



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:21 PM GMT

PDB ID : 4EM4  
Title : Crystal Structure of Staphylococcus aureus bound with the covalent inhibitor Pethyl-VS-CoA  
Authors : Wallace, B.D.; Edwards, J.S.; Claiborne, A.; Redinbo, M.R.  
Deposited on : 2012-04-11  
Resolution : 1.82 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](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 (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

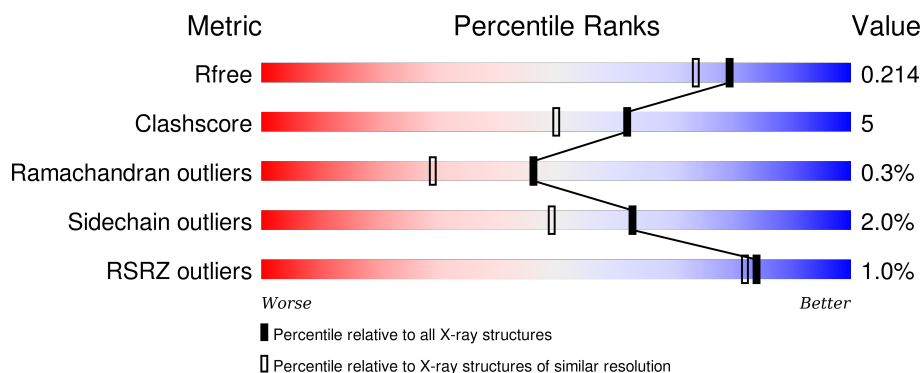
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.82 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	5422 (1.84-1.80)
Clashscore	102246	6347 (1.84-1.80)
Ramachandran outliers	100387	6276 (1.84-1.80)
Sidechain outliers	100360	6276 (1.84-1.80)
RSRZ outliers	91569	5439 (1.84-1.80)

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	437	<div> <div></div> <div>90%10%</div> </div>
1	B	437	<div> <div></div> <div>91%8%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	CA8	A	508	-	-	-	X
5	CA8	B	506	-	-	-	X

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8365 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 Coenzyme A disulfide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	437	Total	C	N	O	S	0	13	0
			3522	2240	597	672	13			
1	B	437	Total	C	N	O	S	0	13	0
			3531	2243	598	677	13			

- 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		

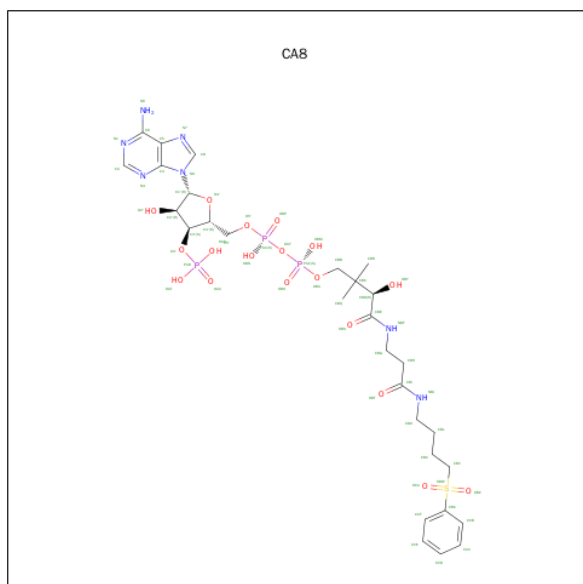
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0
3	A	2	Total Mg 2 2	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	3	Total Cl 3 3	0	0
4	A	4	Total Cl 4 4	0	0

- Molecule 5 is [(2R,3S,4R,5R)-5-(6-AMINOPURIN-9-YL)-4-OXIDANYL-3-PHOSPHONO OXY-OXOLAN-2-YL]METHOXY-OXIDANYL-PHOSPHORYL [(3R)-2,2-DIMETHYL-3-OXIDANYL-4-OXIDANYLIDENE-4-[[3-OXIDANYLIDENE-3-[4-(PHENYLSULFONYL)BUTYLAMINO]PROPYL]AMINO]BUTYL] HYDROGEN PHOSPHATE (three-letter code: CA8) (formula: C<sub>29</sub>H<sub>44</sub>N<sub>7</sub>O<sub>18</sub>P<sub>3</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total 58	C 29	N 7	O 18	P 3	S 1	0	0
5	B	1	Total 58	C 29	N 7	O 18	P 3	S 1	0	0

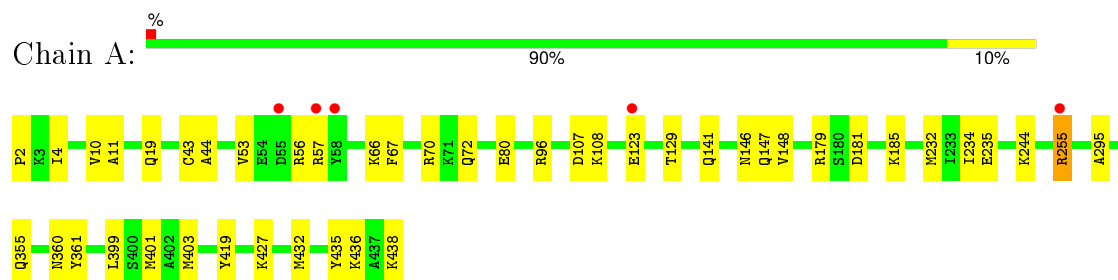
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	564	Total 564	O 564	0	0
6	B	516	Total 516	O 516	0	0

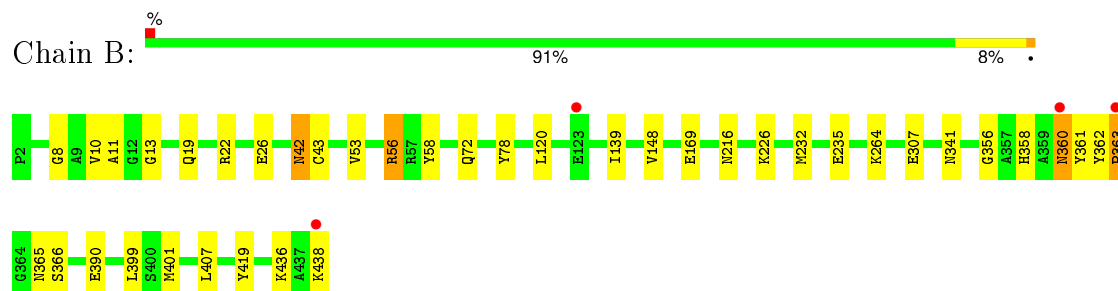
### 3 Residue-property plots [i](#)

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



- Molecule 1: Coenzyme A disulfide reductase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	76.14Å 64.91Å 94.44Å 90.00° 104.62° 90.00°	Depositor
Resolution (Å)	32.45 – 1.82 32.45 – 1.82	Depositor EDS
% Data completeness (in resolution range)	95.8 (32.45-1.82) 95.7 (32.45-1.82)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	10.77 (at 1.82Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1057)	Depositor
R, $R_{free}$	0.153 , 0.202 0.158 , 0.214	Depositor DCC
$R_{free}$ test set	3828 reflections (5.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	20.7	Xtriage
Anisotropy	0.637	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 59.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 76546 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	8365	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, CA8, FAD, CL

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	0.49	0/3637	0.67	0/4920
1	B	0.48	0/3645	0.67	0/4931
All	All	0.49	0/7282	0.67	0/9851

There are no bond length outliers.

There are no bond angle outliers.

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	3522	0	3507	38	0
1	B	3531	0	3510	36	1
2	A	53	0	31	1	0
2	B	53	0	31	4	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
4	A	4	0	0	0	0
4	B	3	0	0	1	0
5	A	58	0	39	8	0
5	B	58	0	38	9	0
6	A	564	0	0	20	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	516	0	0	15	1
All	All	8365	0	7156	78	3

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 (78) 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)
5:A:508:CA8:OBO	6:A:1161:HOH:O	1.52	1.25
1:A:355:GLN:NE2	1:A:432[B]:MET:SD	2.33	1.00
5:A:508:CA8:HBQA	6:A:1160:HOH:O	1.61	0.99
4:B:505:CL:CL	6:B:1030:HOH:O	2.27	0.89
1:B:360:ASN:OD1	6:B:1106:HOH:O	1.96	0.82
1:A:244:LYS:NZ	6:A:855:HOH:O	2.15	0.79
1:A:436:LYS:NZ	6:A:1061:HOH:O	2.17	0.78
1:B:360:ASN:OD1	1:B:360:ASN:N	2.14	0.76
1:B:53:VAL:O	6:B:809:HOH:O	2.11	0.68
1:A:179:ARG:NH1	6:A:840:HOH:O	2.27	0.68
1:B:307:GLU:OE2	6:B:854:HOH:O	2.11	0.68
1:B:72:GLN:OE1	6:B:816:HOH:O	2.11	0.67
1:A:360:ASN:OD1	6:A:1123:HOH:O	2.12	0.67
1:B:120:LEU:O	6:B:949:HOH:O	2.13	0.65
5:A:508:CA8:OCA	1:B:361:TYR:OH	2.15	0.64
1:B:436:LYS:NZ	6:B:885:HOH:O	2.33	0.61
5:A:508:CA8:HCE	6:B:1113:HOH:O	2.02	0.59
1:A:123:GLU:O	6:A:931:HOH:O	2.17	0.59
6:A:1150:HOH:O	5:B:506:CA8:HCC	2.02	0.58
1:B:43:CYS:SG	2:B:501:FAD:C4X	2.92	0.58
1:A:72:GLN:OE1	6:A:1089:HOH:O	2.17	0.58
1:A:147:GLN:O	6:A:1143:HOH:O	2.18	0.56
1:B:22:ARG:HD2	6:B:980:HOH:O	2.06	0.56
1:A:129:THR:HG22	1:A:235[A]:GLU:HG3	1.88	0.55
1:A:419:TYR:CE2	5:B:506:CA8:HBV	2.42	0.55
1:A:438:LYS:HE2	6:A:753:HOH:O	2.06	0.54
1:A:148:VAL:HG11	1:A:232[A]:MET:HE2	1.88	0.54
1:A:295:ALA:HB2	5:A:508:CA8:HBL	1.90	0.53
1:A:53:VAL:HG22	1:B:361:TYR:HB2	1.89	0.53
1:A:419:TYR:HE2	5:B:506:CA8:CBV	2.23	0.51
1:A:361:TYR:HB3	1:B:58:TYR:CG	2.45	0.51
1:A:427[B]:LYS:NZ	6:A:998:HOH:O	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:ALA:HB3	2:B:501:FAD:O5'	2.12	0.50
1:A:80:GLU:OE1	1:A:96[B]:ARG:NH2	2.46	0.49
1:B:78:TYR:OH	6:B:829:HOH:O	2.19	0.49
1:A:419:TYR:CE2	5:B:506:CA8:CBV	2.95	0.49
1:B:139:ILE:HG23	1:B:232[B]:MET:HE1	1.96	0.48
1:B:356:GLY:HA3	1:B:365:ASN:HD21	1.78	0.48
1:B:72:GLN:HG3	6:B:1024:HOH:O	2.13	0.48
1:B:42:ASN:OD1	5:B:506:CA8:HBHA	2.14	0.47
1:A:66:LYS:NZ	6:A:973:HOH:O	2.47	0.47
1:A:355:GLN:HG3	6:A:1120:HOH:O	2.14	0.47
1:A:185:LYS:HG2	6:A:662:HOH:O	2.13	0.47
1:B:264:LYS:NZ	6:B:888:HOH:O	2.46	0.47
1:A:53:VAL:CG2	1:B:361:TYR:HB2	2.46	0.46
1:B:19:GLN:NE2	6:B:980:HOH:O	2.48	0.46
1:A:255:ARG:CZ	1:A:255:ARG:HB2	2.46	0.46
1:B:366:SER:OG	6:B:1105:HOH:O	2.20	0.45
1:A:419:TYR:HE2	5:B:506:CA8:HBV	1.82	0.45
1:A:44:ALA:HB2	1:B:361:TYR:CZ	2.52	0.45
1:A:107:ASP:O	1:A:108:LYS:HD3	2.18	0.44
5:A:508:CA8:HBVA	1:B:419:TYR:CE2	2.53	0.43
1:B:358:HIS:HD2	1:B:390:GLU:OE1	2.01	0.43
1:A:419:TYR:OH	5:B:506:CA8:HBVA	2.18	0.43
2:B:501:FAD:H9	2:B:501:FAD:H1'1	1.82	0.43
1:A:19:GLN:NE2	6:A:1161:HOH:O	2.48	0.43
6:A:1150:HOH:O	5:B:506:CA8:HCB	2.18	0.43
1:A:2:PRO:HG2	1:A:4[B]:ILE:HD11	2.01	0.43
1:A:11:ALA:HB3	2:A:501:FAD:O5'	2.20	0.42
1:B:226:LYS:HB2	1:B:226:LYS:HE2	1.53	0.42
1:B:22:ARG:HH12	5:B:506:CA8:H2'	1.84	0.42
1:B:399:LEU:HD23	1:B:399:LEU:HA	1.90	0.42
1:A:43:CYS:H	5:A:508:CA8:HBUA	1.85	0.42
1:A:123:GLU:HG3	1:A:123:GLU:O	2.18	0.42
1:B:148:VAL:HG11	1:B:232[A]:MET:HE2	2.02	0.42
1:A:438:LYS:O	6:A:746:HOH:O	2.21	0.42
1:B:8:GLY:O	1:B:13:GLY:HA3	2.20	0.41
5:A:508:CA8:HBV	5:A:508:CA8:HBK	1.86	0.41
1:B:169:GLU:HG3	6:B:1082:HOH:O	2.21	0.41
1:A:399:LEU:O	1:A:403:MET:HG3	2.21	0.41
1:A:427[B]:LYS:HE2	1:A:435:TYR:CE2	2.56	0.41
1:B:407:LEU:HD12	1:B:407:LEU:HA	1.86	0.41
1:B:362:TYR:HA	1:B:363:PRO:HD2	1.66	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ASN:ND2	6:A:1134:HOH:O	2.38	0.40
1:A:401:MET:HG3	1:B:401:MET:HG3	2.04	0.40
1:B:43:CYS:SG	2:B:501:FAD:C4	3.10	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:973:HOH:O	6:A:1121:HOH:O[2_646]	2.00	0.20
1:B:26[B]:GLU:OE1	6:A:890:HOH:O[2_556]	2.03	0.17
6:B:1021:HOH:O	6:B:1104:HOH:O[2_647]	2.16	0.04

## 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	448/437 (102%)	437 (98%)	10 (2%)	1 (0%)	52	35
1	B	448/437 (102%)	434 (97%)	12 (3%)	2 (0%)	39	23
All	All	896/874 (102%)	871 (97%)	22 (2%)	3 (0%)	46	29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	B	363	PRO
1	B	10	VAL

### 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	388/376 (103%)	380 (98%)	8 (2%)	61	47
1	B	389/376 (104%)	380 (98%)	9 (2%)	58	42
All	All	777/752 (103%)	760 (98%)	17 (2%)	63	45

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	A	57	ARG
1	A	67	PHE
1	A	70	ARG
1	A	141	GLN
1	A	181	ASP
1	A	234	ILE
1	A	255	ARG
1	B	42	ASN
1	B	56[A]	ARG
1	B	56[B]	ARG
1	B	216	ASN
1	B	235[A]	GLU
1	B	235[B]	GLU
1	B	341	ASN
1	B	360	ASN
1	B	438	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 14 ligands modelled in this entry, 10 are monoatomic - leaving 4 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	FAD	A	501	-	48,58,58	1.19	5 (10%)	54,89,89	2.11	7 (12%)
5	CA8	A	508	1	51,61,61	1.84	13 (25%)	67,91,91	3.04	26 (38%)
2	FAD	B	501	-	48,58,58	1.23	5 (10%)	54,89,89	2.11	7 (12%)
5	CA8	B	506	1	51,61,61	2.97	21 (41%)	67,91,91	3.55	24 (35%)

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	FAD	A	501	-	-	0/30/50/50	0/6/6/6
5	CA8	A	508	1	-	0/55/75/75	0/4/4/4
2	FAD	B	501	-	-	0/30/50/50	0/6/6/6
5	CA8	B	506	1	-	0/55/75/75	0/4/4/4

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	506	CA8	PAR-O3'	-9.05	1.32	1.60
5	B	506	CA8	PAX-O5'	-7.13	1.26	1.59
5	A	508	CA8	CBX-SBW	-5.15	1.66	1.76
5	B	506	CA8	O5'-C5'	-4.14	1.27	1.44
5	B	506	CA8	CBU-CBL	-3.90	1.29	1.51
5	A	508	CA8	C2'-C3'	-3.52	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	508	CA8	C5-C4	-3.27	1.33	1.40
5	A	508	CA8	PAX-OBN	-2.75	1.43	1.54
2	B	501	FAD	C10-N10	-2.50	1.36	1.39
5	B	506	CA8	C5-C4	-2.43	1.35	1.40
5	A	508	CA8	PAR-OAT	-2.41	1.46	1.54
5	B	506	CA8	PAZ-OBM	-2.36	1.44	1.54
5	B	506	CA8	PAX-OBN	-2.29	1.45	1.54
5	B	506	CA8	C2'-C3'	-2.26	1.47	1.53
5	A	508	CA8	OCA-SBW	2.14	1.47	1.44
5	A	508	CA8	PAR-OAU	2.22	1.58	1.51
5	B	506	CA8	PAR-OAU	2.28	1.58	1.51
5	B	506	CA8	CBH-CBI	2.39	1.56	1.51
2	A	501	FAD	C5X-N5	2.40	1.39	1.35
5	B	506	CA8	CBG-NBF	2.41	1.51	1.46
2	A	501	FAD	C4-N3	2.54	1.37	1.33
5	A	508	CA8	C3'-C4'	2.57	1.60	1.52
5	B	506	CA8	C2-N1	2.58	1.38	1.33
2	A	501	FAD	C2A-N1A	2.59	1.38	1.33
5	A	508	CA8	CBH-CBI	2.61	1.56	1.51
2	B	501	FAD	C4-N3	2.75	1.38	1.33
5	B	506	CA8	PAR-OAS	2.75	1.64	1.54
5	A	508	CA8	OBZ-SBW	2.82	1.48	1.44
2	B	501	FAD	C2A-N1A	2.83	1.39	1.33
5	B	506	CA8	O2'-C2'	3.09	1.50	1.43
2	A	501	FAD	C4X-N5	3.22	1.38	1.33
5	A	508	CA8	O4'-C1'	3.23	1.45	1.41
2	B	501	FAD	C4X-N5	3.24	1.38	1.33
5	B	506	CA8	C5'-C4'	3.35	1.62	1.51
5	B	506	CA8	PAZ-OBO	3.52	1.64	1.51
2	A	501	FAD	C2A-N3A	3.60	1.38	1.32
5	A	508	CA8	PAX-OBP	3.64	1.64	1.51
2	B	501	FAD	C2A-N3A	3.71	1.38	1.32
5	A	508	CA8	C5'-C4'	3.76	1.63	1.51
5	B	506	CA8	C2-N3	3.83	1.39	1.32
5	B	506	CA8	C3'-C4'	4.36	1.65	1.52
5	B	506	CA8	CBV-SBW	5.26	1.89	1.78
5	B	506	CA8	O4'-C1'	6.62	1.49	1.41
5	B	506	CA8	CBX-SBW	7.48	1.91	1.76

All (64) bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	508	CA8	N3-C2-N1	-13.07	118.89	128.89
5	B	506	CA8	N3-C2-N1	-12.43	119.38	128.89
2	A	501	FAD	N3A-C2A-N1A	-11.23	120.30	128.89
2	B	501	FAD	N3A-C2A-N1A	-11.18	120.34	128.89
5	A	508	CA8	C5'-C4'-C3'	-7.71	86.54	114.31
5	B	506	CA8	OCA-SBW-OBZ	-7.49	108.51	118.40
5	B	506	CA8	C5'-C4'-C3'	-5.83	93.32	114.31
5	A	508	CA8	OCA-SBW-OBZ	-5.52	111.11	118.40
5	A	508	CA8	CBQ-CBC-CBB	-4.42	102.78	108.50
2	B	501	FAD	P-O3P-PA	-4.18	120.99	132.73
2	A	501	FAD	P-O3P-PA	-3.93	121.69	132.73
5	A	508	CA8	OBZ-SBW-CBV	-3.73	103.37	108.08
2	B	501	FAD	C4X-C4-N3	-3.47	118.84	123.59
5	A	508	CA8	C4-C5-N7	-3.34	106.41	109.48
5	B	506	CA8	OBZ-SBW-CBX	-3.24	104.83	108.38
5	A	508	CA8	CBH-CBI-NBJ	-3.14	111.01	116.46
2	A	501	FAD	C4X-C4-N3	-3.12	119.33	123.59
5	B	506	CA8	OCA-SBW-CBX	-2.65	105.48	108.38
5	B	506	CA8	C4-C5-N7	-2.65	107.04	109.48
2	B	501	FAD	C1B-N9A-C4A	-2.63	122.97	126.94
5	B	506	CA8	C2'-C1'-N9	-2.42	110.60	114.29
5	A	508	CA8	OAY-PAX-O5'	-2.36	96.67	102.94
2	A	501	FAD	C1B-N9A-C4A	-2.34	123.40	126.94
5	B	506	CA8	CBH-CBI-NBJ	-2.34	112.39	116.46
5	B	506	CA8	OAS-PAR-OAU	-2.20	103.49	110.58
5	B	506	CA8	O2'-C2'-C3'	-2.10	105.09	111.16
5	A	508	CA8	CBK-NBJ-CBI	2.02	126.76	122.79
5	A	508	CA8	OCA-SBW-CBV	2.08	110.72	108.08
5	A	508	CA8	C2'-C1'-N9	2.13	117.54	114.29
5	A	508	CA8	OBZ-SBW-CBX	2.14	110.72	108.38
5	A	508	CA8	OBT-CBI-NBJ	2.15	127.21	122.94
5	B	506	CA8	C1'-N9-C4	2.20	130.25	126.94
5	B	506	CA8	OBT-CBI-NBJ	2.31	127.53	122.94
5	A	508	CA8	C4'-O4'-C1'	2.32	112.26	109.72
2	B	501	FAD	C1'-N10-C9A	2.36	121.51	118.86
5	A	508	CA8	O2'-C2'-C3'	2.36	117.98	111.16
5	B	506	CA8	CBG-CBH-CBI	2.38	116.23	112.31
5	A	508	CA8	OCA-SBW-CBX	2.51	111.13	108.38
5	B	506	CA8	CCC-CCB-CBX	2.57	121.74	118.95
2	A	501	FAD	C4X-N5-C5X	2.71	119.89	116.76
5	A	508	CA8	CBQ-CBC-CBD	3.04	114.89	109.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	508	CA8	CBV-SBW-CBX	3.07	109.53	105.09
5	A	508	CA8	OAT-PAR-OAS	3.11	119.22	107.38
2	A	501	FAD	C5X-C9A-N10	3.23	120.07	117.62
5	B	506	CA8	OAY-PAX-O5'	3.24	111.53	102.94
5	B	506	CA8	C3'-C2'-C1'	3.29	107.87	99.98
5	A	508	CA8	O5'-C5'-C4'	3.31	121.31	109.12
5	A	508	CA8	OAY-PAZ-OBA	3.53	112.31	102.94
5	B	506	CA8	CBK-NBJ-CBI	3.63	129.93	122.79
2	B	501	FAD	C5X-C9A-N10	3.72	120.44	117.62
5	B	506	CA8	OBA-CBB-CBC	3.72	116.53	110.55
5	B	506	CA8	O3'-C3'-C2'	3.96	126.91	111.51
5	A	508	CA8	PAR-O3'-C3'	4.32	131.93	121.56
5	A	508	CA8	C3'-C2'-C1'	4.40	110.53	99.98
5	A	508	CA8	CBR-CBC-CBQ	5.21	119.73	109.28
5	A	508	CA8	OBA-CBB-CBC	5.56	119.49	110.55
2	A	501	FAD	C4-N3-C2	5.65	120.13	115.25
2	B	501	FAD	C4-N3-C2	5.99	120.42	115.25
5	B	506	CA8	CBV-SBW-CBX	6.66	114.74	105.09
5	B	506	CA8	OBZ-SBW-CBV	6.92	116.84	108.08
5	B	506	CA8	CBL-CBU-CBV	7.64	135.24	112.83
5	A	508	CA8	O3'-PAR-OAU	8.73	128.90	107.11
5	B	506	CA8	PAR-O3'-C3'	11.19	148.40	121.56
5	B	506	CA8	O3'-PAR-OAU	11.66	136.22	107.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	FAD	1	0
5	A	508	CA8	8	0
2	B	501	FAD	4	0
5	B	506	CA8	9	0

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	437/437 (100%)	-0.52	5 (1%) 82 80	15, 23, 43, 68	0
1	B	437/437 (100%)	-0.51	4 (0%) 85 84	16, 26, 44, 75	0
All	All	874/874 (100%)	-0.52	9 (1%) 84 82	15, 24, 43, 75	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	123	GLU	3.8
1	B	123	GLU	3.7
1	B	438	LYS	3.6
1	A	55	ASP	3.5
1	A	255	ARG	3.2
1	B	363	PRO	3.1
1	A	58	TYR	2.3
1	B	360	ASN	2.2
1	A	57	ARG	2.2

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	CA8	B	506	58/58	0.71	0.29	10.53	23,45,70,85	58
5	CA8	A	508	58/58	0.88	0.20	3.75	17,36,54,66	58
2	FAD	A	501	53/53	0.97	0.10	0.09	11,17,20,23	0
2	FAD	B	501	53/53	0.97	0.08	-0.12	14,19,25,28	0
4	CL	A	504	1/1	0.85	0.05	-3.14	69,69,69,69	0
4	CL	A	507	1/1	0.26	0.20	-	86,86,86,86	0
3	MG	B	502	1/1	0.86	0.15	-	52,52,52,52	0
3	MG	A	502	1/1	0.94	0.06	-	37,37,37,37	0
4	CL	A	506	1/1	0.94	0.24	-	61,61,61,61	0
4	CL	B	504	1/1	0.95	0.04	-	76,76,76,76	0
3	MG	A	503	1/1	0.34	0.17	-	71,71,71,71	0
4	CL	B	505	1/1	0.82	0.14	-	79,79,79,79	0
4	CL	B	503	1/1	0.93	0.09	-	68,68,68,68	0
4	CL	A	505	1/1	0.87	0.12	-	63,63,63,63	0

## 6.5 Other polymers

There are no such residues in this entry.