



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:38 AM GMT

PDB ID : 2XT6
Title : CRYSTAL STRUCTURE OF MYCOBACTERIUM SMEGMATIS ALPHA-KETOGLUTARATE DECARBOXYLASE HOMODIMER (ORTHORHOMBIC FORM)
Authors : Wagner, T.; Bellinzoni, M.; Wehenkel, A.M.; O'Hare, H.M.; Alzari, P.M.
Deposited on : 2010-10-05
Resolution : 2.74 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 ⓘ 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 (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

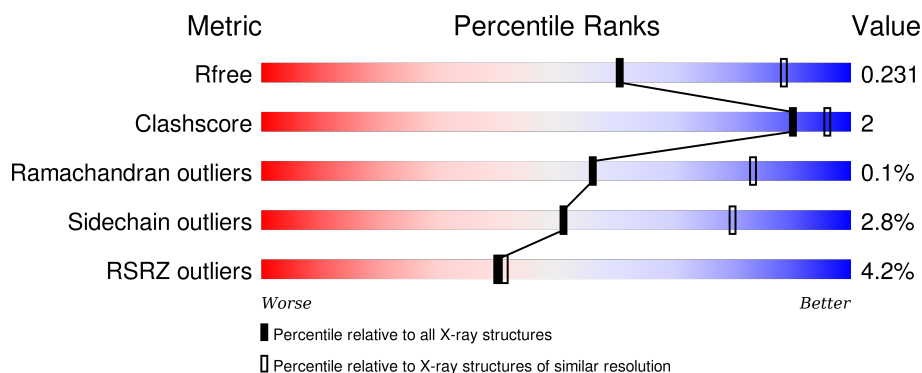
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 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	3050 (2.78-2.70)
Clashscore	102246	3424 (2.78-2.70)
Ramachandran outliers	100387	3367 (2.78-2.70)
Sidechain outliers	100360	3368 (2.78-2.70)
RSRZ outliers	91569	3055 (2.78-2.70)

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 $\leq 5\%$. 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	1113	<div> <div>3%</div> <div>88%</div> <div>7%</div> <div>5%</div> </div>
1	B	1113	<div> <div>5%</div> <div>82%</div> <div>7%</div> <div>11%</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 15902 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.

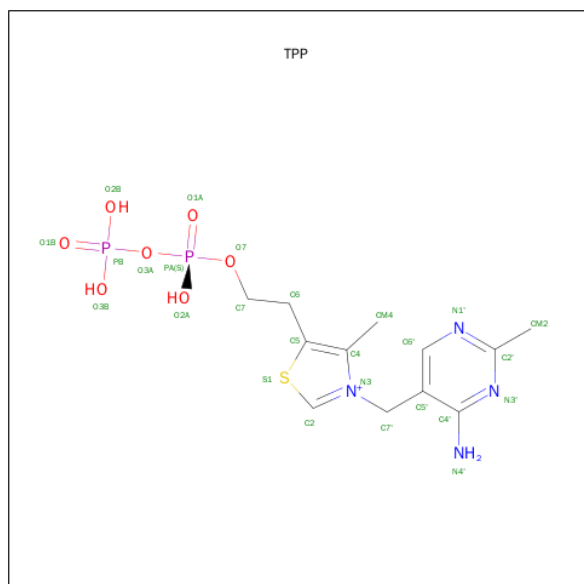
- Molecule 1 is a protein called 2-OXOGLUTARATE DECARBOXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1055	Total	C	N	O	S	0	0	0
			8153	5136	1447	1540	30			
1	B	989	Total	C	N	O	S	0	0	0
			7574	4776	1342	1429	27			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	115	GLY	-	EXPRESSION TAG	UNP A0R2B1
B	115	GLY	-	EXPRESSION TAG	UNP A0R2B1

- Molecule 2 is THIAMINE DIPHOSPHATE (three-letter code: TPP) (formula: C₁₂H₁₉N₄O₇P₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			26	12	4	7	2	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	S	
			26	12	4	7	2	1	
								0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg		
			1	1	0	0
3	A	1	Total	Mg		
			1	1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca		
			1	1	0	0
4	A	1	Total	Ca		
			1	1	0	0

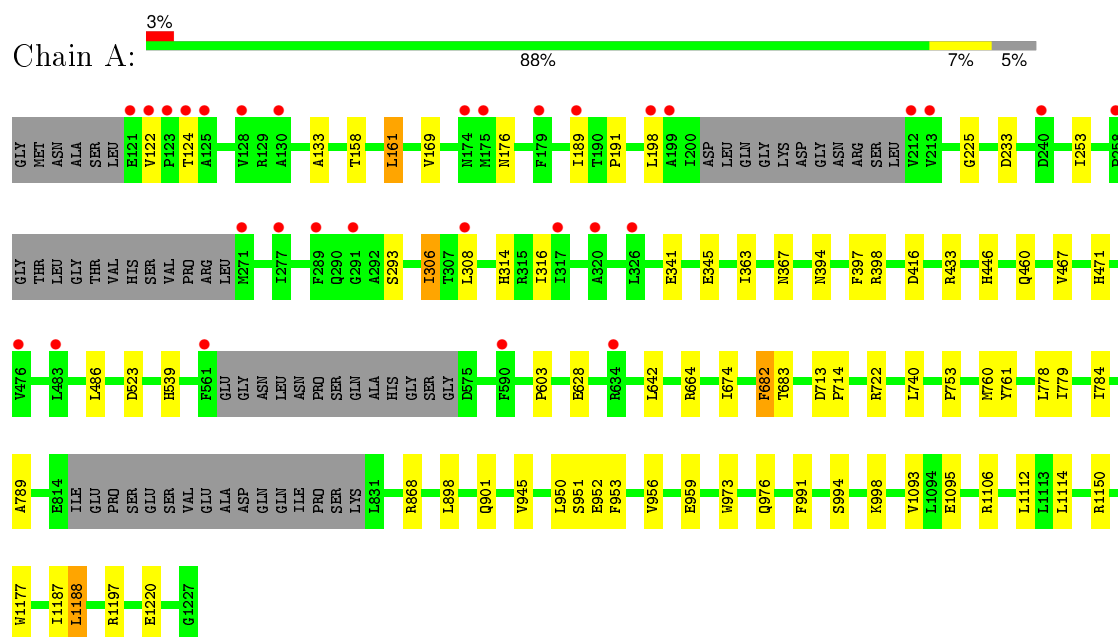
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	56	Total	O		
			56	56	0	0
5	B	63	Total	O		
			63	63	0	0

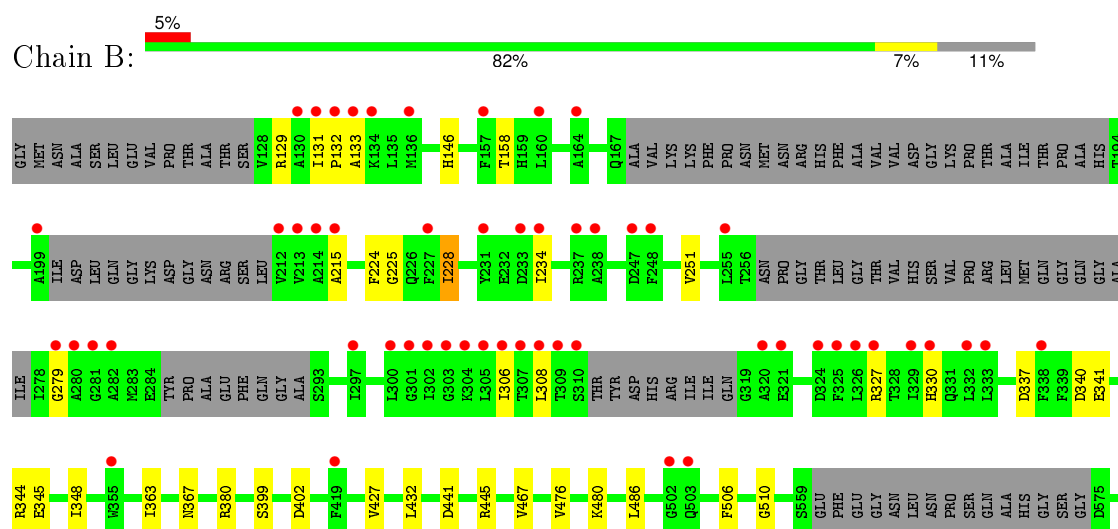
3 Residue-property plots

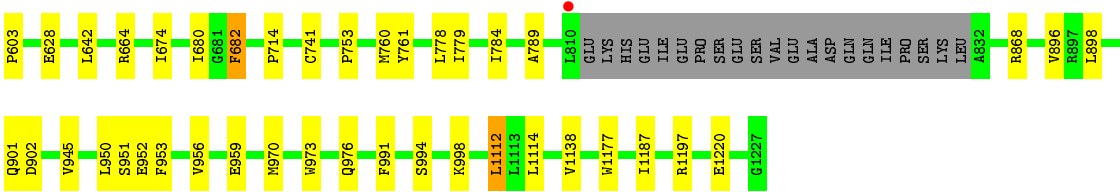
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: 2-OXOGLUTARATE DECARBOXYLASE



• Molecule 1: 2-OXOGLUTARATE DECARBOXYLASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	151.99 Å 247.72 Å 79.98 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.12 – 2.74 72.65 – 2.74	Depositor EDS
% Data completeness (in resolution range)	99.1 (40.12-2.74) 99.1 (72.65-2.74)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.73 Å)	Xtriage
Refinement program	BUSTER-TNT 2.9.3	Depositor
R, R_{free}	0.185 , 0.224 0.192 , 0.231	Depositor DCC
R_{free} test set	4017 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	57.9	Xtriage
Anisotropy	0.620	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 58.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 79783 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15902	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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: CA, MG, TPP

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	0.51	0/8319	0.67	0/11288
1	B	0.51	0/7721	0.67	0/10472
All	All	0.51	0/16040	0.67	0/21760

There are no bond length outliers.

There are no bond angle outliers.

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	8153	0	7892	35	0
1	B	7574	0	7289	38	0
2	A	26	0	16	1	0
2	B	26	0	16	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	56	0	0	0	0
5	B	63	0	0	0	0
All	All	15902	0	15213	71	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 2.

All (71) 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:779:ILE:HD11	1:B:789:ALA:HB2	1.49	0.94
1:A:779:ILE:HD11	1:A:789:ALA:HB2	1.51	0.92
1:A:642:LEU:HD22	1:A:674:ILE:HB	1.75	0.68
1:B:642:LEU:HD22	1:B:674:ILE:HB	1.75	0.68
1:A:1095:GLU:OE1	1:A:1106:ARG:NH1	2.30	0.64
1:B:778:LEU:HB3	1:B:784:ILE:HG12	1.80	0.64
1:A:778:LEU:HB3	1:A:784:ILE:HG12	1.80	0.63
1:A:779:ILE:HD11	1:A:789:ALA:CB	2.27	0.63
1:B:779:ILE:HD11	1:B:789:ALA:CB	2.26	0.63
1:A:416:ASP:CG	1:A:433:ARG:HH12	2.03	0.61
1:A:133:ALA:HB2	1:A:306:ILE:HG13	1.83	0.60
1:A:683:THR:HB	1:B:902:ASP:OD1	2.01	0.60
1:B:225:GLY:H	1:B:345:GLU:HG2	1.65	0.60
1:A:225:GLY:H	1:A:345:GLU:CG	2.18	0.57
1:B:952:GLU:O	1:B:956:VAL:HG23	2.06	0.56
1:A:952:GLU:O	1:A:956:VAL:HG23	2.06	0.55
1:A:523:ASP:OD2	1:A:722:ARG:NH2	2.39	0.55
1:A:225:GLY:H	1:A:345:GLU:HG2	1.72	0.54
1:B:133:ALA:CB	1:B:306:ILE:HD12	2.39	0.52
1:A:1093:VAL:HB	1:A:1150:ARG:HD3	1.92	0.51
1:B:1112:LEU:HD11	1:B:1138:VAL:HG23	1.92	0.51
1:B:132:PRO:HD2	1:B:330:HIS:NE2	2.26	0.51
1:A:603:PRO:HD3	1:A:991:PHE:CZ	2.46	0.51
1:B:603:PRO:HD3	1:B:991:PHE:CZ	2.46	0.51
1:B:680:ILE:HD12	1:B:760:MET:CE	2.41	0.51
1:B:131:ILE:HD12	1:B:308:LEU:HD12	1.93	0.50
1:A:161:LEU:HD21	1:A:308:LEU:HD21	1.92	0.50
1:B:133:ALA:HB3	1:B:306:ILE:HD12	1.93	0.50
1:B:441:ASP:HA	1:B:445:ARG:HG3	1.93	0.50
1:A:397:PHE:O	1:A:398:ARG:HD2	2.11	0.50
1:B:340:ASP:O	1:B:344:ARG:HG2	2.11	0.49
1:B:510:GLY:O	1:B:741:CYS:HB2	2.11	0.49
1:A:901:GLN:OE1	2:B:2001:TPP:H6'	2.13	0.49
1:B:680:ILE:HD12	1:B:760:MET:HE1	1.94	0.49
1:A:898:LEU:O	1:A:945:VAL:HA	2.13	0.48
1:B:215:ALA:HB2	1:B:251:VAL:HG22	1.95	0.48
1:A:959:GLU:HG3	1:A:973:TRP:HB2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:760:MET:HE3	1:B:761:TYR:CE2	2.49	0.48
1:B:279:GLY:O	1:B:308:LEU:HD23	2.13	0.48
1:B:898:LEU:O	1:B:945:VAL:HA	2.15	0.47
1:A:169:VAL:HG13	1:A:176:ASN:HB2	1.96	0.47
1:B:959:GLU:HG3	1:B:973:TRP:HB2	1.96	0.46
1:B:225:GLY:HA2	1:B:228:ILE:HD12	1.95	0.46
1:B:753:PRO:HB2	1:B:761:TYR:CE2	2.50	0.46
1:A:363:ILE:HD12	1:A:363:ILE:H	1.81	0.46
1:B:363:ILE:HD12	1:B:363:ILE:H	1.80	0.46
1:B:476:VAL:HG12	1:B:480:LYS:HD2	1.97	0.45
1:B:337:ASP:HA	1:B:340:ASP:HB2	1.98	0.45
1:A:460:GLN:NE2	1:B:380:ARG:HH11	2.14	0.44
1:A:446:HIS:ND1	1:A:713:ASP:OD2	2.51	0.44
1:A:994:SER:HB2	1:A:998:LYS:HE2	2.00	0.43
1:B:224:PHE:O	1:B:228:ILE:HG13	2.18	0.43
1:A:753:PRO:HB2	1:A:761:TYR:CE2	2.54	0.43
1:A:225:GLY:H	1:A:345:GLU:HG3	1.84	0.43
1:B:486:LEU:HD11	1:B:714:PRO:HG3	1.99	0.43
1:A:760:MET:HE3	1:A:761:TYR:CE2	2.54	0.43
1:A:486:LEU:HD11	1:A:714:PRO:HG3	2.01	0.42
1:A:161:LEU:HD11	1:A:308:LEU:HD11	2.00	0.42
1:B:628:GLU:HG2	1:B:664:ARG:O	2.19	0.42
2:A:2001:TPP:H6'	1:B:901:GLN:OE1	2.20	0.42
1:A:628:GLU:HG2	1:A:664:ARG:O	2.20	0.42
1:A:1177:TRP:CD1	1:A:1197:ARG:HD3	2.56	0.41
1:B:896:VAL:HG22	1:B:970:MET:HB3	2.02	0.41
1:B:129:ARG:NH2	1:B:327:ARG:HD3	2.34	0.41
1:A:539:HIS:H	1:A:539:HIS:CD2	2.37	0.41
1:B:994:SER:HB2	1:B:998:LYS:HE2	2.02	0.41
1:A:189:ILE:O	1:A:191:PRO:HD3	2.21	0.40
1:B:146:HIS:CD2	1:B:348:ILE:HG12	2.56	0.40
1:A:122:VAL:O	1:A:124:THR:HG23	2.21	0.40
1:B:1177:TRP:CD1	1:B:1197:ARG:HD3	2.57	0.40
1:A:1187:ILE:HG22	1:A:1188:LEU:HG	2.03	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	1045/1113 (94%)	1010 (97%)	34 (3%)	1 (0%)	56	83
1	B	973/1113 (87%)	938 (96%)	33 (3%)	2 (0%)	52	80
All	All	2018/2226 (91%)	1948 (96%)	67 (3%)	3 (0%)	56	83

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	399	SER
1	A	682	PHE
1	B	682	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	838/926 (90%)	813 (97%)	25 (3%)	48	78
1	B	767/926 (83%)	747 (97%)	20 (3%)	54	82
All	All	1605/1852 (87%)	1560 (97%)	45 (3%)	51	80

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	158	THR
1	A	161	LEU
1	A	198	LEU

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Mol	Chain	Res	Type
1	A	233	ASP
1	A	253	ILE
1	A	293	SER
1	A	306	ILE
1	A	314	HIS
1	A	316	ILE
1	A	341	GLU
1	A	367	ASN
1	A	394	ASN
1	A	467	VAL
1	A	471	HIS
1	A	682	PHE
1	A	740	LEU
1	A	868	ARG
1	A	950	LEU
1	A	951	SER
1	A	953	PHE
1	A	976	GLN
1	A	1112	LEU
1	A	1114	LEU
1	A	1188	LEU
1	A	1220	GLU
1	B	158	THR
1	B	228	ILE
1	B	234	ILE
1	B	341	GLU
1	B	367	ASN
1	B	402	ASP
1	B	427	VAL
1	B	432	LEU
1	B	467	VAL
1	B	506	PHE
1	B	682	PHE
1	B	868	ARG
1	B	950	LEU
1	B	951	SER
1	B	953	PHE
1	B	976	GLN
1	B	1112	LEU
1	B	1114	LEU
1	B	1187	ILE
1	B	1220	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are 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	$\# Z > 2$
2	TPP	A	2001	3	20,27,27	2.06	3 (15%)	31,40,40	1.92	9 (29%)
2	TPP	B	2001	3	20,27,27	2.30	4 (20%)	31,40,40	1.89	7 (22%)

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
2	TPP	A	2001	3	-	0/16/17/17	0/2/2/2
2	TPP	B	2001	3	-	0/16/17/17	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2001	TPP	C4-N3	-8.62	1.32	1.39
2	A	2001	TPP	C4-N3	-7.42	1.33	1.39
2	B	2001	TPP	PB-O2B	-2.01	1.47	1.54
2	A	2001	TPP	PB-O3B	2.07	1.62	1.54
2	B	2001	TPP	PB-O3B	2.49	1.63	1.54
2	B	2001	TPP	PB-O1B	2.98	1.61	1.51
2	A	2001	TPP	PB-O1B	3.21	1.61	1.51

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2001	TPP	CM4-C4-C5	-4.96	117.76	128.90
2	B	2001	TPP	CM4-C4-C5	-4.74	118.26	128.90
2	A	2001	TPP	N1'-C2'-N3'	-2.60	120.78	125.60
2	B	2001	TPP	N1'-C2'-N3'	-2.45	121.07	125.60
2	B	2001	TPP	C5'-C6'-N1'	-2.43	119.63	123.86
2	A	2001	TPP	C5'-C6'-N1'	-2.22	120.01	123.86
2	A	2001	TPP	CM4-C4-N3	2.04	125.31	122.59
2	A	2001	TPP	O3A-PA-O7	2.04	108.35	102.94
2	A	2001	TPP	C6-C5-C4	2.05	129.40	127.56
2	B	2001	TPP	CM4-C4-N3	2.27	125.62	122.59
2	B	2001	TPP	C6'-N1'-C2'	2.86	120.77	115.77
2	A	2001	TPP	C6'-N1'-C2'	3.19	121.34	115.77
2	B	2001	TPP	C5-C4-N3	3.83	116.12	107.69
2	A	2001	TPP	C5-C4-N3	4.17	116.86	107.69
2	A	2001	TPP	CM2-C2'-N1'	4.66	122.63	117.03
2	B	2001	TPP	CM2-C2'-N1'	5.06	123.10	117.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001	TPP	1	0
2	B	2001	TPP	1	0

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, 95th 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	1055/1113 (94%)	0.16	30 (2%) 56 59	39, 65, 102, 140	0
1	B	989/1113 (88%)	0.24	55 (5%) 28 27	41, 65, 140, 173	0
All	All	2044/2226 (91%)	0.20	85 (4%) 40 41	39, 65, 115, 173	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	214	ALA	8.0
1	B	302	ILE	7.8
1	B	303	GLY	7.3
1	B	282	ALA	7.2
1	B	136	MET	7.1
1	B	325	PHE	7.0
1	B	329	ILE	6.9
1	B	213	VAL	6.5
1	B	281	GLY	6.5
1	B	238	ALA	5.7
1	A	122	VAL	5.0
1	B	308	LEU	4.9
1	A	212	VAL	4.7
1	B	297	ILE	4.7
1	B	304	LYS	4.4
1	B	233	ASP	4.4
1	B	309	THR	4.4
1	B	333	LEU	4.2
1	B	247	ASP	4.2
1	B	327	ARG	4.2
1	B	164	ALA	4.2
1	B	306	ILE	4.1
1	B	307	THR	3.9
1	B	310	SER	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	133	ALA	3.7
1	B	255	LEU	3.7
1	B	280	ALA	3.6
1	B	810	LEU	3.5
1	B	231	TYR	3.5
1	B	503	GLN	3.4
1	A	198	LEU	3.4
1	A	174	ASN	3.4
1	B	212	VAL	3.3
1	B	130	ALA	3.3
1	B	234	ILE	3.3
1	B	132	PRO	3.3
1	A	240	ASP	3.3
1	A	258	PRO	3.2
1	A	590	PHE	3.2
1	A	121	GLU	3.2
1	B	237	ARG	3.1
1	B	134	LYS	3.0
1	B	300	LEU	3.0
1	A	213	VAL	2.9
1	A	476	VAL	2.9
1	A	291	GLY	2.9
1	A	289	PHE	2.8
1	A	123	PRO	2.7
1	B	215	ALA	2.7
1	A	634	ARG	2.7
1	A	125	ALA	2.7
1	A	179	PHE	2.6
1	B	355	TRP	2.6
1	A	128	VAL	2.6
1	A	130	ALA	2.5
1	A	561	PHE	2.5
1	A	326	LEU	2.5
1	B	502	GLY	2.5
1	A	308	LEU	2.4
1	B	248	PHE	2.4
1	A	271	MET	2.4
1	B	330	HIS	2.4
1	B	279	GLY	2.4
1	B	160	LEU	2.3
1	A	199	ALA	2.3
1	A	483	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	305	LEU	2.3
1	B	199	ALA	2.2
1	B	157	PHE	2.2
1	B	301	GLY	2.2
1	B	321	GLU	2.2
1	A	189	ILE	2.2
1	B	419	PHE	2.2
1	A	124	THR	2.2
1	A	317	ILE	2.1
1	B	131	ILE	2.1
1	B	332	LEU	2.1
1	B	326	LEU	2.1
1	A	277	ILE	2.1
1	B	227	PHE	2.1
1	B	338	PHE	2.1
1	B	324	ASP	2.1
1	A	320	ALA	2.1
1	A	175	MET	2.0
1	B	320	ALA	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, 95th 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
2	TPP	A	2001	26/26	0.99	0.18	0.08	35,51,66,74	0
4	CA	B	2003	1/1	0.99	0.14	-0.01	67,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	TPP	B	2001	26/26	0.99	0.18	-0.29	23,41,70,73	0
3	MG	B	2002	1/1	0.98	0.18	-0.51	32,32,32,32	0
3	MG	A	2002	1/1	0.98	0.14	-0.77	41,41,41,41	0
4	CA	A	2003	1/1	0.99	0.07	-2.32	72,72,72,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.