



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:44 am GMT

PDB ID : 1LD8  
Title : Co-crystal structure of Human Farnesyltransferase with farnesyldiphosphate and inhibitor compound 49  
Authors : Taylor, J.S.; Terry, K.L.; Beese, L.S.  
Deposited on : 2002-04-08  
Resolution : 1.80 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.

---

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

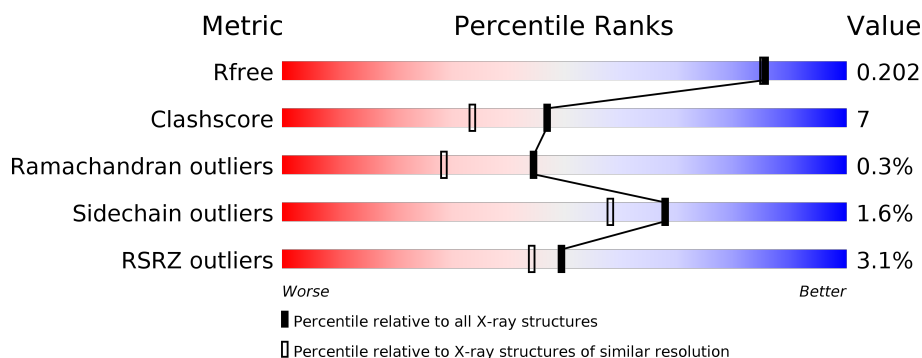
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	100719	4827 (1.80-1.80)
Clashscore	112137	5742 (1.80-1.80)
Ramachandran outliers	110173	5676 (1.80-1.80)
Sidechain outliers	110143	5675 (1.80-1.80)
RSRZ outliers	101464	4906 (1.80-1.80)

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	382	<div> <div>4%</div> <div> <div></div> <div>68%</div> <div>13%</div> <div>•</div> <div>18%</div> </div> </div>
2	B	437	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>15%</div> <div>•</div> <div>7%</div> </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
6	U49	B	1003	X	-	-	-

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6657 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 protein farnesyltransferase alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2670	1704	465	496	5			

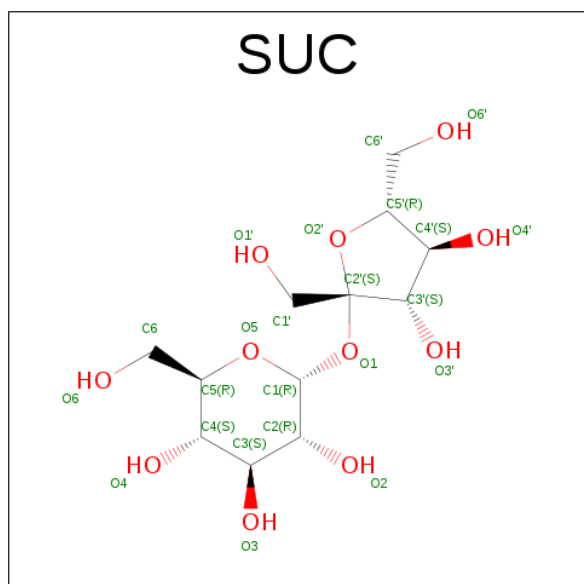
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	380	GLU	-	CLONING ARTIFACT	UNP P49354
A	381	GLU	-	CLONING ARTIFACT	UNP P49354
A	382	PHE	-	CLONING ARTIFACT	UNP P49354

- Molecule 2 is a protein called protein farnesyltransferase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	407	Total	C	N	O	S	0	0	0
			3207	2054	549	582	22			

- Molecule 3 is SUGAR (SUCROSE) (three-letter code: SUC) (formula: C<sub>12</sub>H<sub>22</sub>O<sub>11</sub>).

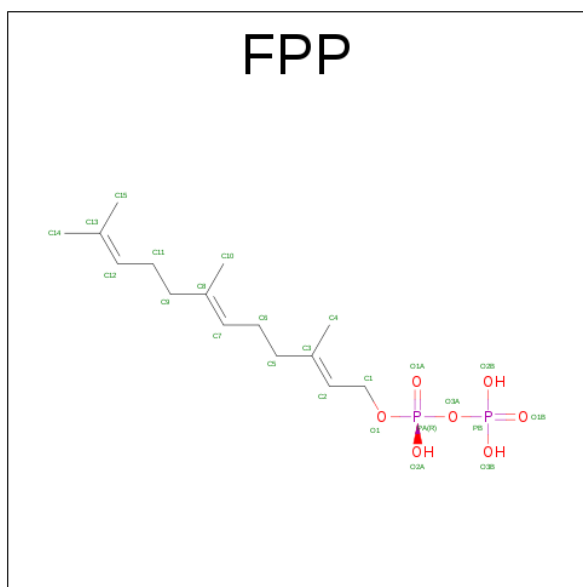


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			23	12	11		

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

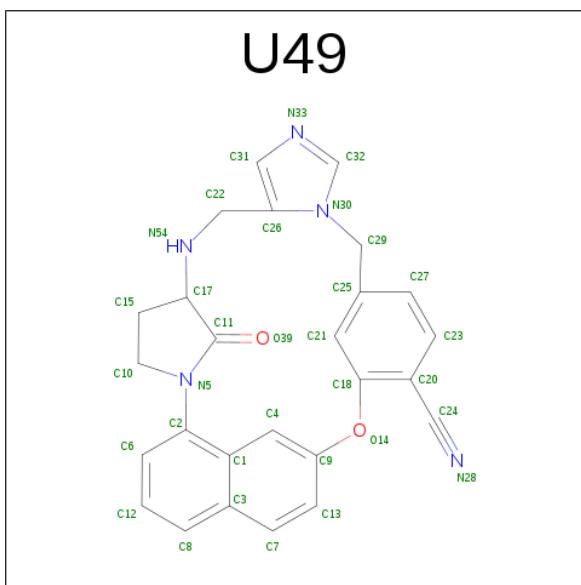
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		

- Molecule 5 is FARNESYL DIPHOSPHATE (three-letter code: FPP) (formula: C<sub>15</sub>H<sub>28</sub>O<sub>7</sub>P<sub>2</sub>).



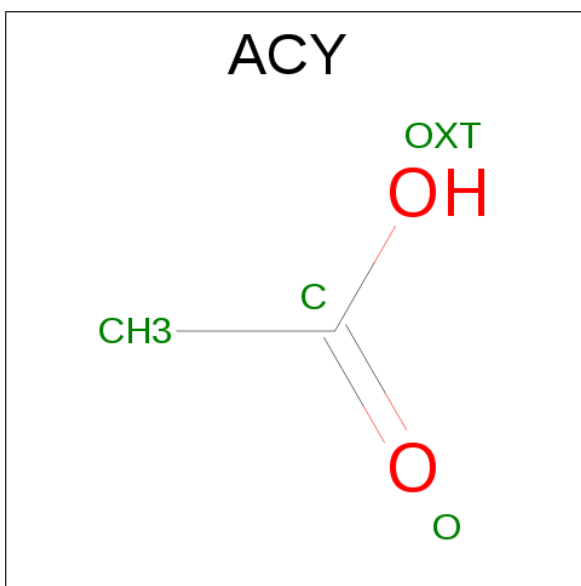
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	O	P	0	0
			24	15	7	2		

- Molecule 6 is (20S)-19,20,21,22-TETRAHYDRO-19-OXO-5H-18,20-ETHANO-12,14-ETHE NO-6,10-METHENO-18H-BENZ[D]IMIDAZO[4,3-K][1,6,9,12]OXATRIAZA-CYCLOOCT ADECOSINE-9-CARBONITRILE (three-letter code: U49) (formula: C<sub>26</sub>H<sub>21</sub>N<sub>5</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			33	26	5	2		

- Molecule 7 is ACETIC ACID (three-letter code: ACY) (formula:  $C_2H_4O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	A	1	Total	C	O	0	0
			4	2	2		

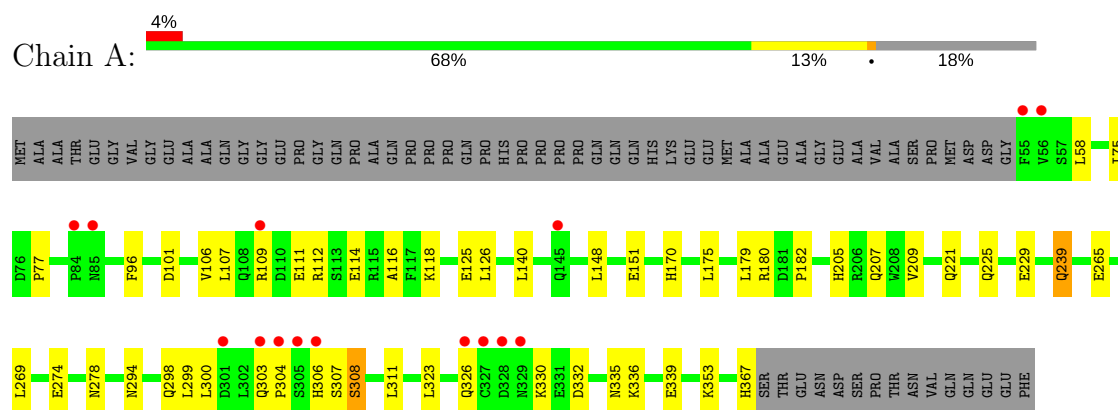
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	315	Total 315	O 315	0	0
8	B	376	Total 376	O 376	0	0

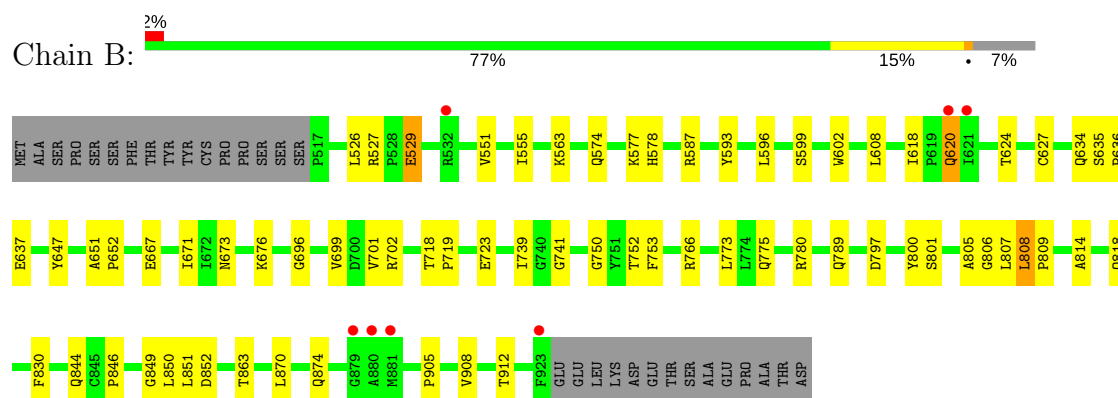
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: protein farnesyltransferase alpha subunit



- Molecule 2: protein farnesyltransferase beta subunit





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	178.13Å 178.13Å 64.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.85 – 1.80 31.02 – 1.80	Depositor EDS
% Data completeness (in resolution range)	80.8 (30.85-1.80) 81.8 (31.02-1.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.06 (at 1.80Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.184 , 0.202 0.185 , 0.202	Depositor DCC
$R_{free}$ test set	4441 reflections (5.04%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.0	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 49.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.020 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6657	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: ACY, ZN, U49, SUC, FPP

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	0/2737	0.50	0/3717
2	B	0.32	0/3296	0.57	1/4479 (0.0%)
All	All	0.31	0/6033	0.54	1/8196 (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	808	LEU	CA-CB-CG	-6.49	100.37	115.30

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	2670	0	2589	37	0
2	B	3207	0	3132	48	0
3	B	23	0	22	0	0
4	B	1	0	0	0	0
5	B	24	0	25	2	0
6	B	33	0	20	1	0
7	A	8	0	6	1	0
8	A	315	0	0	5	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	B	376	0	0	1	0
All	All	6657	0	5794	84	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 7.

All (84) 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)
2:B:578:HIS:HD1	2:B:849:GLY:H	1.12	0.92
2:B:574:GLN:H	2:B:844:GLN:HE22	1.20	0.90
1:A:106:VAL:HG23	1:A:111:GLU:HB3	1.55	0.88
2:B:634:GLN:HE22	2:B:673:ASN:H	1.26	0.84
2:B:627:CYS:HB3	2:B:671:ILE:HD11	1.62	0.82
2:B:780:ARG:HE	2:B:789:GLN:HE21	1.27	0.81
1:A:303:GLN:O	1:A:307:SER:HB2	1.86	0.75
1:A:303:GLN:HB3	1:A:304:PRO:HD3	1.74	0.70
2:B:574:GLN:H	2:B:844:GLN:NE2	1.93	0.65
2:B:801:SER:O	2:B:805:ALA:HB3	1.97	0.64
2:B:608:LEU:HD22	2:B:618:ILE:HD13	1.81	0.62
2:B:577:LYS:HD3	2:B:846:PRO:O	1.99	0.62
2:B:526:LEU:HD13	2:B:563:LYS:HB2	1.82	0.61
2:B:624:THR:HG22	2:B:667:GLU:OE1	2.01	0.60
2:B:808:LEU:HD13	2:B:830:PHE:HB3	1.83	0.60
2:B:739:ILE:HB	2:B:752:THR:HA	1.83	0.60
2:B:780:ARG:HE	2:B:789:GLN:NE2	1.98	0.60
1:A:307:SER:O	1:A:308:SER:HB3	2.03	0.59
2:B:753:PHE:HA	2:B:807:LEU:HD21	1.85	0.58
2:B:634:GLN:NE2	2:B:673:ASN:H	1.98	0.58
2:B:702:ARG:HD2	5:B:1002:FPP:H142	1.86	0.58
2:B:593:TYR:CD2	2:B:596:LEU:HD12	2.39	0.57
2:B:750:GLY:HA3	5:B:1002:FPP:C8	2.35	0.57
1:A:299:LEU:HB3	1:A:311:LEU:HD11	1.87	0.57
2:B:870:LEU:O	2:B:874:GLN:HG3	2.05	0.56
1:A:335:ASN:O	1:A:339:GLU:HG3	2.05	0.56
1:A:323:LEU:HA	1:A:330:LYS:HE3	1.88	0.56
1:A:106:VAL:HG21	1:A:116:ALA:CB	2.36	0.56
1:A:353:LYS:HD3	8:A:1985:HOH:O	2.07	0.55
2:B:620:GLN:OE1	2:B:624:THR:HG23	2.08	0.54
1:A:207:GLN:HE21	1:A:239:GLN:NE2	2.06	0.53
1:A:332:ASP:O	1:A:336:LYS:HG2	2.10	0.52

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:VAL:HG21	1:A:116:ALA:HB1	1.92	0.52
2:B:808:LEU:HD13	2:B:830:PHE:CD2	2.45	0.52
1:A:294:ASN:O	1:A:298:GLN:HG3	2.09	0.52
1:A:58:LEU:HD11	1:A:126:LEU:HG	1.91	0.52
2:B:701:VAL:HG23	2:B:741:GLY:O	2.10	0.52
2:B:578:HIS:HD1	2:B:849:GLY:N	1.95	0.52
1:A:239:GLN:HA	1:A:239:GLN:HE21	1.75	0.52
2:B:574:GLN:N	2:B:844:GLN:HE22	2.00	0.51
1:A:225:GLN:O	1:A:229:GLU:HG3	2.11	0.51
1:A:323:LEU:HD23	1:A:330:LYS:HD2	1.91	0.51
2:B:620:GLN:HE22	2:B:624:THR:HG21	1.74	0.51
1:A:221:GLN:HG3	8:A:1180:HOH:O	2.11	0.50
2:B:806:GLY:O	2:B:809:PRO:HD2	2.14	0.48
1:A:274:GLU:HG2	1:A:278:ASN:ND2	2.29	0.48
1:A:114:GLU:OE2	1:A:118:LYS:HE3	2.15	0.47
7:A:904:ACY:H1	6:B:1003:U49:H152	1.97	0.47
2:B:635:SER:OG	2:B:637:GLU:HG2	2.15	0.46
1:A:170:HIS:HD2	8:A:1243:HOH:O	1.97	0.46
1:A:107:LEU:HD11	2:B:647:TYR:OH	2.16	0.46
2:B:718:THR:HB	2:B:719:PRO:HD2	1.97	0.46
2:B:766:ARG:HG2	2:B:818:GLN:NE2	2.31	0.46
1:A:96:PHE:HA	1:A:126:LEU:HD13	1.99	0.45
2:B:551:VAL:O	2:B:555:ILE:HG12	2.17	0.45
2:B:850:LEU:HB2	2:B:863:THR:HA	1.98	0.45
2:B:814:ALA:O	2:B:818:GLN:HG3	2.17	0.45
2:B:797:ASP:HB3	2:B:800:TYR:CD1	2.51	0.45
2:B:651:ALA:HB3	2:B:652:PRO:CD	2.46	0.44
1:A:170:HIS:HE1	2:B:696:GLY:O	2.00	0.44
1:A:265:GLU:OE2	1:A:269:LEU:HD13	2.16	0.44
1:A:58:LEU:HD12	1:A:125:GLU:HB3	2.00	0.44
2:B:908:VAL:O	2:B:912:THR:HG23	2.17	0.44
2:B:773:LEU:O	2:B:773:LEU:HD23	2.18	0.44
2:B:627:CYS:HB3	2:B:671:ILE:CD1	2.42	0.43
2:B:766:ARG:HG2	2:B:818:GLN:HE21	1.84	0.43
1:A:326:GLN:HE21	1:A:326:GLN:HA	1.84	0.43
1:A:77:PRO:HB2	1:A:101:ASP:HB3	2.01	0.43
2:B:527:ARG:HB3	2:B:529:GLU:OE1	2.19	0.42
2:B:599:SER:O	2:B:602:TRP:HB2	2.19	0.42
2:B:780:ARG:NE	2:B:789:GLN:HE21	2.04	0.42
2:B:852:ASP:OD1	2:B:852:ASP:C	2.58	0.41
2:B:905:PRO:HD2	8:B:1306:HOH:O	2.19	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ARG:HA	1:A:140:LEU:CD2	2.50	0.41
1:A:239:GLN:NE2	1:A:239:GLN:HA	2.36	0.41
1:A:148:LEU:HB2	1:A:179:LEU:HD21	2.02	0.41
8:A:1213:HOH:O	2:B:775:GLN:HG2	2.20	0.41
1:A:205:HIS:O	1:A:209:VAL:HG23	2.21	0.40
1:A:367:HIS:HE1	8:A:1325:HOH:O	2.04	0.40
1:A:75:ILE:HD12	1:A:109:ARG:HH11	1.86	0.40
1:A:151:GLU:HG3	1:A:175:LEU:HD11	2.02	0.40
1:A:207:GLN:HE21	1:A:239:GLN:HE22	1.68	0.40
2:B:636:PRO:O	2:B:676:LYS:HE2	2.21	0.40
1:A:239:GLN:CA	1:A:239:GLN:HE21	2.34	0.40

There are no symmetry-related clashes.

## 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	311/382 (81%)	295 (95%)	14 (4%)	2 (1%)	28	13
2	B	405/437 (93%)	396 (98%)	9 (2%)	0	100	100
All	All	716/819 (87%)	691 (96%)	23 (3%)	2 (0%)	44	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	308	SER
1	A	306	HIS

### 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	292/344 (85%)	288 (99%)	4 (1%)	71	64
2	B	343/370 (93%)	337 (98%)	6 (2%)	66	55
All	All	635/714 (89%)	625 (98%)	10 (2%)	68	58

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	180	ARG
1	A	182	PRO
1	A	239	GLN
1	A	300	LEU
2	B	529	GLU
2	B	587	ARG
2	B	620	GLN
2	B	699	VAL
2	B	723	GLU
2	B	851	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	HIS
1	A	170	HIS
1	A	211	GLN
1	A	218	ASN
1	A	221	GLN
1	A	239	GLN
1	A	246	ASN
1	A	261	GLN
1	A	298	GLN
1	A	303	GLN
1	A	326	GLN
1	A	367	HIS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
2	B	628	GLN
2	B	634	GLN
2	B	789	GLN
2	B	818	GLN
2	B	844	GLN
2	B	875	HIS
2	B	910	GLN

### 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 ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	$\# Z  > 2$
7	ACY	A	904	-	1,3,3	1.55	0	0,3,3	0.00	-
7	ACY	A	905	-	1,3,3	3.34	1 (100%)	0,3,3	0.00	-
5	FPP	B	1002	-	22,23,23	0.62	0	24,31,31	0.72	0
6	U49	B	1003	4	34,38,38	1.72	9 (26%)	44,55,55	2.05	9 (20%)
3	SUC	B	938	-	24,24,24	1.56	3 (12%)	36,36,36	0.95	2 (5%)

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
7	ACY	A	904	-	-	0/0/0/0	0/0/0/0
7	ACY	A	905	-	-	0/0/0/0	0/0/0/0
5	FPP	B	1002	-	-	0/25/25/25	0/0/0/0
6	U49	B	1003	4	1/1/3/4	0/16/32/32	0/2/6/6
3	SUC	B	938	-	-	0/12/51/51	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	1003	U49	C17-N54	-3.35	1.41	1.46
6	B	1003	U49	C17-C11	-2.31	1.48	1.52
6	B	1003	U49	C4-C9	2.06	1.40	1.37
3	B	938	SUC	O5-C1	2.28	1.47	1.41
6	B	1003	U49	C21-C25	2.39	1.43	1.39
3	B	938	SUC	C4-C5	2.47	1.58	1.53
6	B	1003	U49	C27-C23	2.49	1.43	1.38
6	B	1003	U49	C12-C8	2.53	1.42	1.36
6	B	1003	U49	C6-C2	2.96	1.44	1.38
6	B	1003	U49	C7-C13	3.00	1.42	1.36
7	A	905	ACY	CH3-C	3.34	1.53	1.48
6	B	1003	U49	C11-N5	3.55	1.40	1.37
3	B	938	SUC	C3-C2	5.55	1.66	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	1003	U49	C10-C15-C17	-4.12	97.53	104.45
6	B	1003	U49	C22-N54-C17	-3.96	106.40	113.99
6	B	1003	U49	C31-C26-N30	-2.27	104.30	108.06
3	B	938	SUC	C1-C2-C3	-2.20	105.89	109.98
6	B	1003	U49	O14-C18-C21	-2.01	115.99	121.69
3	B	938	SUC	C1-O5-C5	2.05	117.58	113.72
6	B	1003	U49	C9-C4-C1	2.44	123.50	120.05
6	B	1003	U49	O39-C11-N5	2.83	129.84	125.88
6	B	1003	U49	C6-C2-N5	3.14	124.72	119.34
6	B	1003	U49	C15-C17-N54	5.53	126.67	115.48
6	B	1003	U49	C11-C17-N54	7.95	124.88	111.10

All (1) chirality outliers are listed below:



Mol	Chain	Res	Type	Atom
6	B	1003	U49	C17

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	904	ACY	1	0
5	B	1002	FPP	2	0
6	B	1003	U49	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, 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	313/382 (81%)	-0.06	15 (4%) 31 26	15, 24, 39, 54	0
2	B	407/437 (93%)	-0.13	7 (1%) 70 67	13, 19, 31, 40	0
All	All	720/819 (87%)	-0.10	22 (3%) 49 44	13, 21, 34, 54	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	55	PHE	7.8
1	A	306	HIS	6.4
1	A	304	PRO	5.7
2	B	880	ALA	4.9
2	B	923	PHE	4.6
1	A	328	ASP	4.1
1	A	305	SER	3.8
1	A	109	ARG	3.5
2	B	620	GLN	3.4
1	A	329	ASN	3.1
2	B	881	MET	3.0
1	A	303	GLN	2.8
1	A	326	GLN	2.8
1	A	301	ASP	2.8
1	A	56	VAL	2.5
2	B	879	GLY	2.4
2	B	532	ARG	2.2
1	A	145	GLN	2.2
1	A	327	CYS	2.1
1	A	84	PRO	2.0
2	B	621	ILE	2.0
1	A	85	ASN	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, 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( $\text{\AA}^2$ )	Q<0.9
3	SUC	B	938	23/23	0.86	0.13	1.57	27,29,32,33	0
5	FPP	B	1002	24/24	0.96	0.13	0.94	15,17,18,18	0
6	U49	B	1003	33/33	0.94	0.11	0.55	14,17,25,26	0
4	ZN	B	1001	1/1	1.00	0.05	-	18,18,18,18	0
7	ACY	A	905	4/4	0.72	0.29	-	43,43,44,44	0
7	ACY	A	904	4/4	0.85	0.19	-	44,44,44,44	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.