSION TAG	UNP Q51548
B	-4	PRO	-	EXPRESSION TAG	UNP Q51548
B	-3	ARG	-	EXPRESSION TAG	UNP Q51548
B	-2	GLY	-	EXPRESSION TAG	UNP Q51548
B	-1	SER	-	EXPRESSION TAG	UNP Q51548
B	0	HIS	-	EXPRESSION TAG	UNP Q51548

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



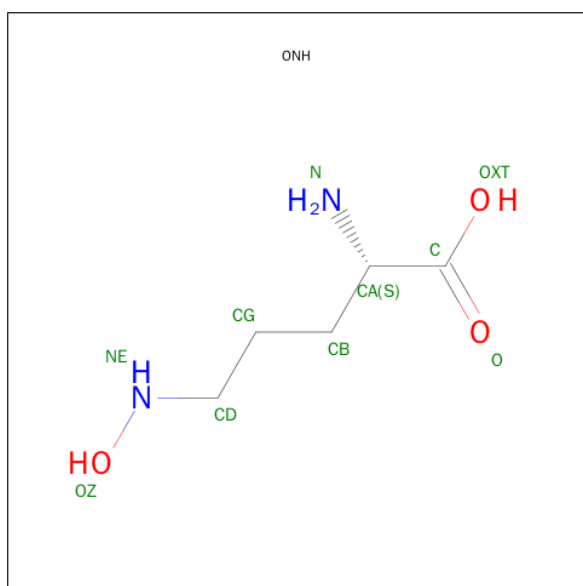
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



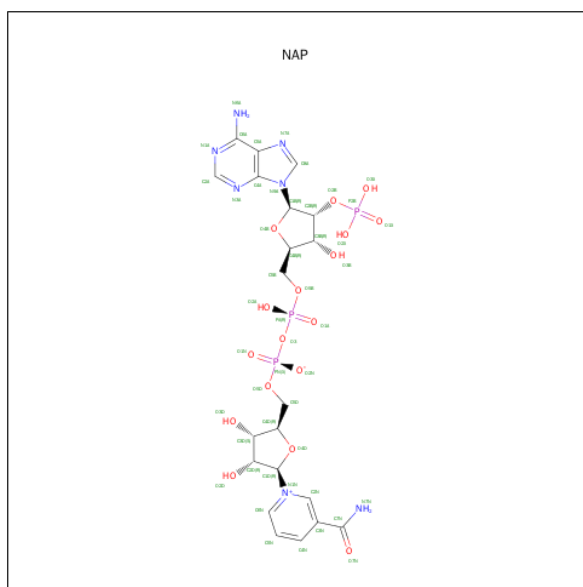
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is N 5 -HYDROXY-L-ORNITHINE (three-letter code: ONH) (formula: C₅H₁₂N₂O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			10	5	2	3		
4	B	1	Total	C	N	O	0	0
			10	5	2	3		

- Molecule 5 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			39	15	5	16	3		
5	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	181	Total	O	0	0
			181	181		
6	B	195	Total	O	0	0
			195	195		

4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	130.94Å 130.94Å 318.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.74 – 1.90 36.74 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.4 (36.74-1.90) 99.4 (36.74-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.192 , 0.217 0.195 , 0.218	Depositor DCC
R_{free} test set	5402 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 41.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 108129 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7145	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.55% 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, 0.375 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: PO4, NAP, FAD, ONH

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	0.71	0/3338	0.68	0/4516
1	B	0.73	0/3361	0.75	2/4552 (0.0%)
All	All	0.72	0/6699	0.72	2/9068 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	106	ARG	NE-CZ-NH2	-11.43	114.58	120.30
1	B	106	ARG	NE-CZ-NH1	8.99	124.80	120.30

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	3267	0	3226	29	0
1	B	3291	0	3236	30	0
2	A	53	0	31	0	0
2	B	53	0	31	0	0
3	A	10	0	0	0	0
3	B	5	0	0	0	0
4	A	10	0	11	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	10	0	11	0	0
5	A	39	0	18	1	0
5	B	31	0	11	0	0
6	A	181	0	0	2	0
6	B	195	0	0	3	0
All	All	7145	0	6575	59	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 4.

All (59) 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:B:235:ALA:H	1:B:311:HIS:HD2	1.16	0.92
1:A:358:GLN:O	1:A:359:LEU:HG	1.69	0.91
1:A:235:ALA:H	1:A:311:HIS:HD2	1.18	0.91
1:A:358:GLN:O	1:A:359:LEU:CG	2.30	0.80
1:B:281:HIS:HE1	6:B:489:HOH:O	1.64	0.80
1:B:56:THR:OG1	1:B:106:ARG:NH2	2.18	0.76
1:B:235:ALA:H	1:B:311:HIS:CD2	2.07	0.71
1:A:358:GLN:O	1:A:359:LEU:CD1	2.40	0.69
1:A:235:ALA:H	1:A:311:HIS:CD2	2.08	0.68
1:A:359:LEU:HD12	1:A:359:LEU:C	2.14	0.67
1:A:358:GLN:O	1:A:359:LEU:HD12	1.98	0.64
1:A:217:GLN:HB3	5:A:452:NAP:H52N	1.84	0.59
1:B:246:PRO:HB3	1:B:290:THR:HG22	1.83	0.59
1:A:189:HIS:CD2	1:A:191:SER:H	2.21	0.59
1:A:359:LEU:HD12	1:A:361:ARG:HD2	1.87	0.57
1:A:206:LYS:HG3	1:A:206:LYS:O	2.06	0.56
1:B:281:HIS:HD2	6:B:614:HOH:O	1.88	0.56
1:B:361:ARG:HG3	1:B:362:GLN:H	1.69	0.56
1:A:358:GLN:C	1:A:359:LEU:HG	2.26	0.55
1:B:190:HIS:HD2	6:B:559:HOH:O	1.89	0.55
1:B:13:LEU:HB3	1:B:41:VAL:HG22	1.88	0.55
1:B:212:ILE:HD11	1:B:226:LEU:HD12	1.89	0.55
1:B:361:ARG:CG	1:B:362:GLN:H	2.20	0.54
1:A:359:LEU:O	1:A:361:ARG:N	2.42	0.53
1:B:189:HIS:CD2	1:B:191:SER:H	2.26	0.53
1:A:168:PRO:HD2	1:A:397:PHE:HE2	1.74	0.52
1:A:359:LEU:O	1:A:359:LEU:HD12	2.10	0.51
1:B:367:LEU:HD13	1:B:392:ILE:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:PRO:HD2	1:B:397:PHE:HE2	1.75	0.51
1:A:359:LEU:CD1	1:A:359:LEU:C	2.78	0.51
1:B:378:ARG:O	1:B:402:HIS:HE1	1.94	0.50
1:A:369:GLU:O	1:A:369:GLU:HG3	2.11	0.50
1:A:281:HIS:HD2	6:A:514:HOH:O	1.93	0.50
1:A:166:VAL:HG12	1:A:168:PRO:HD3	1.94	0.48
1:B:361:ARG:HG3	1:B:362:GLN:N	2.28	0.48
1:A:189:HIS:HD2	1:A:191:SER:H	1.59	0.48
1:A:380:TYR:CE2	1:A:415:ARG:HG3	2.49	0.47
1:B:367:LEU:HD13	1:B:392:ILE:CD1	2.45	0.47
1:B:260:LYS:H	1:B:260:LYS:NZ	2.14	0.46
1:B:10:VAL:HG13	1:B:160:THR:HG22	1.98	0.46
1:B:224:ILE:HG13	1:B:292:LEU:HD11	1.97	0.46
1:A:401:SER:OG	1:A:402:HIS:HD2	1.99	0.46
1:B:189:HIS:HD2	1:B:191:SER:H	1.63	0.46
1:B:401:SER:OG	1:B:402:HIS:HD2	1.99	0.45
1:B:136:MET:SD	1:B:146:ARG:HG3	2.57	0.45
1:B:27:ILE:HD11	1:B:116:VAL:HG11	1.99	0.45
1:A:378:ARG:O	1:A:402:HIS:HE1	2.00	0.44
1:B:380:TYR:CE2	1:B:415:ARG:HG3	2.52	0.44
1:A:227:ASN:OD1	1:A:311:HIS:HE1	2.01	0.43
1:B:361:ARG:CG	1:B:362:GLN:N	2.81	0.43
1:A:105:CYS:HB3	6:A:558:HOH:O	2.17	0.43
1:A:168:PRO:HD2	1:A:397:PHE:CE2	2.53	0.43
1:A:359:LEU:O	1:A:361:ARG:HG2	2.19	0.43
1:B:96:PHE:O	1:B:99:LEU:HB2	2.19	0.42
1:B:227:ASN:OD1	1:B:311:HIS:HE1	2.02	0.42
1:B:189:HIS:HD2	1:B:191:SER:OG	2.02	0.42
1:A:224:ILE:HG13	1:A:292:LEU:HD11	2.03	0.41
1:A:64:GLN:NE2	1:A:217:GLN:HG2	2.35	0.41
1:B:321:ARG:HH11	1:B:323:THR:HG21	1.86	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	404/463 (87%)	394 (98%)	10 (2%)	0	100	100
1	B	410/463 (89%)	401 (98%)	9 (2%)	0	100	100
All	All	814/926 (88%)	795 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	348/386 (90%)	343 (99%)	5 (1%)	74	71
1	B	350/386 (91%)	347 (99%)	3 (1%)	84	83
All	All	698/772 (90%)	690 (99%)	8 (1%)	80	79

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	270	HIS
1	A	359	LEU
1	A	361	ARG
1	A	369	GLU
1	B	260	LYS
1	B	338	SER
1	B	369	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	GLN
1	A	76	ASN

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Mol	Chain	Res	Type
1	A	189	HIS
1	A	281	HIS
1	A	311	HIS
1	A	358	GLN
1	A	402	HIS
1	B	55	ASN
1	B	189	HIS
1	B	190	HIS
1	B	281	HIS
1	B	311	HIS
1	B	402	HIS

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](#)

9 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	PO4	A	444	-	4,4,4	0.36	0	6,6,6	0.29	0
4	ONH	A	445	-	5,9,9	0.36	0	3,10,10	1.36	0
3	PO4	A	446	-	4,4,4	0.37	0	6,6,6	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FAD	A	450	-	48,58,58	1.30	6 (12%)	54,89,89	2.33	9 (16%)
5	NAP	A	452	-	35,42,52	1.18	4 (11%)	46,65,80	2.35	10 (21%)
2	FAD	B	444	-	48,58,58	1.36	8 (16%)	54,89,89	2.45	9 (16%)
5	NAP	B	445	-	27,33,52	1.32	2 (7%)	34,52,80	2.51	4 (11%)
3	PO4	B	446	-	4,4,4	0.37	0	6,6,6	0.27	0
4	ONH	B	447	-	5,9,9	0.19	0	3,10,10	0.64	0

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	PO4	A	444	-	-	0/0/0/0	0/0/0/0
4	ONH	A	445	-	-	0/4/9/9	0/0/0/0
3	PO4	A	446	-	-	0/0/0/0	0/0/0/0
2	FAD	A	450	-	-	0/30/50/50	0/6/6/6
5	NAP	A	452	-	-	0/23/56/67	0/4/4/5
2	FAD	B	444	-	-	0/30/50/50	0/6/6/6
5	NAP	B	445	-	-	0/17/37/67	0/3/3/5
3	PO4	B	446	-	-	0/0/0/0	0/0/0/0
4	ONH	B	447	-	-	0/4/9/9	0/0/0/0

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	452	NAP	PA-O2A	-2.23	1.45	1.54
5	A	452	NAP	O4D-C4D	-2.09	1.40	1.44
5	A	452	NAP	O4B-C4B	-2.02	1.40	1.45
2	B	444	FAD	O4B-C4B	-2.02	1.40	1.45
2	B	444	FAD	C10-N1	2.06	1.39	1.35
2	A	450	FAD	C5X-N5	2.29	1.39	1.35
5	B	445	NAP	C2A-N1A	2.32	1.38	1.33
2	A	450	FAD	C2A-N1A	2.40	1.38	1.33
2	B	444	FAD	C5X-N5	2.47	1.39	1.35
2	B	444	FAD	C2A-N1A	2.50	1.38	1.33
2	A	450	FAD	C4-N3	2.50	1.37	1.33
2	A	450	FAD	C5'-C4'	2.55	1.55	1.51
2	B	444	FAD	C2A-N3A	3.24	1.37	1.32
2	B	444	FAD	C4-N3	3.25	1.39	1.33
2	B	444	FAD	C1'-N10	3.42	1.52	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	450	FAD	C4X-N5	3.74	1.39	1.33
5	A	452	NAP	C2A-N3A	4.04	1.39	1.32
2	B	444	FAD	C4X-N5	4.07	1.39	1.33
2	A	450	FAD	C2A-N3A	4.15	1.39	1.32
5	B	445	NAP	C2A-N3A	4.23	1.39	1.32

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	444	FAD	N3A-C2A-N1A	-12.98	118.96	128.89
5	B	445	NAP	N3A-C2A-N1A	-12.72	119.15	128.89
5	A	452	NAP	N3A-C2A-N1A	-12.55	119.29	128.89
2	A	450	FAD	N3A-C2A-N1A	-12.38	119.42	128.89
2	A	450	FAD	C4X-C4-N3	-3.77	118.43	123.59
2	B	444	FAD	C1B-N9A-C4A	-3.43	121.77	126.94
5	B	445	NAP	C1B-N9A-C4A	-3.28	121.99	126.94
2	A	450	FAD	C1B-N9A-C4A	-3.08	122.30	126.94
2	B	444	FAD	C4X-C4-N3	-3.01	119.47	123.59
2	A	450	FAD	O2'-C2'-C1'	-2.54	103.71	109.94
5	A	452	NAP	O5D-C5D-C4D	-2.49	99.94	109.12
5	A	452	NAP	O2B-P2B-O1X	-2.42	101.07	107.11
5	A	452	NAP	O3-PN-O5D	-2.41	96.55	102.94
5	A	452	NAP	O3-PA-O5B	-2.33	96.76	102.94
2	B	444	FAD	O2'-C2'-C1'	-2.05	104.91	109.94
5	A	452	NAP	O4B-C1B-C2B	-2.04	102.91	106.60
2	B	444	FAD	C4-C4X-C10	-2.01	118.66	119.94
5	B	445	NAP	O4B-C4B-C5B	-2.00	102.15	109.32
5	A	452	NAP	C2A-N1A-C6A	2.04	122.41	118.77
2	A	450	FAD	C2A-N1A-C6A	2.08	122.48	118.77
2	B	444	FAD	C2A-N1A-C6A	2.09	122.50	118.77
5	A	452	NAP	O4B-C1B-N9A	2.15	112.61	108.10
5	B	445	NAP	O4B-C1B-N9A	2.41	113.14	108.10
2	A	450	FAD	C4X-N5-C5X	2.68	119.84	116.76
5	A	452	NAP	O2D-C2D-C3D	2.90	116.79	111.23
2	A	450	FAD	C5X-C9A-N10	3.00	119.90	117.62
2	A	450	FAD	C1'-N10-C9A	3.10	122.34	118.86
5	A	452	NAP	O4D-C4D-C5D	3.12	116.37	109.53
2	B	444	FAD	C1'-N10-C9A	3.20	122.45	118.86
2	B	444	FAD	C5X-C9A-N10	4.27	120.87	117.62
2	B	444	FAD	C4-N3-C2	6.88	121.20	115.25
2	A	450	FAD	C4-N3-C2	6.98	121.28	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	452	NAP	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 ⓘ

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	410/463 (88%)	0.09	20 (4%) 33 36	20, 29, 51, 63	5 (1%)
1	B	414/463 (89%)	0.24	19 (4%) 36 39	19, 30, 50, 63	0
All	All	824/926 (88%)	0.17	39 (4%) 35 38	19, 30, 50, 63	5 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	359	LEU	10.3
1	B	337	GLY	6.7
1	A	206	LYS	6.1
1	B	429	PRO	4.5
1	B	198	ALA	4.1
1	A	386	GLU	3.9
1	A	139	ALA	3.9
1	B	207	PRO	3.7
1	B	9	VAL	3.6
1	B	327	GLN	3.6
1	A	138	SER	3.5
1	B	341	LEU	3.4
1	B	326	ALA	3.2
1	B	271	ALA	3.2
1	B	362	GLN	3.1
1	A	270	HIS	3.0
1	A	141	GLN	2.9
1	A	370	TYR	2.8
1	A	383	GLN	2.8
1	B	340	GLU	2.8
1	A	384	THR	2.7
1	B	199	LYS	2.7
1	A	365	GLU	2.5
1	A	387	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	270	HIS	2.5
1	A	389	LYS	2.4
1	A	388	CYS	2.4
1	A	137	LEU	2.4
1	B	153	ASP	2.3
1	B	361	ARG	2.3
1	A	358	GLN	2.3
1	B	152	ALA	2.2
1	B	321	ARG	2.1
1	A	143	GLU	2.1
1	A	34	GLN	2.1
1	B	155	GLU	2.0
1	B	365	GLU	2.0
1	A	140	GLY	2.0
1	A	9	VAL	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
4	ONH	B	447	10/10	0.94	0.12	0.11	22,24,28,41	0
2	FAD	B	444	53/53	0.98	0.12	0.10	18,24,29,32	0
4	ONH	A	445	10/10	0.96	0.12	0.07	22,24,26,41	0
5	NAP	A	452	39/48	0.96	0.10	0.03	21,27,53,54	0
2	FAD	A	450	53/53	0.98	0.10	-0.19	19,25,30,32	0
5	NAP	B	445	31/48	0.96	0.09	-0.99	26,29,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	PO4	A	446	5/5	0.87	0.32	-	66,66,67,68	0
3	PO4	B	446	5/5	0.73	0.25	-	65,66,67,68	0
3	PO4	A	444	5/5	0.92	0.13	-	50,51,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.