



# Full wwPDB X-ray Structure Validation Report i

Feb 27, 2014 – 01:47 PM GMT

PDB ID : 2CLM  
Title : TRYPTOPHAN SYNTHASE (EXTERNAL ALDIMINE STATE) IN COMPLEX WITH N-(4'-TRIFLUOROMETHOXYBENZOYL)-2-AMINO-1-ETHYLPHOSPHATE(F6F)  
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Deposited on : 2006-04-28  
Resolution : 1.51 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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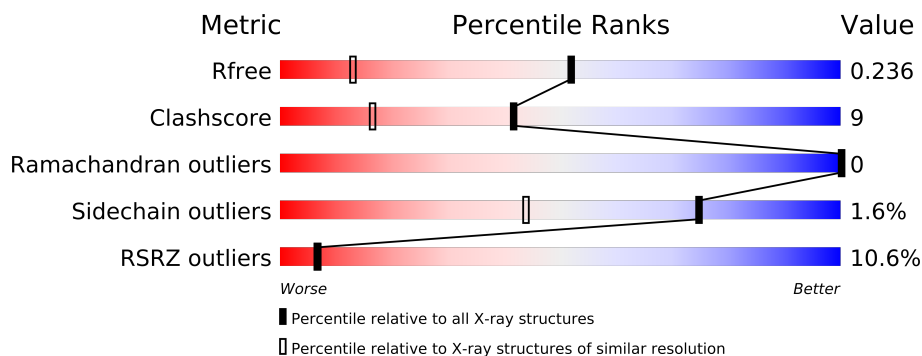
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

# 1 Overall quality at a glance



The reported resolution of this entry is 1.51 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))
$R_{free}$	66092	1879 (1.54-1.50)
Clashscore	79885	2184 (1.54-1.50)
Ramachandran outliers	78287	2123 (1.54-1.50)
Sidechain outliers	78261	2121 (1.54-1.50)
RSRZ outliers	66119	1880 (1.54-1.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	268	
2	B	396	

## 2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 5557 atoms, of which 0 are hydrogen and 0 are deuterium.

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 TRYPTOPHAN SYNTHASE ALPHA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	0	1
			1902	1211	329	355	7			

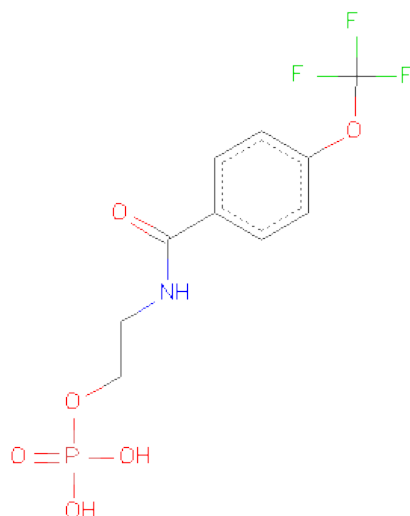
- Molecule 2 is a protein called TRYPTOPHAN SYNTHASE BETA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	393	Total	C	N	O	S	0	0	1
			2968	1865	522	562	19			

There is a discrepancy between the modelled and reference sequences:

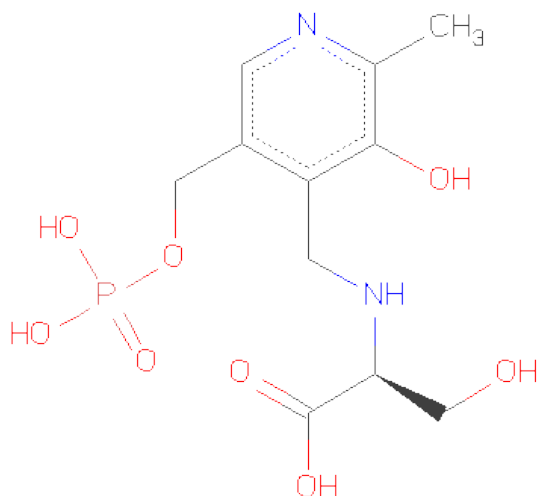
Chain	Residue	Modelled	Actual	Comment	Reference
B	34	ARG	SER	CONFLICT	UNP P0A2K1

- Molecule 3 is 2-[4-(TRIFLUOROMETHOXY)BENZOYL]AMINO}ETHYLDIHYDROGEN PHOSPHATE (three-letter code: F6F) (formula: C<sub>10</sub>H<sub>11</sub>F<sub>3</sub>NO<sub>6</sub>P).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	P	0	0
			21	10	3	1	6	1		

- Molecule 4 is [3-HYDROXY-2-METHYL-5-PHOSPHONOXYMETHYL-PYRIDIN-4-YL METHYL]-SERINE (three-letter code: PLS) (formula: C<sub>11</sub>H<sub>17</sub>N<sub>2</sub>O<sub>8</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	0	0
			22	11	2	8	1		

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Na	0	0
			1	1		

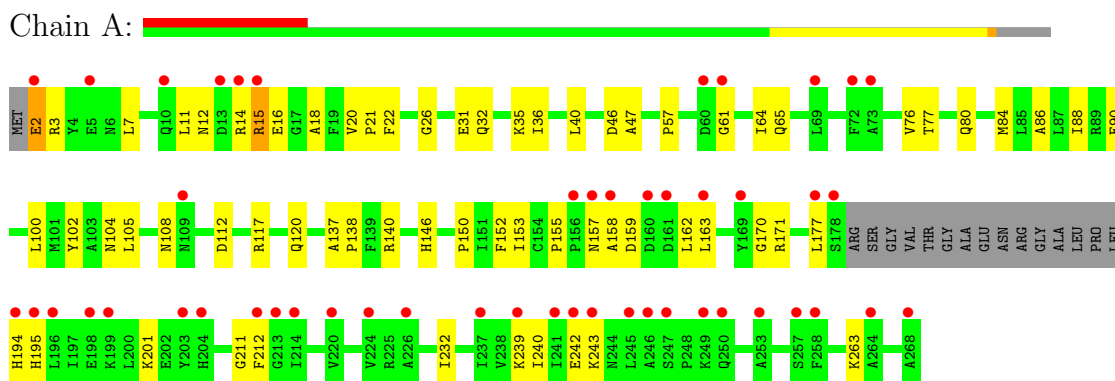
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	201	Total	O	0	0
			201	201		
6	B	442	Total	O	0	0
			442	442		

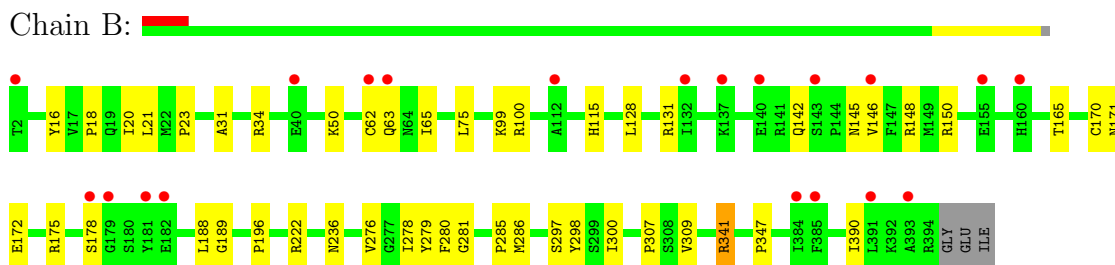
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: TRYPTOPHAN SYNTHASE ALPHA CHAIN



#### • Molecule 2: TRYPTOPHAN SYNTHASE BETA CHAIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	182.51Å 59.54Å 67.71Å 90.00° 94.84° 90.00°	Depositor
Resolution (Å)	19.93 – 1.51 19.92 – 1.51	Depositor EDS
% Data completeness (in resolution range)	89.7 (19.93-1.51) 90.3 (19.92-1.51)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.42 (at 1.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.218 , 0.239 0.217 , 0.236	Depositor DCC
$R_{free}$ test set	5137 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	15.9	Xtriage
Anisotropy	0.547	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 38.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 102735 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5557	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.16% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, PLS, F6F

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	0.29	0/1940	0.54	0/2636
2	B	0.31	0/3026	0.61	0/4090
All	All	0.30	0/4966	0.58	0/6726

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	298	TYR	Sidechain

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1902	0	1903	50	0
2	B	2968	0	2940	42	0
3	A	21	0	9	1	0
4	B	22	0	13	0	0
5	B	1	0	0	0	0
6	A	201	0	0	4	0
6	B	442	0	0	6	1
All	All	5557	0	4865	89	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (89) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:137:ALA:HB3	1:A:138:PRO:HD3	1.70	0.73
1:A:117:ARG:HH11	1:A:120:GLN:NE2	1.87	0.73
1:A:150:PRO:HD2	1:A:171:ARG:HB2	1.73	0.70
1:A:36:ILE:O	1:A:40:LEU:HD13	1.94	0.68
2:B:99:LYS:HD3	2:B:128:LEU:HD21	1.73	0.68
2:B:145:ASN:HD22	2:B:148:ARG:HE	1.44	0.65
2:B:34:ARG:HG3	2:B:100:ARG:HE	1.60	0.65
1:A:140:ARG:HD3	1:A:152:PHE:HZ	1.61	0.64
1:A:159:ASP:O	1:A:163:LEU:HD13	1.99	0.63
2:B:172:GLU:HA	2:B:175:ARG:HD2	1.80	0.63
1:A:240:ILE:HA	1:A:243:LYS:HE3	1.82	0.62
2:B:63:GLN:HG2	6:B:2403:HOH:O	1.98	0.62
2:B:171:ASN:O	2:B:175:ARG:HG2	2.00	0.61
1:A:117:ARG:NH1	1:A:120:GLN:NE2	2.48	0.61
2:B:341:ARG:HH11	2:B:341:ARG:HG3	1.66	0.61
1:A:194:HIS:CD2	1:A:195:HIS:H	2.21	0.59
2:B:150:ARG:NH1	6:B:2244:HOH:O	2.33	0.58
1:A:155:PRO:HG2	1:A:158:ALA:HB2	1.87	0.56
1:A:26:GLY:HA3	1:A:76:VAL:HG21	1.89	0.55
1:A:211:GLY:O	1:A:212:PHE:HB2	2.06	0.55
2:B:62:CYS:SG	2:B:75:LEU:HG	2.47	0.55
2:B:172:GLU:HA	2:B:175:ARG:CG	2.37	0.54
1:A:32:GLN:O	1:A:36:ILE:HG13	2.08	0.54
2:B:99:LYS:CD	2:B:128:LEU:HD21	2.39	0.53
1:A:140:ARG:HD3	1:A:152:PHE:CZ	2.44	0.53
2:B:172:GLU:HA	2:B:175:ARG:CD	2.39	0.53
1:A:155:PRO:HA	1:A:177:LEU:HD12	1.90	0.52
2:B:297:SER:OG	2:B:307:PRO:HA	2.08	0.52
1:A:2:GLU:HG2	1:A:3:ARG:N	2.25	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:34:ARG:HG3	2:B:100:ARG:NE	2.26	0.51
2:B:145:ASN:ND2	2:B:148:ARG:HH21	2.08	0.51
1:A:57:PRO:HA	1:A:102:TYR:CZ	2.45	0.51
1:A:170:GLY:O	1:A:171:ARG:HD2	2.11	0.50
2:B:222:ARG:NH2	6:B:2297:HOH:O	2.45	0.49
1:A:100:LEU:C	1:A:100:LEU:HD13	2.33	0.48
2:B:62:CYS:CB	2:B:65:ILE:HD11	2.43	0.48
1:A:46:ASP:OD2	1:A:263:LYS:HE2	2.13	0.48
1:A:117:ARG:O	1:A:120:GLN:HG2	2.13	0.48
2:B:62:CYS:HB3	2:B:65:ILE:HD11	1.96	0.48
1:A:12:ASN:O	1:A:15:ARG:HD2	2.14	0.47
2:B:341:ARG:NH1	2:B:341:ARG:HG3	2.28	0.47
1:A:158:ALA:CB	1:A:162:LEU:HD23	2.45	0.46
2:B:142:GLN:O	2:B:146:VAL:HG23	2.15	0.46
1:A:15:ARG:N	1:A:15:ARG:HD2	2.31	0.46
2:B:21:LEU:HD21	2:B:178:SER:HA	1.98	0.46
1:A:20:VAL:HG22	1:A:47:ALA:HB3	1.97	0.46
2:B:34:ARG:CG	2:B:100:ARG:HE	2.25	0.46
2:B:31:ALA:O	2:B:34:ARG:HG2	2.16	0.46
2:B:131:ARG:NH1	6:B:2218:HOH:O	2.49	0.46
1:A:22:PHE:CD1	1:A:22:PHE:C	2.89	0.46
2:B:16:TYR:O	2:B:281:GLY:HA2	2.17	0.45
1:A:158:ALA:HB1	1:A:162:LEU:HD23	1.99	0.45
1:A:20:VAL:HB	1:A:232:ILE:HG12	1.99	0.45
1:A:11:LEU:CD1	1:A:18:ALA:HB2	2.46	0.45
1:A:16:GLU:HB3	6:A:2023:HOH:O	2.17	0.45
2:B:31:ALA:HA	2:B:34:ARG:HG2	1.99	0.44
2:B:150:ARG:HH11	2:B:150:ARG:HG2	1.82	0.44
2:B:34:ARG:CD	2:B:100:ARG:HE	2.30	0.43
1:A:61:GLY:O	1:A:65:GLN:HG3	2.18	0.43
1:A:108:ASN:HA	6:A:2099:HOH:O	2.18	0.43
1:A:77:THR:OG1	1:A:80:GLN:HG3	2.18	0.43
1:A:11:LEU:HD23	1:A:14:ARG:HH21	1.83	0.43
1:A:157:ASN:OD1	2:B:20:ILE:HD12	2.19	0.43
2:B:170:CYS:HB3	2:B:280:PHE:CE1	2.53	0.43
1:A:153:ILE:HG23	1:A:177:LEU:HG	1.99	0.43
1:A:157:ASN:ND2	2:B:23:PRO:HG2	2.33	0.43
1:A:7:LEU:O	1:A:11:LEU:HG	2.18	0.43
1:A:239:LYS:HA	1:A:242:GLU:HG2	1.99	0.43
2:B:21:LEU:CD2	2:B:178:SER:HA	2.48	0.42
2:B:276:VAL:HA	2:B:285:PRO:HA	2.01	0.42
2:B:18:PRO:HB2	2:B:20:ILE:HG22	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:86:ALA:O	1:A:90:GLU:HG3	2.19	0.42
1:A:194:HIS:HD2	1:A:195:HIS:H	1.67	0.42
2:B:300:ILE:HD11	2:B:390:ILE:CD1	2.50	0.42
2:B:50:LYS:HG3	6:B:2122:HOH:O	2.18	0.42
2:B:279:TYR:HB3	2:B:286:MET:SD	2.60	0.42
2:B:165:THR:CG2	6:B:2354:HOH:O	2.68	0.42
1:A:112:ASP:OD1	1:A:146:HIS:HE1	2.03	0.42
1:A:31:GLU:O	1:A:35:LYS:HG3	2.20	0.41
2:B:115:HIS:CE1	2:B:189:GLY:HA2	2.54	0.41
1:A:117:ARG:NH1	1:A:120:GLN:HE21	2.15	0.41
1:A:21:PRO:HD2	1:A:47:ALA:O	2.21	0.41
1:A:31:GLU:HG2	6:A:2035:HOH:O	2.19	0.41
1:A:64:ILE:CD1	3:A:1268:F6F:H151	2.51	0.41
2:B:285:PRO:C	2:B:286:MET:HG2	2.42	0.40
1:A:84:MET:O	1:A:88:ILE:HG13	2.21	0.40
1:A:104:ASN:HB2	2:B:278:ILE:O	2.21	0.40
2:B:285:PRO:HG2	2:B:309:VAL:HG12	2.04	0.40
1:A:201:LYS:HA	6:A:2166:HOH:O	2.20	0.40

All (1) 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	Distance(Å)	Clash(Å)
6:B:2125:HOH:O	6:B:2125:HOH:O[2.655]	1.90	0.30

## 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	248/268 (92%)	241 (97%)	7 (3%)	0	100	100
2	B	391/396 (99%)	381 (97%)	10 (3%)	0	100	100
All	All	639/664 (96%)	622 (97%)	17 (3%)	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	197/208 (95%)	194 (98%)	3 (2%)	76	47
2	B	306/310 (99%)	301 (98%)	5 (2%)	75	44
All	All	503/518 (97%)	495 (98%)	8 (2%)	75	44

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

Mol	Chain	Res	Type
1	A	2	GLU
1	A	15	ARG
1	A	105	LEU
2	B	188	LEU
2	B	196	PRO
2	B	236	ASN
2	B	341	ARG
2	B	347	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	120	GLN
1	A	194	HIS
1	A	210	GLN
1	A	250	GLN
2	B	26	ASN
2	B	64	ASN
2	B	142	GLN
2	B	145	ASN
2	B	236	ASN
2	B	317	ASN
2	B	375	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\# Z  > 2$
3	F6F	A	1268	-	21,21,21	1.93	6 (28%)	30,30,30	1.89	5 (16%)
4	PLS	B	1394	-	22,22,22	4.33	13 (59%)	31,31,31	3.13	13 (41%)

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	F6F	A	1268	-	-	0/17/17/17	0/1/1/1
4	PLS	B	1394	-	-	0/17/17/17	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1394	PLS	C3-C2	13.20	1.50	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1394	PLS	C3-C4	8.03	1.54	1.40
4	B	1394	PLS	O-C	7.10	1.46	1.22
4	B	1394	PLS	C4A-N	-4.55	1.33	1.46
4	B	1394	PLS	C5-C4	4.01	1.46	1.40
4	B	1394	PLS	C2-N1	3.96	1.41	1.33
4	B	1394	PLS	P-O3P	-3.81	1.40	1.54
3	A	1268	F6F	C5-C4	3.66	1.45	1.39
4	B	1394	PLS	C6-C5	3.44	1.45	1.37
4	B	1394	PLS	CA-C	3.42	1.62	1.52
3	A	1268	F6F	C3-C4	3.38	1.45	1.39
3	A	1268	F6F	C6-C5	3.31	1.45	1.38
3	A	1268	F6F	C6-C1	3.23	1.45	1.38
3	A	1268	F6F	C2-C1	2.90	1.44	1.38
3	A	1268	F6F	C3-C2	2.88	1.44	1.38
4	B	1394	PLS	P-O2P	-2.86	1.44	1.54
4	B	1394	PLS	C6-N1	2.70	1.40	1.34
4	B	1394	PLS	CA-N	2.27	1.50	1.47
4	B	1394	PLS	P-O4P	-2.15	1.52	1.60

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1394	PLS	C4-C4A-N	8.92	136.56	111.67
3	A	1268	F6F	C4-C12-N13	-8.46	97.18	117.17
4	B	1394	PLS	O3P-P-O4P	7.05	126.10	106.65
4	B	1394	PLS	C4A-N-CA	6.64	128.24	114.34
4	B	1394	PLS	CB-CA-N	5.41	115.64	108.55
4	B	1394	PLS	O3-C3-C2	4.84	126.21	117.61
4	B	1394	PLS	O4P-C5A-C5	3.65	116.69	109.26
4	B	1394	PLS	C6-N1-C2	3.22	126.17	119.28
3	A	1268	F6F	O21-P18-O20	3.11	120.62	110.44
4	B	1394	PLS	OXT-C-O	-2.92	117.47	124.07
4	B	1394	PLS	C4A-C4-C5	-2.90	117.04	119.70
3	A	1268	F6F	O14-C12-N13	-2.52	117.04	122.50
4	B	1394	PLS	C5-C6-N1	-2.44	119.46	123.86
3	A	1268	F6F	O14-C12-C4	-2.33	116.92	121.01
3	A	1268	F6F	P18-O17-C16	-2.32	111.49	118.19
4	B	1394	PLS	OXT-C-CA	2.20	122.17	113.90
4	B	1394	PLS	C4A-C4-C3	2.13	123.01	120.31
4	B	1394	PLS	C3-C2-N1	-2.02	117.89	120.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

### 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, 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	252/268 (94%)	1.13	49 (19%) 2 1	11, 27, 37, 44	0
2	B	393/396 (99%)	0.36	20 (5%) 27 29	9, 15, 32, 55	0
All	All	645/664 (97%)	0.66	69 (10%) 7 6	9, 20, 35, 55	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	195	HIS	7.9
1	A	69	LEU	6.0
2	B	393	ALA	5.2
1	A	245	LEU	5.2
2	B	2	THR	4.9
1	A	13	ASP	4.7
2	B	140	GLU	4.6
1	A	15	ARG	4.6
1	A	157	ASN	4.4
1	A	178	SER	4.2
2	B	385	PHE	4.1
1	A	177	LEU	4.1
1	A	249	LYS	3.9
1	A	194	HIS	3.8
1	A	243	LYS	3.8
1	A	204	HIS	3.8
1	A	158	ALA	3.8
2	B	181	TYR	3.7
1	A	198	GLU	3.7
2	B	63	GLN	3.6
1	A	14	ARG	3.5
2	B	182	GLU	3.5
1	A	247	SER	3.3
1	A	72	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	220	VAL	3.2
1	A	241	ILE	3.2
2	B	62	CYS	3.1
1	A	73	ALA	3.1
1	A	212	PHE	3.1
1	A	213	GLY	3.0
1	A	250	GLN	3.0
1	A	169	TYR	3.0
1	A	199	LYS	3.0
1	A	264	ALA	2.8
1	A	10	GLN	2.8
1	A	246	ALA	2.7
1	A	224	VAL	2.7
1	A	258	PHE	2.6
1	A	163	LEU	2.6
1	A	196	LEU	2.6
2	B	391	LEU	2.6
1	A	60	ASP	2.5
1	A	203	TYR	2.5
1	A	109	ASN	2.5
1	A	61	GLY	2.4
2	B	160	HIS	2.4
1	A	2	GLU	2.4
1	A	161	ASP	2.3
1	A	242	GLU	2.3
2	B	178	SER	2.3
2	B	179	GLY	2.3
2	B	137	LYS	2.2
1	A	237	ILE	2.2
2	B	384	ILE	2.2
1	A	226	ALA	2.2
1	A	239	LYS	2.2
2	B	143	SER	2.2
1	A	257	SER	2.2
1	A	268	ALA	2.2
2	B	146	VAL	2.1
1	A	214	ILE	2.1
1	A	5	GLU	2.1
2	B	132	ILE	2.1
2	B	155	GLU	2.1
2	B	112	ALA	2.1
1	A	160	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	253	ALA	2.0
2	B	40	GLU	2.1
1	A	156	PRO	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	F6F	A	1268	21/21	0.16	0.55	29,31,35,37	0
4	PLS	B	1394	22/22	0.10	0.32	10,16,23,24	0
5	NA	B	1395	1/1	0.08	-0.58	15,15,15,15	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.