Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 07:56 AM GMT

PDB ID	:	3CQ4
Title	:	Histidinol-phosphate aminotransferase from Corynebacterium glutamicum
Authors	:	Sandalova, T.; Marienhagen, J.; Schneider, G.
Deposited on	:	2008-04-02
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:

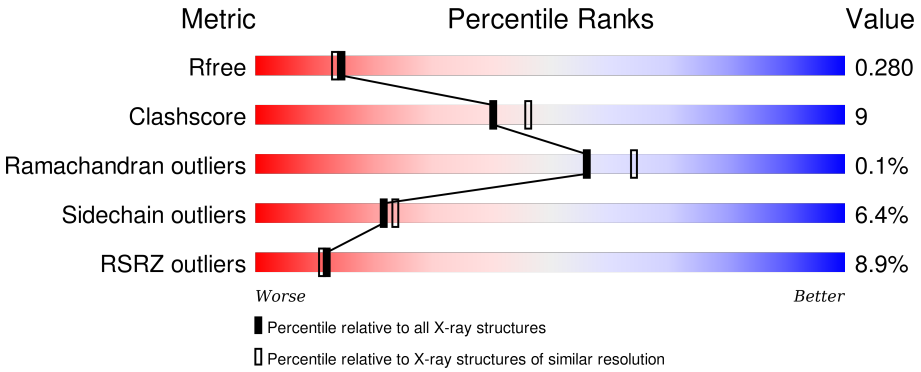
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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	376	5% 79% 13% • 5%
1	B	376	12% 77% 14% • 8%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5636 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.

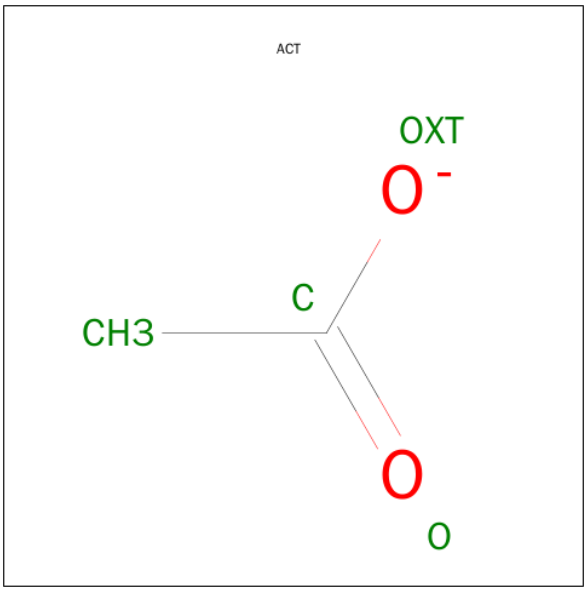
- Molecule 1 is a protein called Histidinol-phosphate aminotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	0	0	0
			2751	1750	470	527	4			
1	B	347	Total	C	N	O	S	0	0	0
			2660	1692	453	511	4			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	367	SER	-	EXPRESSION TAG	UNP Q9KJU4
A	368	ALA	-	EXPRESSION TAG	UNP Q9KJU4
A	369	TRP	-	EXPRESSION TAG	UNP Q9KJU4
A	370	SER	-	EXPRESSION TAG	UNP Q9KJU4
A	371	HIS	-	EXPRESSION TAG	UNP Q9KJU4
A	372	PRO	-	EXPRESSION TAG	UNP Q9KJU4
A	373	GLN	-	EXPRESSION TAG	UNP Q9KJU4
A	374	PHE	-	EXPRESSION TAG	UNP Q9KJU4
A	375	GLU	-	EXPRESSION TAG	UNP Q9KJU4
A	376	LYS	-	EXPRESSION TAG	UNP Q9KJU4
B	367	SER	-	EXPRESSION TAG	UNP Q9KJU4
B	368	ALA	-	EXPRESSION TAG	UNP Q9KJU4
B	369	TRP	-	EXPRESSION TAG	UNP Q9KJU4
B	370	SER	-	EXPRESSION TAG	UNP Q9KJU4
B	371	HIS	-	EXPRESSION TAG	UNP Q9KJU4
B	372	PRO	-	EXPRESSION TAG	UNP Q9KJU4
B	373	GLN	-	EXPRESSION TAG	UNP Q9KJU4
B	374	PHE	-	EXPRESSION TAG	UNP Q9KJU4
B	375	GLU	-	EXPRESSION TAG	UNP Q9KJU4
B	376	LYS	-	EXPRESSION TAG	UNP Q9KJU4

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 3 is water.

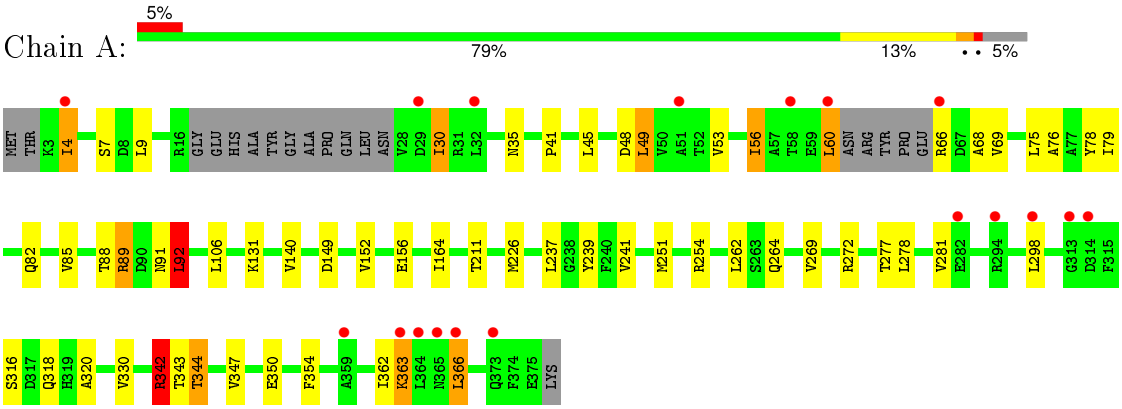
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	106	Total O 106 106	0	0
3	B	111	Total O 111 111	0	0

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.

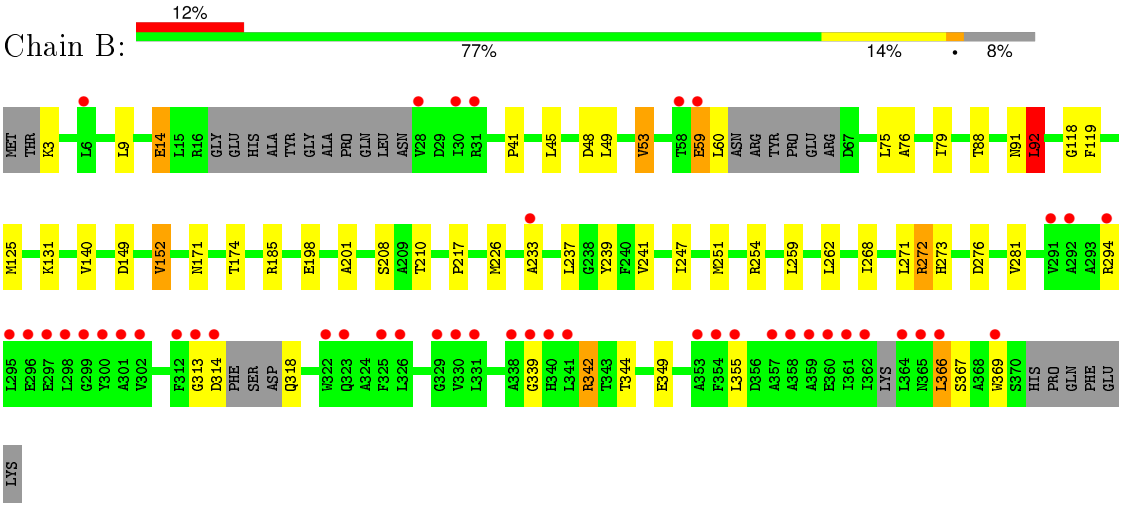
- Molecule 1: Histidinol-phosphate aminotransferase

Chain A:



- Molecule 1: Histidinol-phosphate aminotransferase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	102.34Å 102.34Å 140.11Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.20 54.96 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.3 (25.00-2.20) 99.3 (54.96-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) >$ [1]	4.09 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.237 , 0.277 0.239 , 0.280	Depositor DCC
R_{free} test set	2175 reflections (5.28%)	DCC
Wilson B-factor (Å^2)	36.6	Xtriage
Anisotropy	0.147	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 40.6	EDS
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 43385 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	5636	wwPDB-VP
Average B, all atoms (Å^2)	40.0	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.*

[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: ACT

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	0.56	0/2805	0.72	4/3822 (0.1%)
1	B	0.59	0/2708	0.72	3/3689 (0.1%)
All	All	0.57	0/5513	0.72	7/7511 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	342	ARG	NE-CZ-NH1	9.74	125.17	120.30
1	A	342	ARG	NE-CZ-NH2	-7.36	116.62	120.30
1	A	92	LEU	CA-CB-CG	-6.08	101.31	115.30
1	B	342	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	B	185	ARG	NE-CZ-NH1	5.21	122.91	120.30
1	B	92	LEU	CA-CB-CG	-5.05	103.69	115.30
1	A	342	ARG	CG-CD-NE	5.02	122.34	111.80

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	2751	0	2736	50	0
1	B	2660	0	2653	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	4	0	3	1	0
2	B	4	0	3	0	0
3	A	106	0	0	0	0
3	B	111	0	0	0	0
All	All	5636	0	5395	94	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 9.

All (94) 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:A:79:ILE:HD11	1:A:239:TYR:HE2	1.37	0.90
1:A:75:LEU:HD21	1:A:226:MET:CE	2.04	0.88
1:A:320:ALA:HB1	1:A:366:LEU:HD12	1.58	0.85
1:B:79:ILE:HD12	1:B:92:LEU:HD21	1.58	0.84
1:B:79:ILE:HD11	1:B:239:TYR:HE2	1.43	0.81
1:A:75:LEU:HD21	1:A:226:MET:HE1	1.65	0.77
1:B:49:LEU:O	1:B:53:VAL:HG12	1.88	0.73
1:A:78:TYR:CZ	1:A:82:GLN:HG3	2.24	0.72
1:A:79:ILE:HD13	1:A:241:VAL:CG2	2.21	0.71
1:B:79:ILE:HD13	1:B:241:VAL:CG2	2.20	0.70
1:B:171:ASN:ND2	1:B:174:THR:OG1	2.24	0.70
1:B:41:PRO:HB2	1:B:45:LEU:HD23	1.72	0.70
1:A:48:ASP:OD2	1:A:272:ARG:HD2	1.93	0.69
1:A:75:LEU:HD21	1:A:226:MET:HE2	1.76	0.68
1:A:79:ILE:HD12	1:A:92:LEU:HD21	1.76	0.67
1:B:79:ILE:CD1	1:B:92:LEU:HD21	2.24	0.67
1:B:79:ILE:HD13	1:B:241:VAL:HG21	1.74	0.67
1:B:79:ILE:HD11	1:B:239:TYR:CE2	2.29	0.66
1:A:79:ILE:HD13	1:A:241:VAL:HG21	1.78	0.65
1:A:298:LEU:O	1:A:362:ILE:HD13	1.96	0.65
1:B:247:ILE:O	1:B:251:MET:HG3	1.96	0.65
1:B:60:LEU:HD23	1:B:60:LEU:C	2.18	0.64
1:B:75:LEU:HD21	1:B:226:MET:HE2	1.78	0.64
1:B:226:MET:HE3	1:B:237:LEU:HG	1.79	0.64
1:B:342:ARG:HH11	1:B:342:ARG:HG2	1.63	0.63
1:A:79:ILE:HD11	1:A:239:TYR:CE2	2.26	0.63
1:B:48:ASP:OD2	1:B:273:HIS:HE1	1.82	0.62
1:A:88:THR:H	1:A:91:ASN:HD22	1.47	0.61
1:B:41:PRO:CB	1:B:45:LEU:HD23	2.31	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:ARG:HD3	1:B:355:LEU:HD13	1.86	0.58
1:A:342:ARG:HH11	1:A:342:ARG:HG2	1.69	0.56
1:A:41:PRO:CB	1:A:45:LEU:HD23	2.35	0.56
1:B:198:GLU:HB3	1:B:201:ALA:HB2	1.86	0.56
1:A:79:ILE:HD13	1:A:241:VAL:HG22	1.87	0.56
1:A:49:LEU:HD13	1:A:269:VAL:HG11	1.87	0.55
1:B:171:ASN:HD21	1:B:174:THR:HG23	1.72	0.55
1:A:53:VAL:HG11	1:B:53:VAL:HG11	1.88	0.55
1:A:69:VAL:HG13	1:A:89:ARG:NH1	2.21	0.54
1:B:171:ASN:ND2	1:B:174:THR:H	2.06	0.53
1:A:226:MET:CE	1:A:237:LEU:HD12	2.39	0.53
1:A:149:ASP:O	1:A:152:VAL:HG12	2.09	0.53
1:A:49:LEU:HD13	1:A:269:VAL:CG1	2.39	0.53
1:A:85:VAL:HG21	1:A:211:THR:HA	1.91	0.52
1:A:56:ILE:N	1:A:56:ILE:HD13	2.25	0.51
1:B:75:LEU:HD21	1:B:226:MET:CE	2.41	0.51
1:B:79:ILE:HD13	1:B:241:VAL:HG22	1.93	0.50
1:A:41:PRO:HB2	1:A:45:LEU:HD23	1.94	0.50
1:A:226:MET:HE3	1:A:237:LEU:HD12	1.94	0.49
1:B:49:LEU:HD21	1:B:233:ALA:HB2	1.93	0.49
1:A:277:THR:O	1:A:281:VAL:HG13	2.13	0.49
1:A:79:ILE:CD1	1:A:241:VAL:HG22	2.42	0.49
1:B:226:MET:HE3	1:B:237:LEU:CG	2.43	0.49
1:A:30:ILE:HD11	1:A:330:VAL:HG22	1.94	0.49
1:A:75:LEU:CB	1:A:92:LEU:HD22	2.44	0.48
1:B:149:ASP:OD1	1:B:152:VAL:HG12	2.13	0.48
1:A:76:ALA:HB2	1:A:92:LEU:HD13	1.95	0.48
1:A:79:ILE:CD1	1:A:241:VAL:CG2	2.90	0.47
1:A:69:VAL:CG1	1:A:89:ARG:NH1	2.77	0.47
1:B:59:GLU:HG3	1:B:59:GLU:O	2.13	0.47
1:B:88:THR:H	1:B:91:ASN:HD22	1.61	0.47
1:B:75:LEU:HB3	1:B:92:LEU:HD22	1.97	0.46
1:A:69:VAL:HG13	1:A:89:ARG:CZ	2.46	0.46
1:A:75:LEU:HB3	1:A:92:LEU:CD2	2.46	0.46
1:B:118:GLY:HA3	1:B:125:MET:SD	2.56	0.46
1:B:14:GLU:O	1:B:131:LYS:NZ	2.47	0.45
1:A:342:ARG:NH1	1:A:344:THR:HG22	2.32	0.45
1:B:149:ASP:HB3	1:B:152:VAL:CG1	2.47	0.45
1:A:30:ILE:CD1	1:A:330:VAL:HG22	2.46	0.45
1:B:268:ILE:O	1:B:272:ARG:HG2	2.17	0.45
1:B:366:LEU:HD12	1:B:367:SER:H	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:LEU:HD22	1:A:60:LEU:C	2.37	0.44
1:B:208:SER:OG	1:B:210:THR:HG23	2.17	0.44
1:B:119:PHE:HA	1:B:140:VAL:O	2.16	0.44
1:A:69:VAL:HG13	1:A:89:ARG:HD3	2.00	0.44
1:A:347:VAL:HG22	1:A:350:GLU:OE1	2.18	0.44
1:B:226:MET:HE1	1:B:271:LEU:HD11	2.00	0.43
1:B:149:ASP:HB3	1:B:152:VAL:HG13	2.00	0.43
1:B:226:MET:HE2	1:B:237:LEU:HD12	1.99	0.43
1:A:35:ASN:HB2	2:A:401:ACT:H2	2.00	0.43
1:A:4:ILE:HD13	1:B:217:PRO:HB2	2.01	0.43
1:B:313:GLY:O	1:B:314:ASP:HB2	2.19	0.43
1:A:278:LEU:O	1:A:281:VAL:HG22	2.18	0.43
1:A:106:LEU:HD23	1:A:164:ILE:CD1	2.49	0.43
1:B:75:LEU:CB	1:B:92:LEU:HD22	2.49	0.42
1:B:76:ALA:HB2	1:B:92:LEU:HD13	2.01	0.42
1:A:75:LEU:HB2	1:A:92:LEU:HD22	2.01	0.42
1:B:314:ASP:N	1:B:339:GLY:O	2.52	0.42
1:B:251:MET:HB3	1:B:251:MET:HE2	1.82	0.41
1:A:88:THR:H	1:A:91:ASN:ND2	2.17	0.41
1:A:68:ALA:HA	1:A:264:GLN:NE2	2.36	0.41
1:A:363:LYS:HE3	1:A:363:LYS:HA	2.02	0.41
1:A:140:VAL:HG22	1:A:156:GLU:HG2	2.02	0.41
1:B:149:ASP:CG	1:B:152:VAL:HG12	2.41	0.40
1:A:343:THR:HG21	1:A:354:PHE:CZ	2.56	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	351/376 (93%)	339 (97%)	12 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	337/376 (90%)	327 (97%)	9 (3%)	1 (0%)	46 50
All	All	688/752 (92%)	666 (97%)	21 (3%)	1 (0%)	56 64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	366	LEU

5.3.2 Protein sidechains (i)

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	293/308 (95%)	273 (93%)	20 (7%)	20 21
1	B	283/308 (92%)	266 (94%)	17 (6%)	24 26
All	All	576/616 (94%)	539 (94%)	37 (6%)	22 24

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE
1	A	7	SER
1	A	9	LEU
1	A	30	ILE
1	A	49	LEU
1	A	56	ILE
1	A	60	LEU
1	A	66	ARG
1	A	89	ARG
1	A	92	LEU
1	A	131	LYS
1	A	251	MET
1	A	254	ARG
1	A	262	LEU
1	A	316	SER

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Mol	Chain	Res	Type
1	A	318	GLN
1	A	342	ARG
1	A	344	THR
1	A	363	LYS
1	A	366	LEU
1	B	3	LYS
1	B	9	LEU
1	B	14	GLU
1	B	53	VAL
1	B	59	GLU
1	B	92	LEU
1	B	152	VAL
1	B	254	ARG
1	B	259	LEU
1	B	262	LEU
1	B	272	ARG
1	B	276	ASP
1	B	281	VAL
1	B	318	GLN
1	B	344	THR
1	B	349	GLU
1	B	369	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	ASN
1	A	99	ASN
1	A	103	GLN
1	A	188	ASN
1	A	264	GLN
1	A	273	HIS
1	A	319	HIS
1	A	371	HIS
1	B	82	GLN
1	B	91	ASN
1	B	99	ASN
1	B	103	GLN
1	B	171	ASN
1	B	264	GLN
1	B	273	HIS

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

2 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	$\#\|Z\| > 2$
2	ACT	A	401	-	1,3,3	0.91	0	0,3,3	0.00	-
2	ACT	B	401	-	1,3,3	0.85	0	0,3,3	0.00	-

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	ACT	A	401	-	-	0/0/0/0	0/0/0/0
2	ACT	B	401	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	ACT	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, 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	357/376 (94%)	0.42	18 (5%) 32 32	22, 35, 66, 75	0
1	B	347/376 (92%)	0.92	45 (12%) 5 4	21, 34, 67, 76	0
All	All	704/752 (93%)	0.67	63 (8%) 12 11	21, 35, 66, 76	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	358	ALA	8.9
1	B	361	ILE	8.3
1	A	366	LEU	7.6
1	B	298	LEU	6.9
1	B	353	ALA	6.6
1	B	357	ALA	6.3
1	B	329	GLY	5.9
1	B	331	LEU	5.9
1	B	299	GLY	5.8
1	B	355	LEU	5.7
1	B	313	GLY	5.2
1	B	300	TYR	5.1
1	B	294	ARG	5.0
1	B	292	ALA	4.8
1	B	339	GLY	4.5
1	B	354	PHE	4.4
1	B	314	ASP	4.4
1	A	313	GLY	4.0
1	B	295	LEU	4.0
1	A	294	ARG	3.9
1	B	312	PHE	3.8
1	A	60	LEU	3.8
1	B	364	LEU	3.7
1	B	301	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	364	LEU	3.6
1	B	291	VAL	3.5
1	B	341	LEU	3.5
1	B	330	VAL	3.5
1	B	30	ILE	3.5
1	B	302	VAL	3.4
1	B	58	THR	3.4
1	B	326	LEU	3.3
1	B	340	HIS	3.3
1	B	323	GLN	3.2
1	B	28	VAL	3.2
1	A	51	ALA	3.1
1	A	363	LYS	3.1
1	B	360	GLU	3.0
1	B	59	GLU	2.9
1	B	322	TRP	2.9
1	B	362	ILE	2.8
1	A	32	LEU	2.7
1	A	314	ASP	2.7
1	B	359	ALA	2.5
1	B	297	GLU	2.5
1	A	29	ASP	2.4
1	B	325	PHE	2.4
1	B	369	TRP	2.4
1	A	4	ILE	2.4
1	A	58	THR	2.3
1	A	373	GLN	2.2
1	A	365	ASN	2.2
1	A	66	ARG	2.2
1	A	359	ALA	2.2
1	B	366	LEU	2.2
1	B	365	ASN	2.1
1	B	6	LEU	2.1
1	B	338	ALA	2.1
1	B	233	ALA	2.1
1	B	31	ARG	2.1
1	B	296	GLU	2.0
1	A	282	GLU	2.0
1	A	298	LEU	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^{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
2	ACT	B	401	4/4	0.93	0.20	0.27	62,63,63,63	0
2	ACT	A	401	4/4	0.85	0.13	-0.96	48,49,49,50	0

6.5 Other polymers (i)

There are no such residues in this entry.