Full wwPDB X-ray Structure Validation Report

Feb 1, 2016 – 01:32 PM GMT

PDB ID	:	3TZY
Title	:	Crystal structure of a fragment containing the acyltransferase domain of Pks13 from Mycobacterium tuberculosis in the palmitoylated form at 2.2 A
Authors	:	Bergeret, F.; Pedelacq, J.D.; Mourey, L.; Bon, C.
Deposited on	:	2011-09-28
Resolution	:	2.20 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
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

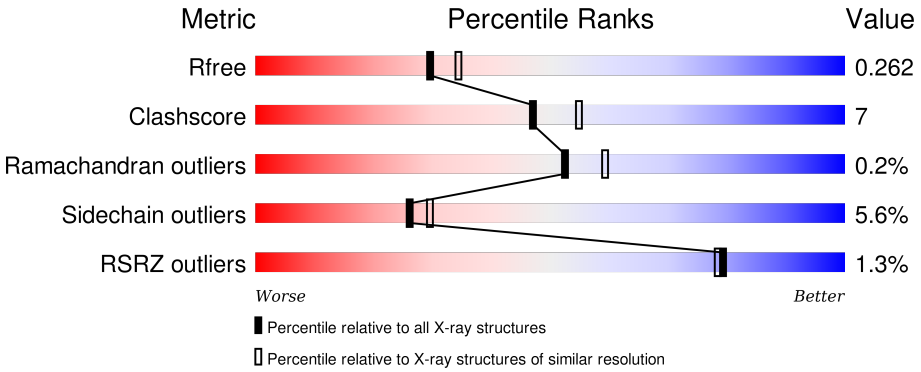
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (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 >=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 <=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	491	2% 79% 14% • 5%
1	B	491	82% 12% • 5%
2	C	12	8% 42% 33% 8% 17%
2	D	12	17% 50% 33% 17%

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
3	PLM	A	2001	-	-	-	X
3	PLM	B	2001	-	-	-	X
4	GOL	A	2004	-	-	-	X
4	GOL	A	2005	-	-	-	X
4	GOL	B	2002	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7427 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 Polyketide synthase PKS13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3481	2204	600	665	12			
1	B	466	Total	C	N	O	S	0	0	0
			3478	2202	602	663	11			

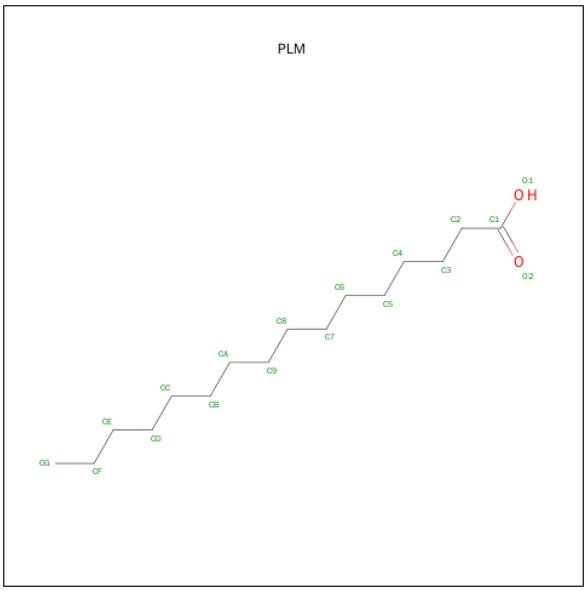
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	572	GLY	-	EXPRESSION TAG	UNP O53579
A	573	SER	-	EXPRESSION TAG	UNP O53579
A	574	HIS	-	EXPRESSION TAG	UNP O53579
A	575	MET	-	EXPRESSION TAG	UNP O53579
B	572	GLY	-	EXPRESSION TAG	UNP O53579
B	573	SER	-	EXPRESSION TAG	UNP O53579
B	574	HIS	-	EXPRESSION TAG	UNP O53579
B	575	MET	-	EXPRESSION TAG	UNP O53579

- Molecule 2 is a protein called 12-mer peptide.

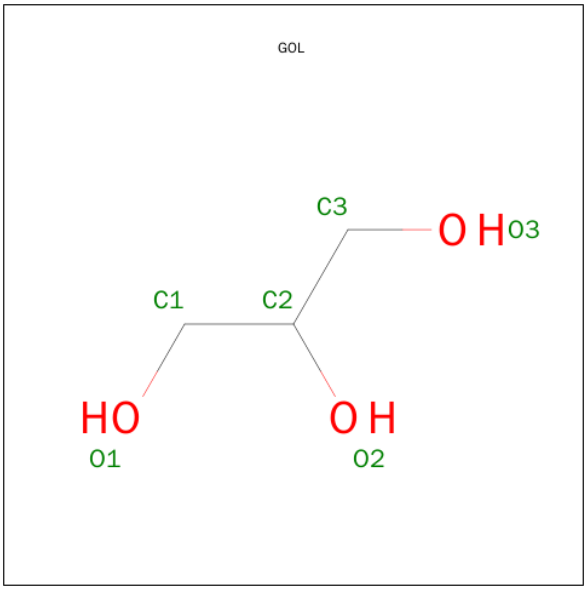
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	10	Total	C	N	O	S	0	0	0
			83	54	13	15	1			
2	D	10	Total	C	N	O	S	0	0	0
			65	41	12	11	1			

- Molecule 3 is PALMITIC ACID (three-letter code: PLM) (formula: $C_{16}H_{32}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 17 16 1	0	0
3	B	1	Total C O 17 16 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0

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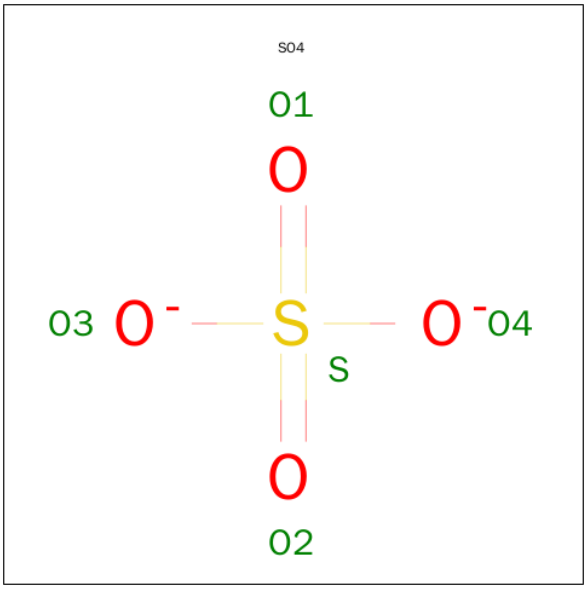
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	B	1	Total C O 6 3 3	0	0
4	B	1	Total C O 6 3 3	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	113	Total O 113 113	0	0
6	B	120	Total O 120 120	0	0

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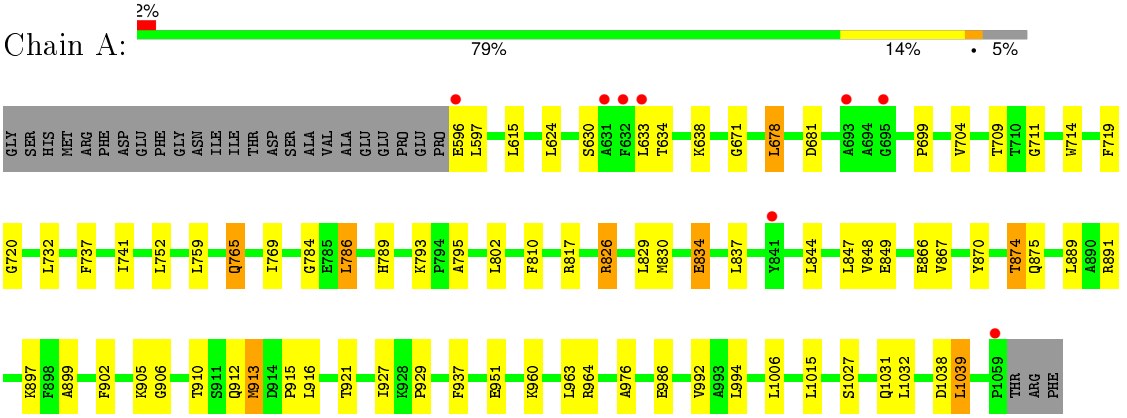
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	1	Total O 1 1	0	0

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.

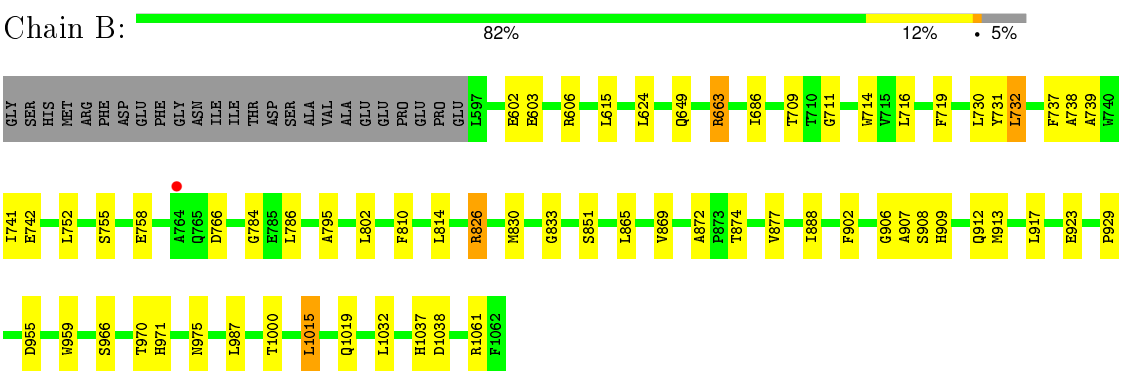
- Molecule 1: Polyketide synthase PKS13

Chain A:



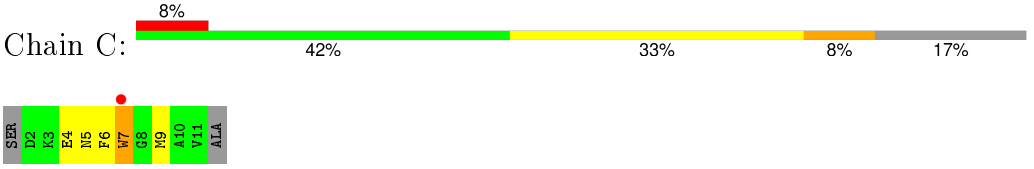
- Molecule 1: Polyketide synthase PKS13

Chain B:



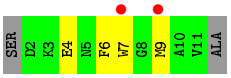
- Molecule 2: 12-mer peptide

Chain C:



- Molecule 2: 12-mer peptide

Chain D:


SER
D2
K3
E4
N5
F6
W7
G8
M9
A10
V11
ALA

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	105.89Å 105.89Å 259.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 49.04 – 2.20	Depositor EDS
% Data completeness (in resolution range)	91.1 (50.00-2.20) 91.1 (49.04-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	3.21 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.223 , 0.265 0.219 , 0.262	Depositor DCC
R_{free} test set	3787 reflections (5.78%)	DCC
Wilson B-factor (Å^2)	43.9	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 42.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Outliers	0 of 69008 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7427	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, PLM, SO4

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	0.58	1/3554 (0.0%)	0.69	0/4835
1	B	0.64	2/3551 (0.1%)	0.75	2/4834 (0.0%)
2	C	1.24	1/85 (1.2%)	0.89	0/113
2	D	0.87	0/66	0.87	0/89
All	All	0.63	4/7256 (0.1%)	0.72	2/9871 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	714	TRP	CD2-CE2	5.22	1.47	1.41
2	C	7	TRP	NE1-CE2	-5.18	1.30	1.37
1	B	714	TRP	CD2-CE2	5.13	1.47	1.41
1	B	959	TRP	CD2-CE2	5.05	1.47	1.41

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	663	ARG	NE-CZ-NH2	-7.24	116.68	120.30
1	B	955	ASP	CB-CA-C	-5.01	100.37	110.40

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	3481	0	3416	57	0
1	B	3478	0	3387	46	0
2	C	83	0	73	5	0
2	D	65	0	44	5	0
3	A	17	0	31	7	0
3	B	17	0	31	8	0
4	A	30	0	40	2	0
4	B	12	0	16	2	0
5	B	5	0	0	0	0
5	C	5	0	0	1	0
6	A	113	0	0	1	0
6	B	120	0	0	3	0
6	D	1	0	0	0	0
All	All	7427	0	7038	106	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 (106) 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)
1:B:913:MET:HE1	3:B:2001:PLM:HD2	1.60	0.83
1:A:834:GLU:HB3	3:A:2001:PLM:HC1	1.60	0.81
1:A:634:THR:HG22	1:A:638:LYS:HE3	1.65	0.79
1:B:906:GLY:HA3	3:B:2001:PLM:HA1	1.67	0.77
1:A:784:GLY:HA3	1:A:810:PHE:CZ	2.21	0.75
1:A:769:ILE:HG13	1:A:834:GLU:HG2	1.68	0.75
1:A:720:GLY:HA2	4:A:2004:GOL:H32	1.70	0.72
1:B:907:ALA:H	1:B:912:GLN:HE21	1.41	0.68
1:B:624:LEU:HD11	2:C:7:TRP:HZ3	1.59	0.68
1:B:602:GLU:CD	1:B:602:GLU:H	1.98	0.67
1:B:826:ARG:NH1	3:B:2001:PLM:O2	2.24	0.67
1:B:872:ALA:HA	1:B:970:THR:CG2	2.27	0.65
1:B:624:LEU:CD1	2:C:7:TRP:HZ3	2.09	0.65
1:A:837:LEU:O	1:A:905:LYS:HE2	1.97	0.64
1:A:913:MET:HG2	1:A:963:LEU:HG	1.81	0.62
1:A:906:GLY:HA3	3:A:2001:PLM:HC2	1.82	0.62
1:B:755:SER:OG	1:B:758:GLU:HG3	2.00	0.61
1:A:624:LEU:HD11	2:D:7:TRP:HZ3	1.64	0.61
1:A:719:PHE:HD1	3:A:2001:PLM:H32	1.65	0.61
1:A:847:LEU:HD11	1:A:875:GLN:HB3	1.83	0.61
1:A:927:ILE:O	1:A:929:PRO:HD3	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:865:LEU:HD11	1:B:888:ILE:CD1	2.32	0.59
1:A:711:GLY:HA3	1:A:795:ALA:HB2	1.86	0.58
1:B:731:TYR:HD2	1:B:732:LEU:HD13	1.70	0.56
1:A:624:LEU:CD1	2:D:7:TRP:HZ3	2.19	0.56
1:B:970:THR:HG22	1:B:1000:THR:OG1	2.06	0.55
1:B:731:TYR:OH	1:B:742:GLU:OE1	2.22	0.55
1:A:597:LEU:HD21	1:A:817:ARG:NH2	2.21	0.55
1:A:681:ASP:HB2	4:A:2006:GOL:H2	1.88	0.55
1:B:907:ALA:H	1:B:912:GLN:NE2	2.04	0.55
1:B:663:ARG:NH2	6:B:2111:HOH:O	2.40	0.55
1:A:867:VAL:HG11	1:A:870:TYR:CZ	2.42	0.55
1:B:1019:GLN:OE1	2:C:6:PHE:HB3	2.07	0.54
1:B:737:PHE:CE1	1:B:741:ILE:HD11	2.43	0.54
1:B:624:LEU:HD11	2:C:7:TRP:CZ3	2.42	0.54
1:A:719:PHE:CD1	3:A:2001:PLM:H32	2.43	0.54
2:D:4:GLU:HA	2:D:9:MET:O	2.08	0.54
1:A:624:LEU:HD21	1:A:678:LEU:HG	1.90	0.53
1:B:649:GLN:HG3	1:B:686:ILE:HD13	1.89	0.53
1:A:826:ARG:O	1:A:830:MET:HG3	2.09	0.53
1:A:912:GLN:O	1:A:915:PRO:CD	2.58	0.52
1:A:802:LEU:HB3	3:A:2001:PLM:H31	1.93	0.51
1:B:865:LEU:HD11	1:B:888:ILE:HD13	1.92	0.51
1:B:1037:HIS:ND1	4:B:2002:GOL:O2	2.44	0.51
1:B:971:HIS:O	1:B:975:ASN:HB2	2.11	0.51
1:B:909:HIS:HE2	3:B:2001:PLM:C1	2.23	0.51
1:A:913:MET:HG3	1:A:916:LEU:HD12	1.93	0.51
1:A:847:LEU:O	1:A:899:ALA:HA	2.10	0.51
1:A:834:GLU:HA	1:A:837:LEU:HD12	1.92	0.50
1:B:755:SER:CB	1:B:758:GLU:HG3	2.42	0.50
1:A:759:LEU:HD22	1:A:765:GLN:HG2	1.94	0.49
1:B:869:VAL:HB	1:B:877:VAL:HB	1.94	0.49
1:A:634:THR:CG2	1:A:638:LYS:HE3	2.39	0.48
1:A:786:LEU:O	1:A:789:HIS:HB3	2.13	0.48
1:A:866:GLU:OE1	1:A:910:THR:HB	2.13	0.48
1:A:848:VAL:HG23	1:A:899:ALA:HB2	1.96	0.48
1:A:630:SER:HG	1:A:671:GLY:H	1.59	0.48
1:A:737:PHE:CE2	1:A:741:ILE:HD12	2.48	0.48
1:B:1038:ASP:H	4:B:2002:GOL:H11	1.79	0.47
1:A:1038:ASP:O	1:A:1039:LEU:HB2	2.14	0.47
1:A:844:LEU:HB3	6:A:2171:HOH:O	2.14	0.47
1:B:906:GLY:CA	3:B:2001:PLM:HA1	2.41	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:826:ARG:NH2	3:A:2001:PLM:O2	2.43	0.47
1:A:912:GLN:O	1:A:915:PRO:HD3	2.14	0.47
1:B:830:MET:HE1	1:B:908:SER:HB3	1.96	0.47
1:B:711:GLY:HA3	1:B:795:ALA:HB2	1.95	0.47
1:B:865:LEU:HD11	1:B:888:ILE:HD11	1.95	0.46
1:B:716:LEU:HD11	1:B:987:LEU:HD12	1.96	0.46
1:B:802:LEU:C	1:B:802:LEU:HD12	2.36	0.46
1:A:874:THR:CG2	1:A:874:THR:O	2.64	0.46
1:B:833:GLY:HA3	3:B:2001:PLM:HE2	1.97	0.46
1:B:826:ARG:O	1:B:830:MET:HG3	2.16	0.45
1:A:849:GLU:HB3	1:A:897:LYS:HB3	1.98	0.45
1:A:624:LEU:HD11	2:D:7:TRP:CZ3	2.48	0.45
1:A:699:PRO:HB2	2:D:7:TRP:CD1	2.52	0.45
1:A:837:LEU:O	1:A:905:LYS:CE	2.65	0.45
1:A:769:ILE:HG13	1:A:834:GLU:CG	2.41	0.44
1:A:902:PHE:CD2	1:A:902:PHE:N	2.85	0.44
1:B:784:GLY:HA3	1:B:810:PHE:CZ	2.52	0.44
1:A:937:PHE:CZ	1:A:976:ALA:HA	2.52	0.44
1:B:814:LEU:HD23	1:B:929:PRO:HA	2.00	0.44
1:A:834:GLU:HA	1:A:837:LEU:CD1	2.48	0.44
1:A:737:PHE:CE2	1:A:741:ILE:CD1	3.02	0.43
1:A:1027:SER:O	1:A:1031:GLN:HG3	2.18	0.43
1:B:906:GLY:HA3	3:B:2001:PLM:HC2	2.01	0.43
1:A:759:LEU:CD2	1:A:765:GLN:HG2	2.47	0.43
2:C:5:ASN:HB2	5:C:101:SO4:O1	2.18	0.43
1:B:755:SER:HB3	1:B:758:GLU:HG3	2.01	0.43
1:A:921:THR:OG1	1:A:960:LYS:HD3	2.19	0.43
1:B:830:MET:CE	1:B:908:SER:HB3	2.49	0.43
1:B:872:ALA:CA	1:B:970:THR:CG2	2.95	0.42
1:A:912:GLN:O	1:A:915:PRO:HD2	2.18	0.42
1:B:603:GLU:OE1	1:B:603:GLU:HA	2.19	0.42
1:B:906:GLY:HA2	1:B:912:GLN:NE2	2.35	0.42
1:B:719:PHE:CE2	1:B:902:PHE:CE2	3.08	0.42
3:B:2001:PLM:H61	6:B:2207:HOH:O	2.20	0.41
1:A:910:THR:O	1:A:964:ARG:HD2	2.20	0.41
1:A:802:LEU:HD23	3:A:2001:PLM:H31	2.01	0.41
1:A:829:LEU:HD23	1:A:829:LEU:HA	1.80	0.41
1:A:902:PHE:HD2	1:A:902:PHE:N	2.18	0.41
1:B:738:ALA:O	1:B:739:ALA:C	2.58	0.41
1:A:826:ARG:HG3	1:A:830:MET:SD	2.61	0.41
1:A:681:ASP:OD1	1:A:681:ASP:C	2.59	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:986:GLU:HB2	1:A:994:LEU:HG	2.02	0.41
1:B:730:LEU:HD22	1:B:1015:LEU:HD11	2.02	0.41
1:B:755:SER:HA	6:B:2121:HOH:O	2.20	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	462/491 (94%)	446 (96%)	15 (3%)	1 (0%)	52 59
1	B	464/491 (94%)	445 (96%)	19 (4%)	0	100 100
2	C	8/12 (67%)	6 (75%)	2 (25%)	0	100 100
2	D	8/12 (67%)	6 (75%)	1 (12%)	1 (12%)	0 0
All	All	942/1006 (94%)	903 (96%)	37 (4%)	2 (0%)	52 59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1039	LEU
2	D	6	PHE

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	353/387 (91%)	331 (94%)	22 (6%)	23 25
1	B	346/387 (89%)	330 (95%)	16 (5%)	33 40
2	C	8/9 (89%)	6 (75%)	2 (25%)	1 0
2	D	3/9 (33%)	3 (100%)	0	100 100
All	All	710/792 (90%)	670 (94%)	40 (6%)	26 29

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	596	GLU
1	A	615	LEU
1	A	633	LEU
1	A	678	LEU
1	A	704	VAL
1	A	709	THR
1	A	732	LEU
1	A	752	LEU
1	A	765	GLN
1	A	786	LEU
1	A	793	LYS
1	A	826	ARG
1	A	834	GLU
1	A	874	THR
1	A	889	LEU
1	A	891	ARG
1	A	913	MET
1	A	951	GLU
1	A	992	VAL
1	A	1006	LEU
1	A	1015	LEU
1	A	1032	LEU
1	B	606	ARG
1	B	615	LEU
1	B	709	THR
1	B	732	LEU
1	B	752	LEU
1	B	766	ASP
1	B	786	LEU
1	B	826	ARG
1	B	851	SER
1	B	874	THR
1	B	917	LEU

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Mol	Chain	Res	Type
1	B	923	GLU
1	B	966	SER
1	B	1015	LEU
1	B	1032	LEU
1	B	1061	ARG
2	C	4	GLU
2	C	9	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	654	GLN
1	A	1019	GLN
1	B	680	HIS
1	B	749	GLN
1	B	912	GLN
1	B	1031	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

11 ligands are modelled in this entry.

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	PLM	A	2001	1	16,16,17	0.57	0	14,15,17	0.27	0
4	GOL	A	2002	-	5,5,5	0.38	0	5,5,5	0.68	0
4	GOL	A	2003	-	5,5,5	0.33	0	5,5,5	0.26	0
4	GOL	A	2004	-	5,5,5	0.29	0	5,5,5	0.45	0
4	GOL	A	2005	-	5,5,5	0.43	0	5,5,5	0.30	0
4	GOL	A	2006	-	5,5,5	0.29	0	5,5,5	0.36	0
3	PLM	B	2001	1	16,16,17	0.47	0	14,15,17	0.74	1 (7%)
4	GOL	B	2002	-	5,5,5	0.62	0	5,5,5	0.58	0
4	GOL	B	2003	-	5,5,5	0.32	0	5,5,5	0.71	0
5	SO4	B	2004	-	4,4,4	0.41	0	6,6,6	0.13	0
5	SO4	C	101	-	4,4,4	0.35	0	6,6,6	0.29	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	PLM	A	2001	1	-	0/13/14/15	0/0/0/0
4	GOL	A	2002	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2003	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2004	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2005	-	-	0/4/4/4	0/0/0/0
4	GOL	A	2006	-	-	0/4/4/4	0/0/0/0
3	PLM	B	2001	1	-	0/13/14/15	0/0/0/0
4	GOL	B	2002	-	-	0/4/4/4	0/0/0/0
4	GOL	B	2003	-	-	0/4/4/4	0/0/0/0
5	SO4	B	2004	-	-	0/0/0/0	0/0/0/0
5	SO4	C	101	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	2001	PLM	C3-C2-C1	-2.21	107.00	113.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2001	PLM	7	0
4	A	2004	GOL	1	0
4	A	2006	GOL	1	0
3	B	2001	PLM	8	0
4	B	2002	GOL	2	0
5	C	101	SO4	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^{th} 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^2)	Q<0.9
1	A	464/491 (94%)	-0.14	8 (1%) 73 72	31, 44, 61, 94	0
1	B	466/491 (94%)	-0.19	1 (0%) 95 95	26, 41, 65, 87	0
2	C	10/12 (83%)	0.28	1 (10%) 9 8	48, 53, 64, 69	0
2	D	10/12 (83%)	1.22	2 (20%) 1 1	74, 84, 93, 97	0
All	All	950/1006 (94%)	-0.14	12 (1%) 79 78	26, 43, 65, 97	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	7	TRP	5.3
1	A	631	ALA	3.5
1	A	1059	PRO	2.9
1	B	764	ALA	2.9
1	A	632	PHE	2.9
2	D	9	MET	2.8
1	A	633	LEU	2.7
1	A	695	GLY	2.7
1	A	841	TYR	2.5
2	C	7	TRP	2.5
1	A	596	GLU	2.3
1	A	693	ALA	2.2

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^{th} 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^2)	Q<0.9
3	PLM	A	2001	17/18	0.88	0.26	8.74	42,53,75,75	0
3	PLM	B	2001	17/18	0.85	0.27	6.18	41,55,65,68	0
4	GOL	B	2002	6/6	0.81	0.20	5.46	47,54,58,62	0
4	GOL	A	2005	6/6	0.85	0.27	3.68	55,66,69,72	0
4	GOL	A	2004	6/6	0.87	0.18	2.39	64,66,70,75	0
4	GOL	B	2003	6/6	0.85	0.13	-	51,55,56,57	0
5	SO4	B	2004	5/5	0.96	0.18	-	89,96,98,100	0
4	GOL	A	2002	6/6	0.83	0.17	-	53,55,60,66	0
4	GOL	A	2006	6/6	0.73	0.21	-	58,71,76,79	0
4	GOL	A	2003	6/6	0.78	0.21	-	63,67,69,69	0
5	SO4	C	101	5/5	0.98	0.10	-	68,71,72,73	0

6.5 Other polymers

There are no such residues in this entry.