



# Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 11:50 PM GMT

PDB ID : 3SFY  
Title : Cryptococcus neoformans protein farnesyltransferase in complex with FPT-II and ethylenediamine inhibitor 2  
Authors : Hast, M.A.; Beese, L.S.  
Deposited on : 2011-06-14  
Resolution : 2.10 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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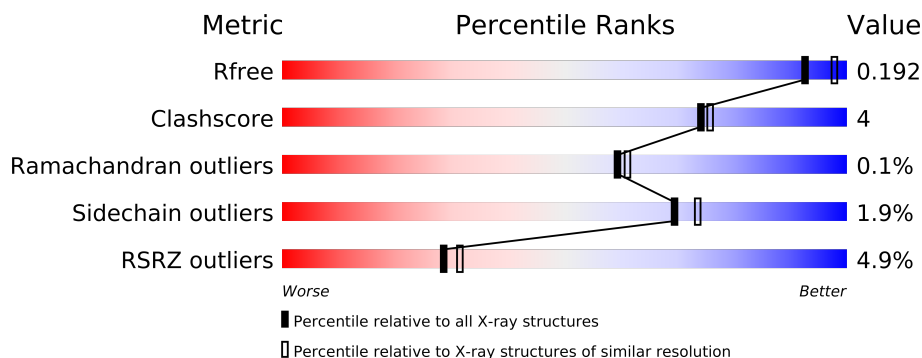
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.10 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	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

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
4	3FX	B	523	-	X
4	3FX	B	524	-	X

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7079 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 *Cryptococcus neoformans* protein farnesyltransferase alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	315	Total	C	N	O	S	0	0	0
			2621	1691	446	473	11			

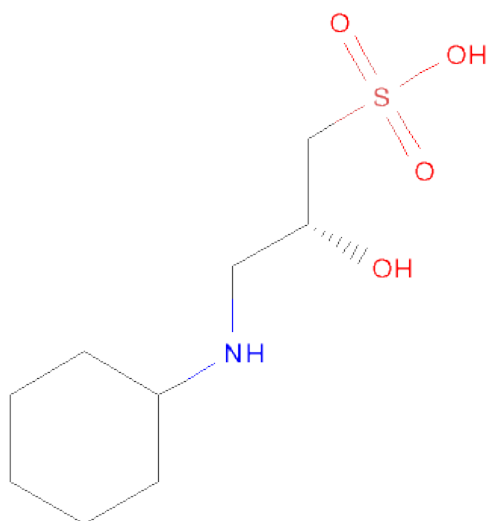
- Molecule 2 is a protein called *Cryptococcus neoformans* protein farnesyltransferase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	482	Total	C	N	O	S	0	1	0
			3702	2345	646	696	15			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

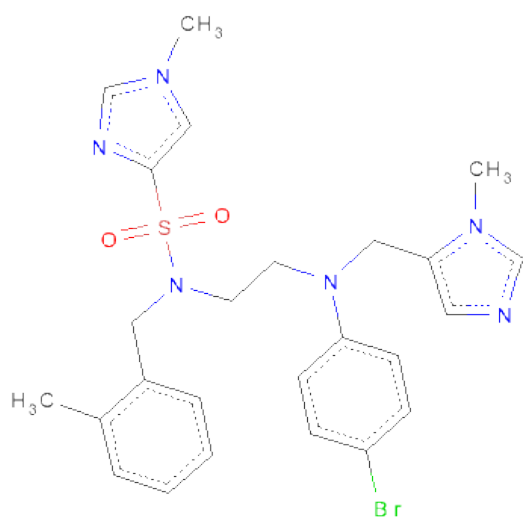
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		

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



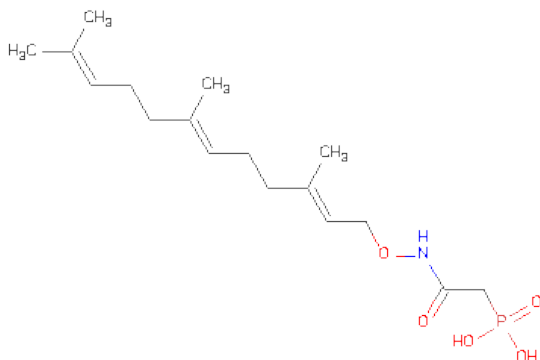
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			15	9	1	4	1		
4	B	1	Total	C	N	O	S	0	0
			15	9	1	4	1		
4	B	1	Total	C	N	O	S	0	0
			15	9	1	4	1		

- Molecule 5 is N-(2-{(4-BROMOPHENYL)[(1-METHYL-1H-IMIDAZOL-5-YL)METHYL]AMINO}ETHYL)-1-METHYL-N-(2-METHYLBENZYL)-1H-IMIDAZOLE-4-SULFONAMIDE (three-letter code: 3FY) (formula: C<sub>25</sub>H<sub>29</sub>BrN<sub>6</sub>O<sub>2</sub>S).



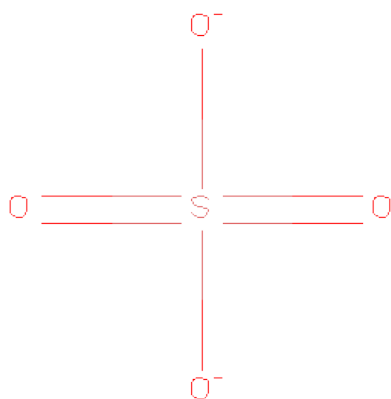
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	B	1	Total	Br	C	N	O	S	0	0
			35	1	25	6	2	1		

- Molecule 6 is [(3,7,11-TRIMETHYL-DODECA-2,6,10-TRIENYLOXYCARBAMOYL)-METHYL]-PHOSPHONICACID (three-letter code: FII) (formula: C<sub>17</sub>H<sub>30</sub>NO<sub>5</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	P	0	0
			24	17	1	5	1		

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 8 is water.

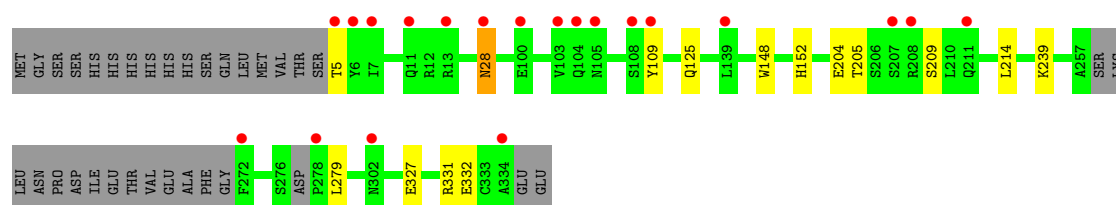
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	228	Total	O	0	0
			228	228		
8	B	418	Total	O	0	0
			418	418		

### 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.

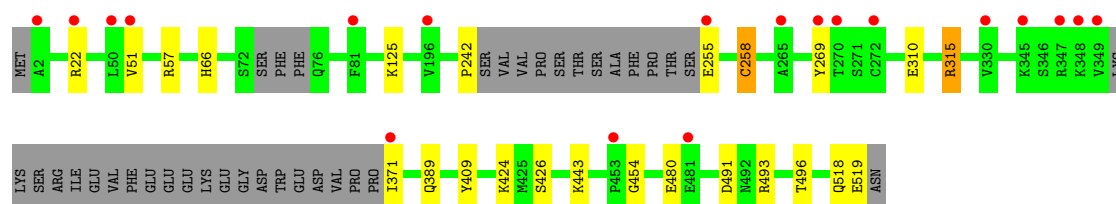
- Molecule 1: *Cryptococcus neoformans* protein farnesyltransferase alpha subunit

Chain A: 



- Molecule 2: *Cryptococcus neoformans* protein farnesyltransferase beta subunit

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	141.16Å 141.16Å 129.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.57 – 2.10 49.91 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.57-2.10) 99.9 (49.91-2.05)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.55 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.178 , 0.203 0.191 , 0.192	Depositor DCC
$R_{free}$ test set	4123 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.0	Xtriage
Anisotropy	0.468	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 29.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 82454 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7079	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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, FII, SO4, 3FX, 3FY

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	0.52	0/2703	0.56	0/3680
2	B	0.59	0/3796	0.63	2/5155 (0.0%)
All	All	0.56	0/6499	0.60	2/8835 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	315	ARG	NE-CZ-NH2	-9.65	115.47	120.30
2	B	315	ARG	NE-CZ-NH1	8.00	124.30	120.30

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	2621	0	0	8	0
2	B	3702	0	0	15	0
3	B	1	0	0	0	0
4	B	45	0	57	3	0
5	B	35	0	29	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	24	0	28	3	0
7	B	5	0	0	0	0
8	A	228	0	0	3	0
8	B	418	0	0	10	0
All	All	7079	0	114	27	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 4.

All (27) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:258:CYS:CB	8:B:2540:HOH:O	2.39	0.69
1:A:28:ASN:N	8:A:2714:HOH:O	2.30	0.64
1:A:28:ASN:CB	8:B:2794:HOH:O	2.44	0.64
4:B:522:3FX:HAJA	8:B:1900:HOH:O	2.03	0.58
2:B:255:GLU:CA	8:B:2757:HOH:O	2.55	0.54
1:A:239:LYS:NZ	8:A:1838:HOH:O	2.41	0.54
2:B:496:THR:CG2	8:B:2053:HOH:O	2.57	0.53
1:A:125:GLN:NE2	8:A:2721:HOH:O	2.41	0.52
2:B:310:GLU:OE1	2:B:315:ARG:NH2	2.43	0.51
2:B:426:SER:OG	2:B:480:GLU:OE1	2.30	0.50
2:B:125:LYS:N	4:B:524:3FX:HOAB	2.10	0.50
2:B:66:HIS:NE2	2:B:389:GLN:NE2	2.60	0.50
2:B:57:ARG:NH1	8:B:1608:HOH:O	2.47	0.47
1:A:109:TYR:CE2	5:B:525:3FY:HAR	2.51	0.46
5:B:525:3FY:HAAB	8:B:2916:HOH:O	2.16	0.46
2:B:269:TYR:CE2	6:B:526:FII:H302	2.52	0.45
2:B:409:TYR:CE2	5:B:525:3FY:HAO	2.53	0.44
1:A:148:TRP:O	1:A:152:HIS:CD2	2.70	0.44
2:B:454:GLY:N	8:B:2621:HOH:O	2.50	0.43
6:B:526:FII:H111	6:B:526:FII:H221	1.67	0.43
2:B:491:ASP:OD2	4:B:522:3FX:NAL	2.52	0.43
2:B:518:GLN:O	2:B:519:GLU:C	2.57	0.43
2:B:493:ARG:NH1	8:B:1846:HOH:O	2.52	0.42
6:B:526:FII:H181	6:B:526:FII:H232	1.80	0.42
2:B:242:PRO:C	8:B:2634:HOH:O	2.58	0.42
1:A:109:TYR:CD2	5:B:525:3FY:HAR	2.55	0.42
1:A:327:GLU:OE2	1:A:331:ARG:NE	2.54	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	309/349 (88%)	298 (96%)	11 (4%)	0	100	100
2	B	475/520 (91%)	466 (98%)	8 (2%)	1 (0%)	56	57
All	All	784/869 (90%)	764 (97%)	19 (2%)	1 (0%)	59	61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	VAL

### 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	285/316 (90%)	277 (97%)	8 (3%)	56	59
2	B	401/436 (92%)	396 (99%)	5 (1%)	82	87
All	All	686/752 (91%)	673 (98%)	13 (2%)	69	73

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	28	ASN
1	A	204	GLU
1	A	205	THR
1	A	209	SER
1	A	214	LEU
1	A	279	LEU
1	A	332	GLU

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Mol	Chain	Res	Type
2	B	22	ARG
2	B	258	CYS
2	B	371	ILE
2	B	424	LYS
2	B	443	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 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 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\# Z  > 2$
4	3FX	B	522	-	15,15,15	1.28	2 (13%)	20,20,20	1.54	2 (10%)
4	3FX	B	523	-	15,15,15	1.93	1 (6%)	20,20,20	1.23	1 (5%)
4	3FX	B	524	-	15,15,15	1.76	1 (6%)	20,20,20	0.98	1 (5%)
5	3FY	B	525	3	38,38,38	1.78	8 (21%)	52,54,54	1.82	11 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	FII	B	526	-	23,23,23	2.35	9 (39%)	29,29,29	1.48	6 (20%)
7	SO4	B	527	-	4,4,4	0.49	0	6,6,6	0.31	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
4	3FX	B	522	-	-	0/10/18/18	0/1/1/1
4	3FX	B	523	-	-	0/10/18/18	0/1/1/1
4	3FX	B	524	-	-	0/10/18/18	0/1/1/1
5	3FY	B	525	3	-	0/23/29/29	0/4/4/4
6	FII	B	526	-	-	0/24/24/24	0/0/0/0
7	SO4	B	527	-	-	0/0/0/0	0/0/0/0

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	523	3FX	CAK-SAO	6.46	1.84	1.77
4	B	524	3FX	CAK-SAO	5.75	1.84	1.77
5	B	525	3FY	OAE-SBI	5.48	1.50	1.43
6	B	526	FII	P46-O49	5.31	1.62	1.50
5	B	525	3FY	OAD-SBI	5.09	1.50	1.43
6	B	526	FII	P46-O50	3.89	1.62	1.54
6	B	526	FII	C23-C22	-3.82	1.39	1.50
6	B	526	FII	C11-C10	-3.80	1.39	1.50
5	B	525	3FY	CBC-NBH	-3.64	1.33	1.39
4	B	522	3FX	CAK-SAO	3.52	1.81	1.77
6	B	526	FII	C34-C27	3.22	1.39	1.32
6	B	526	FII	C22-C15	3.19	1.39	1.32
6	B	526	FII	P46-O51	-3.15	1.48	1.54
6	B	526	FII	C35-C34	-2.81	1.39	1.49
5	B	525	3FY	CAR-NBG	-2.54	1.33	1.37
5	B	525	3FY	CAP-NBH	-2.47	1.32	1.36
6	B	526	FII	C10-C2	2.22	1.39	1.32
5	B	525	3FY	CAR-CBD	-2.19	1.34	1.37
4	B	522	3FX	OAB-SAO	2.06	1.51	1.46
5	B	525	3FY	CAO-CBC	-2.03	1.33	1.37
5	B	525	3FY	CBD-NAX	-2.01	1.34	1.38

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	525	3FY	OAE-SBI-OAD	-6.38	107.99	119.38
4	B	523	3FX	OAB-SAO-CAK	4.39	111.29	105.64
5	B	525	3FY	CAU-NBF-SBI	-4.31	107.28	117.36
4	B	522	3FX	OAB-SAO-CAK	4.09	110.91	105.64
5	B	525	3FY	CAS-CAT-NBF	-3.68	106.09	112.34
4	B	522	3FX	CAM-CAK-SAO	3.32	119.21	114.79
5	B	525	3FY	CAT-NBF-SBI	-3.31	110.74	117.84
6	B	526	FII	O51-P46-C45	3.17	113.37	106.85
5	B	525	3FY	CAU-NBF-CAT	-3.15	111.72	116.99
6	B	526	FII	C30-C27-C24	3.04	120.01	115.39
6	B	526	FII	C18-C15-C12	3.02	119.98	115.39
5	B	525	3FY	NAW-CAP-NBH	-2.89	106.56	112.02
6	B	526	FII	O50-P46-O49	-2.71	105.32	112.56
6	B	526	FII	O49-P46-C45	-2.43	105.35	110.22
5	B	525	3FY	OAD-SBI-NBF	2.42	109.17	106.71
5	B	525	3FY	CBC-CAV-NBE	2.38	117.44	113.73
5	B	525	3FY	OAE-SBI-NBF	2.24	109.00	106.71
4	B	524	3FX	OAD-SAO-CAK	2.21	110.91	107.03
5	B	525	3FY	NAX-CAQ-NBG	-2.18	107.40	112.55
6	B	526	FII	C6-C2-C1	2.12	120.02	114.62
5	B	525	3FY	CBD-SBI-NBF	2.09	111.00	106.19

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	315/349 (90%)	0.10	20 (6%) 19 22	24, 39, 62, 73	0
2	B	482/520 (92%)	0.10	19 (3%) 37 42	21, 30, 51, 82	0
All	All	797/869 (91%)	0.10	39 (4%) 28 31	21, 34, 59, 82	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	349	VAL	6.1
1	A	5	THR	5.8
2	B	51	VAL	5.1
1	A	334	ALA	4.6
1	A	103	VAL	4.6
2	B	348	LYS	4.1
2	B	347	ARG	3.5
2	B	2	ALA	3.3
1	A	7	ILE	3.2
1	A	139	LEU	3.1
1	A	208	ARG	3.1
1	A	13	ARG	3.0
2	B	255	GLU	3.0
2	B	453	PRO	2.9
1	A	207	SER	2.9
1	A	278	PRO	2.8
1	A	28	ASN	2.7
1	A	100	GLU	2.6
2	B	371	ILE	2.6
1	A	272	PHE	2.5
1	A	109	TYR	2.4
2	B	481	GLU	2.3
1	A	104	GLN	2.3
2	B	196	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	22	ARG	2.3
1	A	6	TYR	2.2
2	B	50	LEU	2.2
1	A	302	ASN	2.2
2	B	270	THR	2.2
1	A	105	ASN	2.2
2	B	330	VAL	2.2
2	B	265	ALA	2.1
2	B	269	TYR	2.1
1	A	108	SER	2.1
1	A	211	GLN	2.0
1	A	11	GLN	2.0
2	B	345	LYS	2.0
2	B	81	PHE	2.0
2	B	272	CYS	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
4	3FX	B	523	15/15	0.22	7.99	40,52,58,58	0
4	3FX	B	524	15/15	0.21	4.66	55,62,72,72	0
6	FII	B	526	24/24	0.24	1.88	28,33,35,38	0
5	3FY	B	525	35/35	0.24	1.35	37,54,62,62	35
4	3FX	B	522	15/15	0.11	1.08	28,29,32,33	0
7	SO4	B	527	5/5	0.17	0.81	51,53,56,56	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZN	B	521	1/1	0.11	-0.64	28,28,28,28	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.