



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 07:10 AM GMT

PDB ID : 3Q7F  
Title : Cryptococcus neoformans protein farnesyltransferase in complex with FPP  
and ethylenediamine inhibitor 1  
Authors : Hast, M.A.; Beese, L.S.  
Deposited on : 2011-01-04  
Resolution : 2.20 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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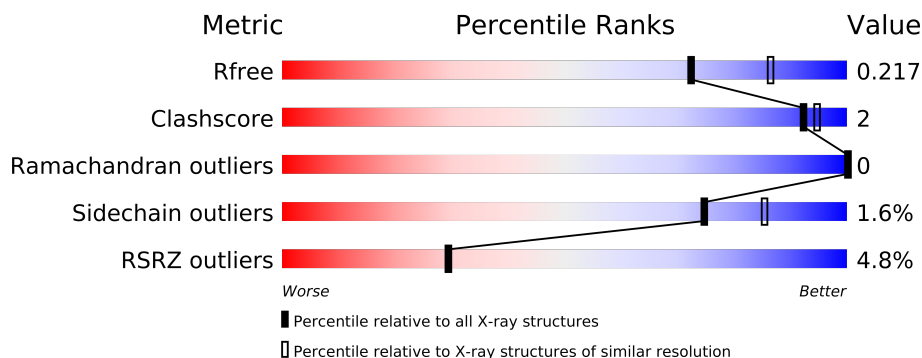
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.20 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	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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	349	
2	B	520	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	SUC	A	3010	-	X
3	SUC	B	3010	-	X
5	3CX	B	524	-	X
8	SO4	B	526	-	X

## 2 Entry composition i

There are 9 unique types of molecules in this entry. The entry contains 7003 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.

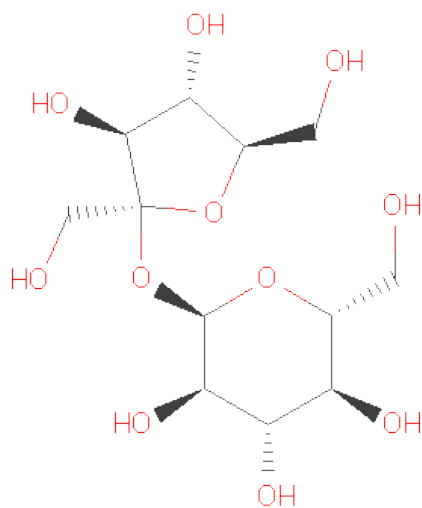
- Molecule 1 is a protein called Farnesyltransferase alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2606	1680	444	471	11			

- Molecule 2 is a protein called Farnesyltransferase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	485	Total	C	N	O	S	0	1	0
			3730	2366	649	700	15			

- Molecule 3 is SUGAR (SUCROSE) (three-letter code: SUC) (formula:  $C_{12}H_{22}O_{11}$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			23	12	11		
3	B	1	Total	C	O	0	0
			23	12	11		

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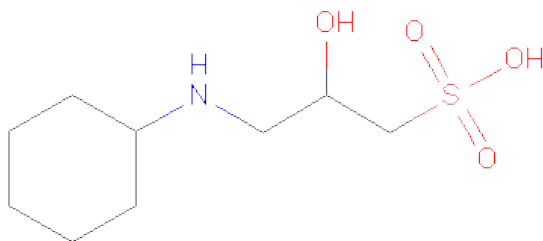
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			23	12	11		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

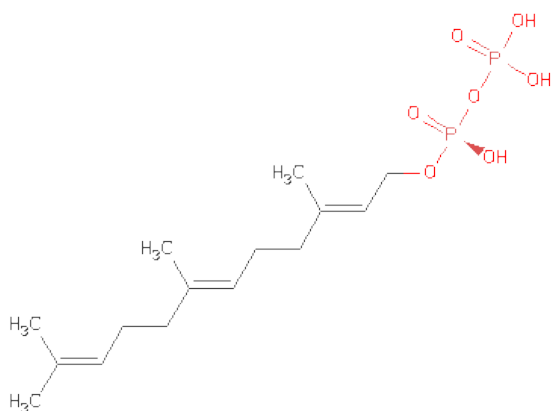
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		

- Molecule 5 is (2S)-3-(CYCLOHEXYLAMINO)-2-HYDROXYPROPANE-1-SULFONIC ACID (three-letter code: 3CX) (formula: C<sub>9</sub>H<sub>19</sub>NO<sub>4</sub>S).



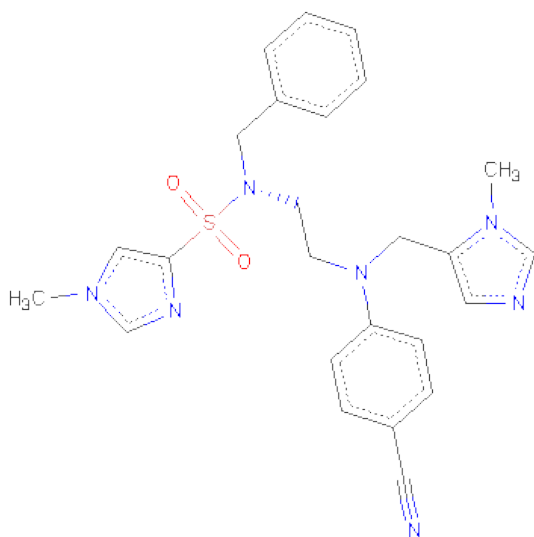
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			15	9	1	4	1		
5	B	1	Total	C	N	O	S	0	0
			15	9	1	4	1		
5	B	1	Total	C	N	O	S	0	0
			15	9	1	4	1		

- Molecule 6 is FARNESYL DIPHOSPHATE (three-letter code: FPP) (formula: C<sub>15</sub>H<sub>28</sub>O<sub>7</sub>P<sub>2</sub>).



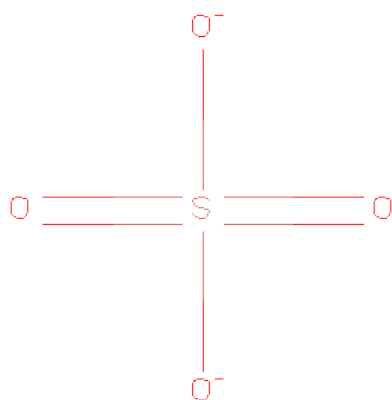
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	O	P	0	0
			24	15	7	2		

- Molecule 7 is N-BENZYL-N-(2-{(4-CYANOPHENYL)[(1-METHYL-1H-IMIDAZOL-5-YL)METHYL]AMINO}ETHYL)-1-METHYL-1H-IMIDAZOLE-4-SULFONAMIDE (three-letter code: ED2) (formula: C<sub>25</sub>H<sub>27</sub>N<sub>7</sub>O<sub>2</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	S	0	0
			35	25	7	2	1		

- Molecule 8 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

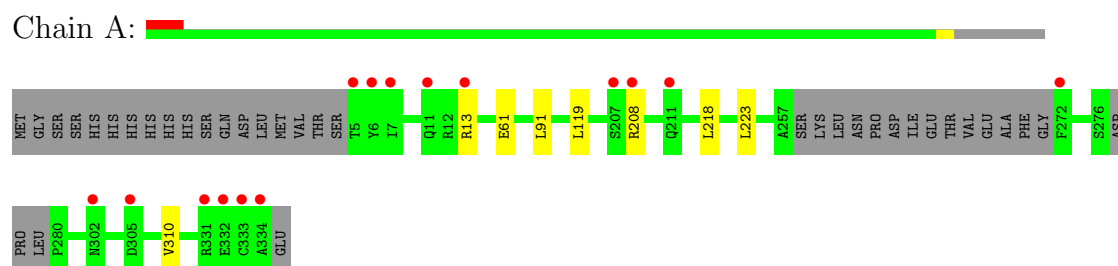
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	171	Total	O	0	0
			171	171		
9	B	317	Total	O	0	0
			317	317		

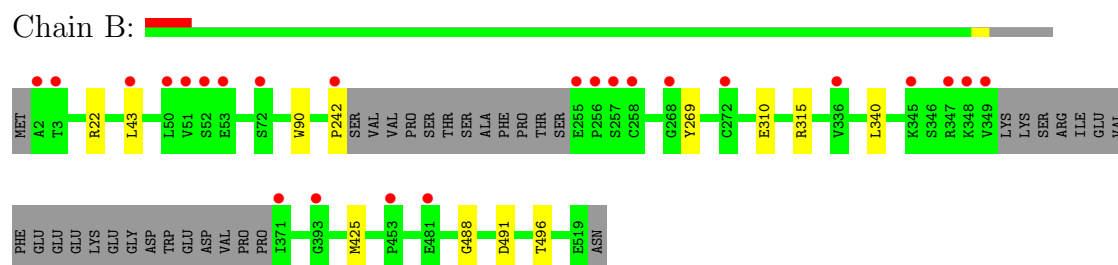
### 3 Residue-property plots

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: Farnesyltransferase alpha subunit



- Molecule 2: Farnesyltransferase beta subunit



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	143.58Å 143.58Å 130.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.94 – 2.20 44.94 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (44.94-2.20) 99.1 (44.94-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.55 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.191 , 0.213 0.196 , 0.217	Depositor DCC
$R_{free}$ test set	3447 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.7	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 32.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 69057 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7003	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% 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: ZN, SUC, ED2, SO4, FPP, 3CX

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	0.35	0/2687	0.47	0/3657
2	B	0.37	0/3827	0.50	1/5198 (0.0%)
All	All	0.36	0/6514	0.48	1/8855 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	242	PRO	CA-N-CD	-8.53	99.56	111.50

There are no chirality outliers.

There are no planarity outliers.

### 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	2606	0	0	0	0
2	B	3730	0	2	7	0
3	A	23	0	22	0	0
3	B	46	0	44	2	0
4	B	1	0	0	0	0
5	B	45	0	48	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	24	0	25	3	0
7	B	35	0	27	5	0
8	B	5	0	0	0	0
9	A	171	0	0	0	0
9	B	317	0	0	1	0
All	All	7003	0	168	12	0

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 2.

All (12) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
6:B:525:FPP:H102	7:B:1003:ED2:NAY	2.15	0.60
2:B:90:TRP:CD1	7:B:1003:ED2:HAG	2.37	0.60
6:B:525:FPP:PA	7:B:1003:ED2:HAAB	2.55	0.47
2:B:491:ASP:OD2	5:B:522:3CX:NAL	2.47	0.46
2:B:269:TYR:CE2	6:B:525:FPP:H41	2.53	0.43
2:B:315:ARG:NE	9:B:1828:HOH:O	2.52	0.43
7:B:1003:ED2:HAVA	7:B:1003:ED2:HATA	1.70	0.42
3:B:3010:SUC:H1'2	3:B:3010:SUC:H1	1.76	0.42
2:B:425:MET:SD	2:B:496:THR:CG2	3.08	0.42
2:B:488:GLY:N	3:B:1002:SUC:HO4	2.18	0.42
7:B:1003:ED2:CAN	7:B:1003:ED2:HAU	2.50	0.41
2:B:310:GLU:OE1	2:B:315:ARG:NH2	2.54	0.41

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	307/349 (88%)	298 (97%)	9 (3%)	0	100	100
2	B	480/520 (92%)	472 (98%)	8 (2%)	0	100	100
All	All	787/869 (91%)	770 (98%)	17 (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	283/316 (90%)	275 (97%)	8 (3%)	56	67
2	B	404/436 (93%)	401 (99%)	3 (1%)	91	96
All	All	687/752 (91%)	676 (98%)	11 (2%)	75	85

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

Mol	Chain	Res	Type
1	A	13	ARG
1	A	61	GLU
1	A	91	LEU
1	A	119	LEU
1	A	208	ARG
1	A	218	LEU
1	A	223	LEU
1	A	310	VAL
2	B	22	ARG
2	B	43	LEU
2	B	340	LEU

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 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 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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	SUC	A	3010	-	24,24,24	1.04	1 (4%)	36,36,36	0.92	1 (2%)
3	SUC	B	1002	-	24,24,24	1.03	1 (4%)	36,36,36	1.01	2 (5%)
7	ED2	B	1003	4	38,38,38	1.86	7 (18%)	51,53,53	1.58	7 (13%)
3	SUC	B	3010	-	24,24,24	1.05	1 (4%)	36,36,36	0.98	1 (2%)
5	3CX	B	522	-	15,15,15	1.83	1 (6%)	20,20,20	1.81	6 (30%)
5	3CX	B	523	-	15,15,15	1.87	1 (6%)	20,20,20	1.85	5 (25%)
5	3CX	B	524	-	15,15,15	1.90	1 (6%)	20,20,20	1.78	4 (20%)
6	FPP	B	525	-	23,23,23	1.18	2 (8%)	31,31,31	1.29	6 (19%)
8	SO4	B	526	-	4,4,4	0.24	0	6,6,6	0.08	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	SUC	A	3010	-	4/4/9/9	0/12/51/51	1/2/2/2
3	SUC	B	1002	-	4/4/9/9	0/12/51/51	0/2/2/2
7	ED2	B	1003	4	-	0/25/31/31	0/4/4/4
3	SUC	B	3010	-	4/4/9/9	0/12/51/51	0/2/2/2
5	3CX	B	522	-	1/1/3/4	0/10/18/18	0/1/1/1
5	3CX	B	523	-	1/1/3/4	0/10/18/18	0/1/1/1
5	3CX	B	524	-	1/1/3/4	0/10/18/18	0/1/1/1
6	FPP	B	525	-	-	1/25/25/25	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	SO4	B	526	-	-	0/0/0/0	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	524	3CX	OAC-CAM	-6.76	1.22	1.43
5	B	523	3CX	OAC-CAM	-6.75	1.22	1.43
5	B	522	3CX	OAC-CAM	-6.46	1.23	1.43
7	B	1003	ED2	CAL-CAN	5.65	1.49	1.38
7	B	1003	ED2	CAS-CBD	4.36	1.43	1.37
7	B	1003	ED2	CAM-CAZ	4.20	1.48	1.39
7	B	1003	ED2	CAR-NBG	3.22	1.41	1.36
7	B	1003	ED2	CAS-NBG	3.00	1.42	1.37
7	B	1003	ED2	CBD-SBI	-2.58	1.76	1.78
7	B	1003	ED2	CAZ-CAF	2.51	1.51	1.44
6	B	525	FPP	C7-C8	2.35	1.37	1.32
3	B	1002	SUC	O2-C2	-2.29	1.37	1.43
3	A	3010	SUC	O2-C2	-2.28	1.37	1.43
6	B	525	FPP	C2-C3	2.27	1.37	1.32
3	B	3010	SUC	O2-C2	-2.22	1.37	1.43

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	1003	ED2	OAD-SBI-OAE	-5.78	109.06	119.38
7	B	1003	ED2	OAE-SBI-NBF	4.54	111.34	106.71
5	B	524	3CX	CAJ-CAM-CAK	3.94	120.87	111.15
5	B	524	3CX	OAD-SAO-CAK	3.74	110.46	105.64
5	B	523	3CX	CAJ-CAM-CAK	3.54	119.89	111.15
5	B	522	3CX	OAC-CAM-CAJ	3.50	120.88	109.11
5	B	522	3CX	CAM-CAK-SAO	3.29	119.18	114.79
7	B	1003	ED2	CAR-NAY-CBD	3.27	110.34	104.75
5	B	522	3CX	OAC-CAM-CAK	3.21	120.88	110.19
5	B	523	3CX	CAM-CAK-SAO	3.20	119.05	114.79
5	B	523	3CX	OAC-CAM-CAJ	3.13	119.66	109.11
3	B	3010	SUC	O5-C5-C4	3.05	115.40	109.76
5	B	523	3CX	OAC-CAM-CAK	3.00	120.16	110.19
5	B	522	3CX	OAD-SAO-CAK	2.97	109.47	105.64
5	B	524	3CX	OAC-CAM-CAJ	2.92	118.96	109.11
5	B	524	3CX	OAC-CAM-CAK	2.89	119.81	110.19
7	B	1003	ED2	CAW-CBC-NBH	2.88	126.36	123.29
7	B	1003	ED2	OAE-SBI-CBD	2.82	111.71	107.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	525	FPP	C4-C3-C2	-2.80	117.97	123.52
3	A	3010	SUC	O5-C5-C4	2.75	114.86	109.76
3	B	1002	SUC	C1-O5-C5	-2.45	108.96	113.73
5	B	522	3CX	CAJ-CAM-CAK	2.39	117.04	111.15
3	B	1002	SUC	C1-C2-C3	2.36	114.59	110.00
5	B	523	3CX	OAD-SAO-CAK	2.28	108.58	105.64
6	B	525	FPP	PA-O3A-PB	-2.27	125.04	131.68
6	B	525	FPP	C1-C2-C3	-2.24	122.20	126.19
6	B	525	FPP	C9-C8-C7	-2.15	116.95	121.08
7	B	1003	ED2	CAV-NBF-SBI	-2.08	112.49	117.36
7	B	1003	ED2	CAU-NBF-SBI	-2.08	113.39	117.84
6	B	525	FPP	C10-C8-C9	2.06	118.52	115.39
6	B	525	FPP	C6-C7-C8	-2.04	123.41	127.80
5	B	522	3CX	CAM-CAJ-NAL	2.02	118.41	112.15

All (15) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	522	3CX	CAM
5	B	524	3CX	CAM
5	B	523	3CX	CAM
3	B	3010	SUC	C2
3	B	3010	SUC	C4'
3	B	3010	SUC	C3
3	B	3010	SUC	C5'
3	A	3010	SUC	C2
3	A	3010	SUC	C4'
3	A	3010	SUC	C3
3	A	3010	SUC	C5'
3	B	1002	SUC	C2
3	B	1002	SUC	C4'
3	B	1002	SUC	C3
3	B	1002	SUC	C5'

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	525	FPP	C1-C2-C3-C4

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	3010	SUC	C1-C2-C3-C4-C5-O5

## 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	313/349 (89%)	-0.03	15 (4%) 29 29	26, 41, 62, 71	0
2	B	485/520 (93%)	0.17	24 (4%) 28 28	21, 33, 49, 71	0
All	All	798/869 (91%)	0.09	39 (4%) 29 28	21, 36, 59, 71	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	349	VAL	8.6
1	A	5	THR	6.3
2	B	257	SER	5.8
2	B	255	GLU	5.7
2	B	2	ALA	5.0
1	A	333	CYS	4.8
2	B	51	VAL	4.5
2	B	256	PRO	4.2
2	B	242	PRO	4.2
2	B	371	ILE	3.9
2	B	348	LYS	3.9
1	A	334	ALA	3.8
2	B	481	GLU	3.6
2	B	347	ARG	3.1
1	A	302	ASN	3.1
1	A	207	SER	3.1
1	A	7	ILE	2.9
1	A	13	ARG	2.7
2	B	53	GLU	2.6
2	B	52	SER	2.6
1	A	211	GLN	2.5
2	B	50	LEU	2.5
1	A	6	TYR	2.5
1	A	272	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	332	GLU	2.4
2	B	393	GLY	2.4
1	A	208	ARG	2.3
2	B	336	VAL	2.3
2	B	345	LYS	2.3
2	B	272	CYS	2.2
2	B	268	GLY	2.1
2	B	43	LEU	2.1
1	A	305	ASP	2.1
2	B	3	THR	2.1
2	B	72	SER	2.1
2	B	258	CYS	2.1
2	B	453	PRO	2.1
1	A	331	ARG	2.0
1	A	11	GLN	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
8	SO4	B	526	5/5	0.23	4.43	79,79,79,79	0
3	SUC	A	3010	23/23	0.30	3.58	79,80,81,81	0
3	SUC	B	3010	23/23	0.25	2.86	72,72,73,73	0
5	3CX	B	524	15/15	0.21	2.63	64,67,73,73	0
5	3CX	B	523	15/15	0.15	1.52	46,47,48,50	0
7	ED2	B	1003	35/35	0.21	1.30	45,52,62,62	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	FPP	B	525	24/24	0.21	0.53	33,37,41,42	0
3	SUC	B	1002	23/23	0.15	0.04	44,47,47,48	0
4	ZN	B	521	1/1	0.13	-0.08	27,27,27,27	0
5	3CX	B	522	15/15	0.11	-0.24	27,28,30,31	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.