



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 02:09 pm GMT

PDB ID : 3H1Q
Title : Crystal structure of ethanolamine utilization protein EutJ from *Carboxydothermus hydrogenoformans*
Authors : Chang, C.; Tesar, C.; Jedrzejczak, R.; Kinney, J.; Kerfeld, C.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2009-04-13
Resolution : 2.80 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

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

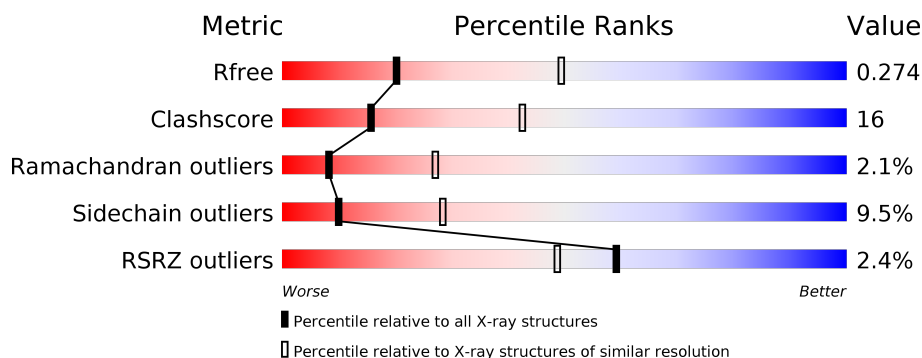
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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 ≥ 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	272	<div> <div>4%</div> <div> <div></div> <div>56%</div> <div>38%</div> <div>5% •</div> </div> </div>
1	B	272	<div> <div>75%</div> <div>21%</div> <div>•</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3978 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 Ethanolamine utilization protein EutJ.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	Se	0	0	0
			1912	1231	322	356	1	2			
1	B	272	Total	C	N	O	S	Se	0	0	0
			1977	1272	330	371	1	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q3AE93
A	-1	ASN	-	EXPRESSION TAG	UNP Q3AE93
A	0	ALA	-	EXPRESSION TAG	UNP Q3AE93
A	201	ARG	PRO	SEE REMARK 999	UNP Q3AE93
B	-2	SER	-	EXPRESSION TAG	UNP Q3AE93
B	-1	ASN	-	EXPRESSION TAG	UNP Q3AE93
B	0	ALA	-	EXPRESSION TAG	UNP Q3AE93
B	201	ARG	PRO	SEE REMARK 999	UNP Q3AE93

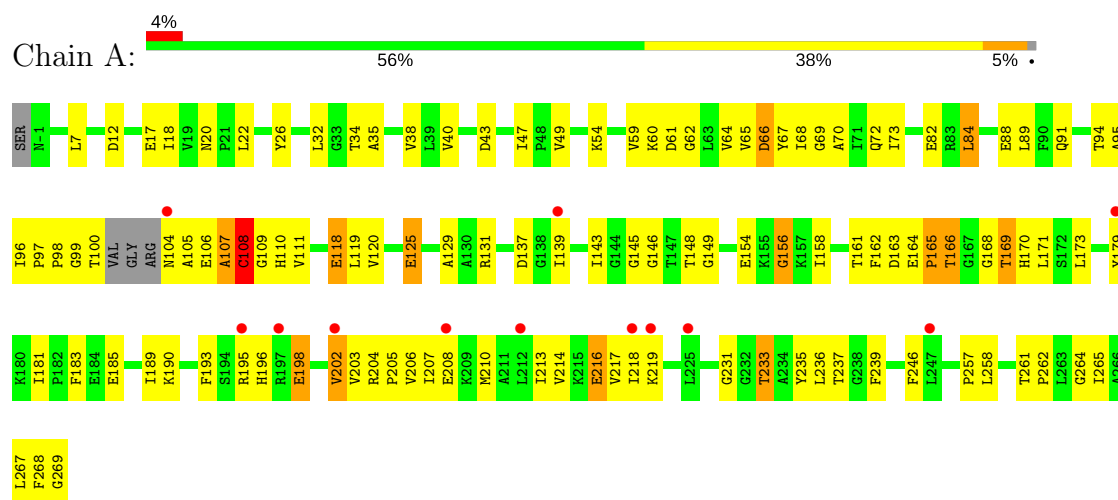
- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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: Ethanolamine utilization protein EutJ



• Molecule 1: Ethanolamine utilization protein EutJ



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	128.91Å 128.91Å 104.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 48.59 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-2.80) 99.5 (48.59-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.57 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0054	Depositor
R, R_{free}	0.227 , 0.278 0.227 , 0.274	Depositor DCC
R_{free} test set	1136 reflections (5.40%)	DCC
Wilson B-factor (Å ²)	60.2	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 80.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3978	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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, 0.333 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: ATP

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	0.71	1/1942 (0.1%)	0.79	1/2649 (0.0%)
1	B	0.75	0/2007	0.82	0/2737
All	All	0.73	1/3949 (0.0%)	0.80	1/5386 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	108	CYS	CB-SG	-6.17	1.71	1.82

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	ASP	CB-CG-OD1	5.07	122.86	118.30

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	1912	0	1888	88	0
1	B	1977	0	2002	42	0
2	A	31	0	12	2	0
2	B	30	0	11	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	11	0	0	2	0
3	B	17	0	0	0	0
All	All	3978	0	3913	129	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 16.

All (129) 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:34:THR:HG21	1:A:146:GLY:HA3	1.53	0.90
1:B:182:PRO:HD2	1:B:185:GLU:OE1	1.77	0.85
1:B:30:VAL:HG22	1:B:39:LEU:HD12	1.61	0.82
2:B:301:ATP:N3	2:B:301:ATP:N1	2.22	0.81
1:B:71:ILE:HD13	1:B:114:GLY:HA3	1.60	0.81
1:B:37:ILE:HD11	1:B:58:VAL:HG11	1.63	0.80
1:A:166:THR:HA	1:A:170:HIS:CE1	2.19	0.78
1:B:171:LEU:HD21	1:B:207:ILE:CD1	2.19	0.73
1:A:204:ARG:N	1:A:205:PRO:HD2	2.05	0.71
1:A:100:THR:OG1	1:A:105:ALA:HB2	1.92	0.70
1:B:58:VAL:HG23	1:B:59:VAL:HG23	1.71	0.69
1:A:95:ALA:O	1:A:97:PRO:HD3	1.93	0.69
1:A:100:THR:OG1	1:A:105:ALA:CB	2.40	0.69
1:A:54:LYS:HG2	1:A:73:ILE:HD12	1.76	0.68
1:A:120:VAL:HG11	1:A:269:GLY:HA2	1.77	0.67
1:A:181:ILE:CB	1:A:185:GLU:OE2	2.43	0.67
1:B:143:ILE:HA	1:B:148:THR:HB	1.77	0.66
1:A:17:GLU:HB3	1:A:84:LEU:HD21	1.78	0.65
1:A:213:ILE:O	1:A:217:VAL:HG23	1.98	0.64
1:B:95:ALA:HB2	1:B:265:ILE:CD1	2.27	0.64
1:B:113:ALA:HA	1:B:117:LEU:O	1.98	0.63
1:A:34:THR:HG21	1:A:146:GLY:CA	2.26	0.63
1:A:131:ARG:NH1	1:A:268:PHE:CD1	2.68	0.62
1:B:171:LEU:HD21	1:B:207:ILE:HD11	1.81	0.61
1:A:67:TYR:HD2	1:A:68:ILE:HD13	1.64	0.61
1:B:166:THR:HB	1:B:209:LYS:HZ3	1.65	0.61
1:A:95:ALA:C	1:A:97:PRO:HD3	2.21	0.60
1:B:264:GLY:O	1:B:267:LEU:HB2	2.03	0.59
1:B:264:GLY:HA2	1:B:267:LEU:HD12	1.85	0.59
1:B:204:ARG:HB3	1:B:205:PRO:HD3	1.86	0.58
1:B:243:PHE:HD2	1:B:251:VAL:HG21	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:GLU:C	1:A:156:GLY:H	2.08	0.57
1:A:38:VAL:HG12	1:A:40:VAL:HG23	1.86	0.56
1:A:173:LEU:HG	1:A:183:PHE:HE1	1.70	0.55
1:A:106:GLU:C	1:A:108:CYS:N	2.59	0.55
1:A:139:ILE:HD13	1:A:218:ILE:HD12	1.88	0.55
1:A:166:THR:HA	1:A:170:HIS:HE1	1.68	0.55
1:A:214:VAL:O	1:A:218:ILE:HG12	2.06	0.55
1:B:170:HIS:CD2	1:B:209:LYS:NZ	2.75	0.55
1:A:98:PRO:CB	3:A:408:HOH:O	2.55	0.55
1:A:206:VAL:O	1:A:210:MSE:HG3	2.07	0.54
1:A:109:GLY:HA3	1:A:119:LEU:CD2	2.39	0.53
1:B:91:GLN:HB3	1:B:120:VAL:HG23	1.90	0.53
1:A:67:TYR:CD2	1:A:68:ILE:HD13	2.43	0.53
1:A:60:LYS:HB2	1:A:65:VAL:HG21	1.91	0.53
1:A:261:THR:HB	1:A:262:PRO:HD3	1.90	0.53
1:A:12:ASP:OD2	1:B:18:ILE:HD13	2.09	0.53
1:A:82:GLU:HG2	1:A:88:GLU:HA	1.91	0.53
1:A:168:GLY:HA2	1:A:233:THR:HG23	1.91	0.53
1:A:169:THR:HG23	1:A:183:PHE:HZ	1.74	0.52
1:A:35:ALA:N	2:A:301:ATP:O1B	2.42	0.52
1:A:193:PHE:HA	1:A:196:HIS:CD2	2.45	0.52
1:A:20:ASN:HB3	1:A:49:VAL:HA	1.92	0.51
1:B:179:TYR:OH	1:B:198:GLU:OE2	2.18	0.51
1:A:95:ALA:O	1:A:108:CYS:SG	2.69	0.51
1:A:143:ILE:HG22	1:A:233:THR:OG1	2.11	0.51
1:A:32:LEU:HD12	1:A:108:CYS:SG	2.51	0.51
1:B:1:MSE:O	1:B:5:GLN:HG3	2.09	0.51
1:B:165:PRO:HD2	1:B:166:THR:HG23	1.92	0.50
1:B:106:GLU:HG2	1:B:122:LEU:HD11	1.93	0.50
1:A:173:LEU:HG	1:A:183:PHE:CE1	2.47	0.50
1:A:109:GLY:HA3	1:A:119:LEU:HD21	1.94	0.50
1:A:171:LEU:HD11	1:A:233:THR:HG22	1.92	0.50
1:A:216:GLU:O	1:A:219:LYS:HB3	2.11	0.50
1:B:20:ASN:OD1	1:B:22:LEU:HD13	2.12	0.50
1:A:107:ALA:O	1:A:111:VAL:HG23	2.12	0.50
1:A:66:ASP:OD2	1:A:69:GLY:HA3	2.11	0.50
1:A:164:GLU:O	1:A:166:THR:N	2.39	0.49
1:A:236:LEU:HB3	1:A:239:PHE:HB2	1.93	0.49
1:A:206:VAL:HG12	1:A:210:MSE:HE3	1.94	0.49
1:A:154:GLU:O	1:A:156:GLY:N	2.33	0.49
1:A:106:GLU:C	1:A:108:CYS:H	2.15	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:GLY:HA3	1:A:99:GLY:O	2.13	0.49
1:A:98:PRO:HB3	3:A:408:HOH:O	2.13	0.49
1:A:202:VAL:HG22	1:A:203:VAL:HG23	1.93	0.49
1:B:91:GLN:HB3	1:B:120:VAL:CG2	2.43	0.48
1:A:43:ASP:OD1	1:A:47:ILE:N	2.45	0.48
1:B:170:HIS:CD2	1:B:209:LYS:HZ2	2.31	0.48
1:A:70:ALA:HB1	1:A:111:VAL:HG11	1.95	0.48
1:A:179:TYR:OH	1:A:198:GLU:OE1	2.27	0.48
1:A:60:LYS:O	1:A:61:ASP:HB2	2.13	0.48
1:A:98:PRO:HD3	1:A:125:GLU:HG3	1.95	0.48
1:B:95:ALA:HB2	1:B:265:ILE:HD13	1.96	0.47
1:A:34:THR:HB	2:A:301:ATP:O1B	2.14	0.47
1:B:71:ILE:O	1:B:75:ARG:HB2	2.16	0.46
1:A:143:ILE:HG12	1:A:148:THR:HG23	1.96	0.46
1:B:185:GLU:O	1:B:189:ILE:HG12	2.15	0.46
1:A:161:THR:C	1:A:162:PHE:CD2	2.88	0.46
1:A:231:GLY:O	1:A:257:PRO:HB2	2.16	0.46
1:A:97:PRO:HA	1:A:98:PRO:HD3	1.81	0.46
1:B:78:LYS:HE2	1:B:89:LEU:HB2	1.99	0.45
1:A:109:GLY:CA	1:A:119:LEU:HD22	2.46	0.45
1:B:166:THR:HB	1:B:209:LYS:NZ	2.30	0.45
1:B:223:GLN:HA	1:B:223:GLN:NE2	2.32	0.44
1:A:204:ARG:N	1:A:205:PRO:CD	2.79	0.44
1:B:212:LEU:O	1:B:213:ILE:C	2.56	0.44
1:A:189:ILE:O	1:A:190:LYS:C	2.56	0.44
1:A:20:ASN:O	1:A:22:LEU:N	2.48	0.44
1:A:145:GLY:O	1:A:169:THR:OG1	2.31	0.44
1:A:148:THR:N	1:A:166:THR:O	2.51	0.44
1:B:190:LYS:NZ	2:B:301:ATP:H1'	2.32	0.44
1:A:219:LYS:HG2	1:A:219:LYS:O	2.18	0.44
1:A:189:ILE:HD12	1:A:195:ARG:NH1	2.33	0.44
1:A:109:GLY:CA	1:A:119:LEU:CD2	2.96	0.43
1:A:258:LEU:O	1:A:258:LEU:HD23	2.18	0.43
1:A:208:GLU:HG2	1:A:246:PHE:CZ	2.54	0.43
1:A:12:ASP:OD2	1:B:18:ILE:CD1	2.67	0.43
1:B:153:ILE:HD12	1:B:153:ILE:N	2.34	0.43
1:B:212:LEU:O	1:B:216:GLU:HG2	2.19	0.43
1:B:58:VAL:HG23	1:B:59:VAL:CG2	2.46	0.43
1:A:171:LEU:HD23	1:A:206:VAL:HB	2.01	0.42
1:A:22:LEU:O	1:A:26:TYR:OH	2.31	0.42
1:A:164:GLU:HA	1:A:165:PRO:HD3	1.74	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:LEU:HA	1:A:236:LEU:HD12	1.88	0.42
1:B:214:VAL:O	1:B:218:ILE:HG12	2.20	0.42
1:A:148:THR:HG22	1:A:149:GLY:H	1.84	0.42
1:A:149:GLY:HA2	1:A:163:ASP:HA	2.01	0.42
1:B:55:TRP:O	1:B:56:ALA:HB2	2.20	0.42
1:A:95:ALA:HB2	1:A:265:ILE:HD11	2.02	0.42
1:B:208:GLU:HG3	1:B:246:PHE:CE1	2.55	0.42
1:B:150:ILE:HB	1:B:162:PHE:HB2	2.02	0.41
1:A:154:GLU:C	1:A:156:GLY:N	2.73	0.41
1:A:264:GLY:O	1:A:267:LEU:HB2	2.21	0.41
1:A:91:GLN:HB2	1:A:118:GLU:O	2.21	0.41
1:A:107:ALA:HA	1:A:110:HIS:HB2	2.03	0.41
1:A:59:VAL:CG1	1:A:62:GLY:HA2	2.51	0.41
1:A:162:PHE:O	1:A:163:ASP:HB3	2.21	0.40
1:B:38:VAL:HG12	1:B:40:VAL:HG23	2.02	0.40
1:A:171:LEU:HD11	1:A:233:THR:CG2	2.52	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	264/272 (97%)	218 (83%)	38 (14%)	8 (3%)	5	17
1	B	270/272 (99%)	247 (92%)	20 (7%)	3 (1%)	17	47
All	All	534/544 (98%)	465 (87%)	58 (11%)	11 (2%)	8	27

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	107	ALA
1	A	108	CYS

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Mol	Chain	Res	Type
1	A	165	PRO
1	A	96	ILE
1	B	178	SER
1	A	129	ALA
1	B	56	ALA
1	A	64	VAL
1	B	65	VAL
1	A	156	GLY
1	A	158	ILE

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	186/216 (86%)	166 (89%)	20 (11%)	7	22
1	B	202/216 (94%)	185 (92%)	17 (8%)	13	35
All	All	388/432 (90%)	351 (90%)	37 (10%)	10	28

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

Mol	Chain	Res	Type
1	A	7	LEU
1	A	18	ILE
1	A	72	GLN
1	A	84	LEU
1	A	89	LEU
1	A	94	THR
1	A	104	ASN
1	A	108	CYS
1	A	118	GLU
1	A	125	GLU
1	A	137	ASP
1	A	166	THR
1	A	169	THR
1	A	198	GLU

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Mol	Chain	Res	Type
1	A	202	VAL
1	A	207	ILE
1	A	216	GLU
1	A	233	THR
1	A	235	TYR
1	A	237	THR
1	B	1	MSE
1	B	16	ARG
1	B	18	ILE
1	B	31	ASP
1	B	36	ASP
1	B	45	GLU
1	B	106	GLU
1	B	135	ILE
1	B	143	ILE
1	B	148	THR
1	B	166	THR
1	B	178	SER
1	B	199	ILE
1	B	203	VAL
1	B	223	GLN
1	B	235	TYR
1	B	265	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	GLN
1	B	44	GLN
1	B	170	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	ATP	A	301	-	27,33,33	1.12	1 (3%)	25,52,52	1.89	3 (12%)
2	ATP	B	301	-	22,31,33	1.43	2 (9%)	18,49,52	1.18	1 (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
2	ATP	A	301	-	-	0/18/38/38	0/3/3/3
2	ATP	B	301	-	-	0/18/38/38	0/2/2/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	ATP	C5-C6	-4.49	1.44	1.53
2	A	301	ATP	C5-C4	3.53	1.48	1.40
2	B	301	ATP	PG-O3B	3.81	1.66	1.60

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	ATP	N3-C2-N1	-7.51	122.31	128.86
2	A	301	ATP	C4-C5-N7	-3.23	106.29	109.41
2	B	301	ATP	O5'-C5'-C4'	-2.58	99.84	109.00
2	A	301	ATP	C2-N1-C6	2.38	122.93	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	ATP	2	0
2	B	301	ATP	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	265/272 (97%)	0.16	12 (4%) 34 24	30, 49, 63, 66	0
1	B	269/272 (98%)	-0.04	1 (0%) 92 90	26, 40, 52, 61	0
All	All	534/544 (98%)	0.06	13 (2%) 59 49	26, 43, 61, 66	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	104	ASN	4.5
1	A	208	GLU	3.1
1	A	218	ILE	2.5
1	A	197	ARG	2.4
1	A	219	LYS	2.4
1	A	195	ARG	2.4
1	A	212	LEU	2.4
1	A	225	LEU	2.4
1	A	179	TYR	2.4
1	A	247	LEU	2.2
1	A	139	ILE	2.1
1	A	202	VAL	2.1
1	B	115	ALA	2.1

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	ATP	B	301	30/31	0.95	0.20	0.75	45,52,62,65	0
2	ATP	A	301	31/31	0.94	0.18	0.48	55,66,79,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.