



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 2, 2017 – 09:37 AM EDT

PDB ID : 1Y0G
Title : CRYSTAL STRUCTURE OF THE ESCHERICHIA COLI YCEI PROTEIN,
STRUCTURAL GENOMICS
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New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : unknown
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 ⓘ 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 : rb-20030345
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) : rb-20030345

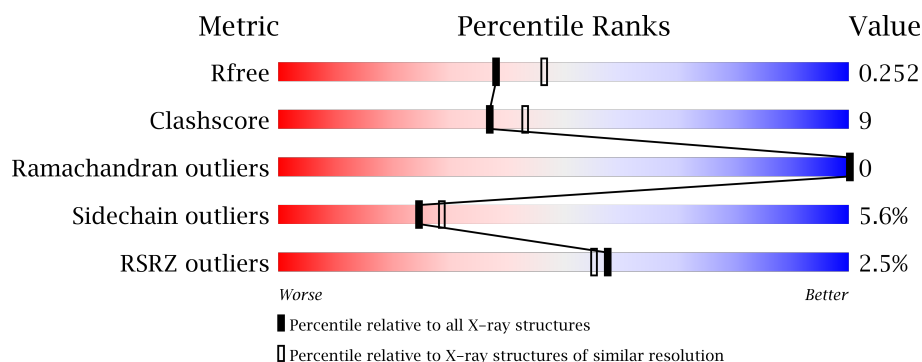
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (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 $\leq 5\%$. 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	191	<div> <div>4%</div> <div> <div></div> <div>73%</div> <div>13%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	191	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>15%</div> <div>•</div> <div>12%</div> </div> </div>
1	C	191	<div> <div>%</div> <div> <div></div> <div>71%</div> <div>16%</div> <div>•</div> <div>12%</div> </div> </div>
1	D	191	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>13%</div> <div>•</div> <div>12%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	8PP	A	320	-	-	-	X
2	8PP	B	321	-	-	-	X
2	8PP	C	322	-	-	-	X
2	8PP	D	323	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5796 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 yceI.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	169	Total	C	N	O	0	0	0
			1323	834	223	266			
1	B	169	Total	C	N	O	0	0	0
			1323	834	223	266			
1	C	169	Total	C	N	O	0	0	0
			1323	834	223	266			
1	D	169	Total	C	N	O	0	0	0
			1323	834	223	266			

- Molecule 2 is 2-[(2E,6E,10E,14E,18E,22E,26E)-3,7,11,15,19,23,27,31-OCTAMETHYL DOTRIACONTA-2,6,10,14,18,22,26,30-OCTAENYL]PHENOL (three-letter code: 8PP) (formula: C₄₆H₇₀O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			47	46	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			47	46	1		
2	C	1	Total	C	O	0	0
			47	46	1		
2	D	1	Total	C	O	0	0
			47	46	1		

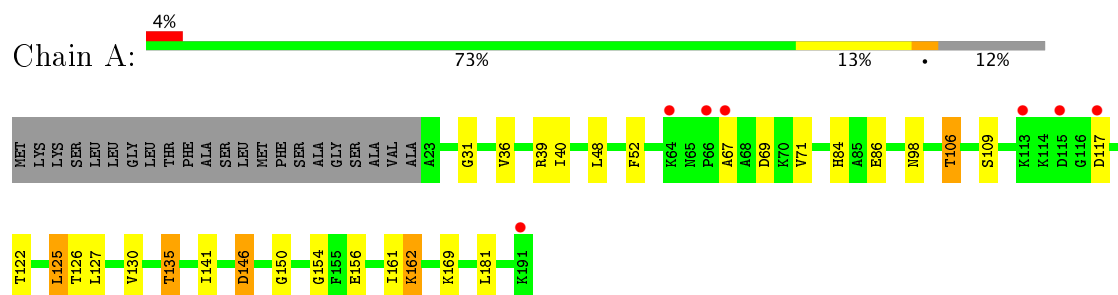
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	78	Total	O	0	0
			78	78		
3	B	88	Total	O	0	0
			88	88		
3	C	80	Total	O	0	0
			80	80		
3	D	70	Total	O	0	0
			70	70		

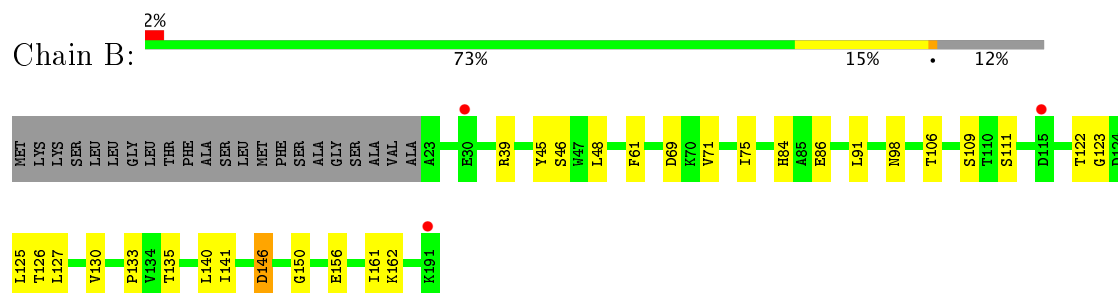
3 Residue-property plots [i](#)

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.

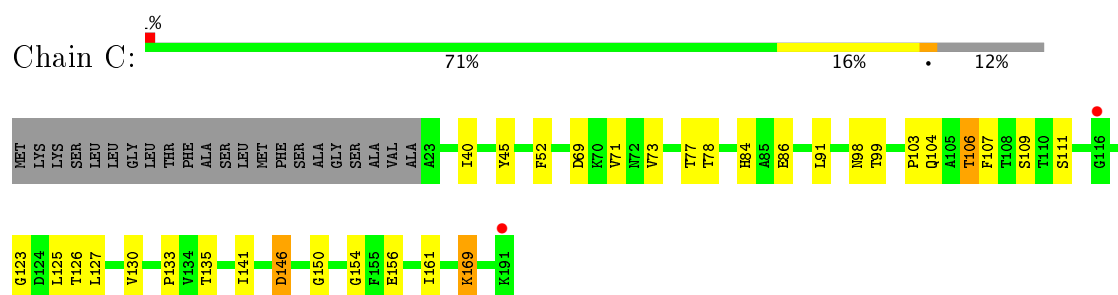
• Molecule 1: Protein yceI



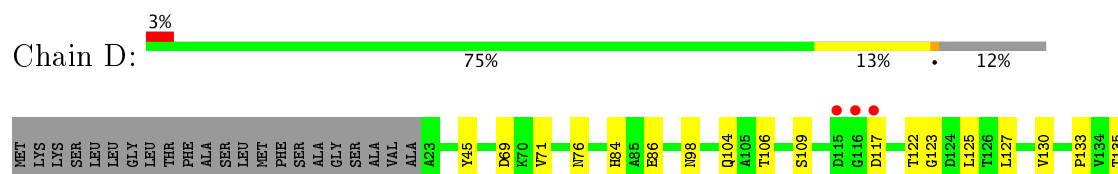
• Molecule 1: Protein yceI

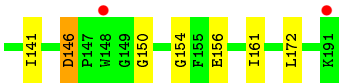


• Molecule 1: Protein yceI



• Molecule 1: Protein yceI





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	188.17Å 41.93Å 130.25Å 90.00° 126.33° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 19.96 – 2.20	Depositor EDS
% Data completeness (in resolution range)	91.1 (20.00-2.20) 96.4 (19.96-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.19Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.228 , 0.248 0.233 , 0.252	Depositor DCC
R_{free} test set	1234 reflections (3.03%)	DCC
Wilson B-factor (Å ²)	31.5	Xtriage
Anisotropy	0.324	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5796	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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: 8PP

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	0.41	0/1349	0.71	0/1822
1	B	0.42	0/1349	0.71	0/1822
1	C	0.41	0/1349	0.68	0/1822
1	D	0.43	0/1349	0.68	0/1822
All	All	0.42	0/5396	0.69	0/7288

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 [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	1323	0	1284	27	0
1	B	1323	0	1284	23	0
1	C	1323	0	1284	28	0
1	D	1323	0	1284	22	0
2	A	47	0	69	8	0
2	B	47	0	69	8	0
2	C	47	0	69	6	0
2	D	47	0	69	10	0
3	A	78	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	88	0	0	2	0
3	C	80	0	0	5	0
3	D	70	0	0	0	0
All	All	5796	0	5412	97	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 (97) 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:D:172:LEU:HD22	2:D:323:8PP:H103	1.37	1.01
1:D:172:LEU:HD13	2:D:323:8PP:H101	1.56	0.84
1:B:71:VAL:HG12	1:B:109:SER:HB3	1.62	0.78
1:C:40:ILE:HG21	2:C:322:8PP:H13	1.65	0.77
1:A:67:ALA:HB1	1:C:104:GLN:OE1	1.89	0.72
1:D:71:VAL:HG12	1:D:109:SER:HB3	1.73	0.71
1:B:106:THR:HG22	1:B:126:THR:HB	1.71	0.71
1:C:91:LEU:HD21	2:C:322:8PP:H122	1.74	0.70
1:B:106:THR:HG23	3:B:362:HOH:O	1.93	0.69
1:D:69:ASP:HB3	2:D:323:8PP:H451	1.76	0.68
1:D:172:LEU:CD2	2:D:323:8PP:H103	2.21	0.67
1:D:84:HIS:CE1	1:D:86:GLU:HB3	2.31	0.65
1:B:146:ASP:HB2	1:B:150:GLY:H	1.60	0.65
1:C:71:VAL:HG12	1:C:109:SER:HB3	1.76	0.65
1:A:146:ASP:HB2	1:A:150:GLY:H	1.62	0.65
1:A:146:ASP:CB	1:A:150:GLY:H	2.11	0.63
1:D:172:LEU:HD13	2:D:323:8PP:C10	2.28	0.62
1:A:146:ASP:HB2	1:A:150:GLY:O	1.99	0.61
1:A:106:THR:HG22	1:A:126:THR:HB	1.83	0.61
1:B:69:ASP:HB3	2:B:321:8PP:H451	1.83	0.61
1:C:78:THR:HG22	3:C:357:HOH:O	2.00	0.60
1:B:84:HIS:CE1	1:B:86:GLU:HB2	2.37	0.60
1:A:84:HIS:CE1	1:A:86:GLU:HB2	2.36	0.59
1:B:146:ASP:CB	1:B:150:GLY:H	2.16	0.58
1:C:69:ASP:HB3	2:C:322:8PP:H451	1.86	0.57
1:D:141:ILE:HD11	1:D:156:GLU:HB3	1.86	0.56
1:C:154:GLY:HA3	1:D:45:TYR:O	2.04	0.56
1:B:71:VAL:HG21	2:B:321:8PP:H38	1.88	0.56
1:D:71:VAL:HG21	2:D:323:8PP:H38	1.87	0.56
1:C:146:ASP:CB	1:C:150:GLY:H	2.19	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:146:ASP:HB2	1:C:150:GLY:H	1.70	0.55
1:A:40:ILE:HG21	2:A:320:8PP:H13	1.88	0.55
1:C:45:TYR:O	1:D:154:GLY:HA3	2.06	0.55
1:C:52:PHE:CE1	2:C:322:8PP:H222	2.42	0.54
1:A:161:ILE:CD1	2:A:320:8PP:H23	2.38	0.53
1:B:161:ILE:CD1	2:B:321:8PP:H23	2.39	0.52
1:D:161:ILE:CD1	2:D:323:8PP:H23	2.39	0.52
1:D:146:ASP:CB	1:D:150:GLY:H	2.23	0.52
1:B:146:ASP:HB2	1:B:150:GLY:O	2.10	0.52
1:B:61:PHE:CZ	1:B:140:LEU:HD13	2.45	0.50
1:A:71:VAL:HG12	1:A:109:SER:HB3	1.94	0.50
1:B:123:GLY:O	1:B:133:PRO:HA	2.11	0.49
1:B:146:ASP:HB2	1:B:150:GLY:N	2.26	0.49
1:C:169:LYS:N	1:C:169:LYS:HD2	2.26	0.49
1:B:141:ILE:HD11	1:B:156:GLU:HB3	1.94	0.49
1:C:146:ASP:HB2	1:C:150:GLY:O	2.12	0.49
1:A:69:ASP:HB3	2:A:320:8PP:H451	1.94	0.49
1:A:154:GLY:HA3	1:B:45:TYR:O	2.13	0.49
1:D:146:ASP:HB2	1:D:150:GLY:O	2.13	0.48
1:A:146:ASP:HB2	1:A:150:GLY:N	2.28	0.48
1:C:161:ILE:CD1	2:C:322:8PP:H23	2.44	0.48
1:A:31:GLY:HA2	3:B:322:HOH:O	2.14	0.47
1:C:123:GLY:O	1:C:133:PRO:HA	2.14	0.47
1:A:161:ILE:HD12	1:A:181:LEU:HD21	1.97	0.47
1:C:111:SER:HB2	3:C:377:HOH:O	2.14	0.47
1:B:122:THR:OG1	1:B:135:THR:HG22	2.15	0.47
1:D:122:THR:OG1	1:D:135:THR:HG22	2.15	0.47
1:A:36:VAL:HG22	2:A:320:8PP:H353	1.97	0.47
1:B:84:HIS:HE1	1:B:86:GLU:HB2	1.80	0.46
1:A:106:THR:HG23	3:A:372:HOH:O	2.16	0.46
1:C:106:THR:CG2	3:C:375:HOH:O	2.64	0.46
1:A:52:PHE:CE1	2:A:320:8PP:H222	2.50	0.46
1:C:141:ILE:HD11	1:C:156:GLU:HB3	1.98	0.46
1:C:73:VAL:HG21	2:C:322:8PP:H321	1.97	0.46
1:A:122:THR:OG1	1:A:135:THR:HG23	2.16	0.46
1:B:91:LEU:HD21	2:B:321:8PP:H122	1.99	0.45
1:D:146:ASP:HB2	1:D:150:GLY:H	1.81	0.45
1:D:76:ASN:HD22	1:D:104:GLN:HG2	1.82	0.44
1:C:73:VAL:HG22	1:C:107:PHE:HB3	1.99	0.44
1:D:122:THR:HA	1:D:135:THR:HG22	2.00	0.44
1:A:141:ILE:HD11	1:A:156:GLU:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ILE:CD1	2:B:321:8PP:H261	2.49	0.43
1:A:71:VAL:HG21	2:A:320:8PP:H38	1.99	0.43
1:C:77:THR:HG21	3:C:372:HOH:O	2.18	0.43
1:A:161:ILE:HD12	2:A:320:8PP:H23	2.01	0.43
1:A:39:ARG:HA	1:A:48:LEU:O	2.19	0.42
1:A:84:HIS:HE1	1:A:86:GLU:HB2	1.83	0.42
1:C:106:THR:HG22	1:C:126:THR:HB	2.02	0.42
1:C:103:PRO:O	1:C:104:GLN:HG3	2.19	0.42
1:D:71:VAL:HG21	2:D:323:8PP:C38	2.47	0.42
1:B:39:ARG:HA	1:B:48:LEU:O	2.20	0.42
2:B:321:8PP:H121	2:B:321:8PP:H152	1.70	0.42
1:C:146:ASP:HB2	1:C:150:GLY:N	2.34	0.42
1:D:71:VAL:CG2	2:D:323:8PP:H38	2.49	0.41
1:A:146:ASP:HB3	1:A:150:GLY:H	1.85	0.41
1:C:77:THR:OG1	1:C:99:THR:HG22	2.20	0.41
1:B:71:VAL:HG21	2:B:321:8PP:C38	2.50	0.41
1:C:84:HIS:CE1	1:C:86:GLU:HB2	2.56	0.41
1:D:123:GLY:O	1:D:133:PRO:HA	2.21	0.41
1:A:125:LEU:HG	2:A:320:8PP:H303	2.02	0.40
2:D:323:8PP:H101	2:D:323:8PP:H122	1.95	0.40
1:B:71:VAL:CG2	2:B:321:8PP:H38	2.50	0.40
1:D:146:ASP:HB3	1:D:150:GLY:H	1.87	0.40
1:A:141:ILE:HD12	1:B:46:SER:HA	2.03	0.40
1:A:162:LYS:HE3	1:A:162:LYS:HB2	1.88	0.40
1:C:106:THR:HG23	3:C:375:HOH:O	2.22	0.40
1:C:71:VAL:HG12	1:C:109:SER:CB	2.50	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	167/191 (87%)	158 (95%)	9 (5%)	0	100	100
1	B	167/191 (87%)	157 (94%)	10 (6%)	0	100	100
1	C	167/191 (87%)	155 (93%)	12 (7%)	0	100	100
1	D	167/191 (87%)	156 (93%)	11 (7%)	0	100	100
All	All	668/764 (87%)	626 (94%)	42 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	142/158 (90%)	132 (93%)	10 (7%)	18	19
1	B	142/158 (90%)	135 (95%)	7 (5%)	29	35
1	C	142/158 (90%)	134 (94%)	8 (6%)	25	29
1	D	142/158 (90%)	135 (95%)	7 (5%)	29	35
All	All	568/632 (90%)	536 (94%)	32 (6%)	25	29

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	106	THR
1	A	117	ASP
1	A	125	LEU
1	A	127	LEU
1	A	130	VAL
1	A	135	THR
1	A	146	ASP
1	A	162	LYS
1	A	169	LYS
1	B	98	ASN
1	B	111	SER
1	B	125	LEU

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Mol	Chain	Res	Type
1	B	127	LEU
1	B	130	VAL
1	B	146	ASP
1	B	162	LYS
1	C	98	ASN
1	C	106	THR
1	C	125	LEU
1	C	127	LEU
1	C	130	VAL
1	C	135	THR
1	C	146	ASP
1	C	169	LYS
1	D	98	ASN
1	D	106	THR
1	D	117	ASP
1	D	125	LEU
1	D	127	LEU
1	D	130	VAL
1	D	146	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	76	ASN
1	A	98	ASN
1	A	104	GLN
1	B	37	ASN
1	B	76	ASN
1	B	90	HIS
1	B	98	ASN
1	B	104	GLN
1	C	37	ASN
1	C	76	ASN
1	C	98	ASN
1	D	37	ASN
1	D	76	ASN
1	D	98	ASN
1	D	104	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 ⓘ

4 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	8PP	A	320	-	47,47,47	2.22	8 (17%)	56,57,57	2.62	17 (30%)
2	8PP	B	321	-	47,47,47	1.94	9 (19%)	56,57,57	2.65	20 (35%)
2	8PP	C	322	-	47,47,47	1.86	8 (17%)	56,57,57	2.84	19 (33%)
2	8PP	D	323	-	47,47,47	1.98	9 (19%)	56,57,57	2.49	16 (28%)

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	8PP	A	320	-	-	0/47/47/47	0/1/1/1
2	8PP	B	321	-	-	0/47/47/47	0/1/1/1
2	8PP	C	322	-	-	0/47/47/47	0/1/1/1
2	8PP	D	323	-	-	0/47/47/47	0/1/1/1

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	321	8PP	O-C1	-6.26	1.23	1.36
2	D	323	8PP	O-C1	-6.05	1.24	1.36
2	A	320	8PP	O-C1	-5.98	1.24	1.36
2	C	322	8PP	O-C1	-5.91	1.24	1.36
2	B	321	8PP	C3-C2	-3.39	1.33	1.39
2	A	320	8PP	C3-C2	-2.63	1.35	1.39
2	C	322	8PP	C3-C2	-2.27	1.35	1.39
2	C	322	8PP	C42-C43	-2.21	1.42	1.50
2	D	323	8PP	C42-C43	-2.07	1.43	1.50
2	D	323	8PP	C18-C19	2.04	1.38	1.33
2	B	321	8PP	C30-C29	2.11	1.56	1.50
2	B	321	8PP	C18-C19	2.34	1.38	1.33
2	B	321	8PP	C1-C2	2.55	1.43	1.40
2	C	322	8PP	C23-C24	2.70	1.39	1.33
2	A	320	8PP	C23-C24	2.70	1.39	1.33
2	D	323	8PP	C23-C24	2.72	1.39	1.33
2	B	321	8PP	C23-C24	2.88	1.40	1.33
2	C	322	8PP	C1-C2	2.90	1.44	1.40
2	D	323	8PP	C11-C9	2.95	1.57	1.51
2	B	321	8PP	C5-C4	3.20	1.45	1.38
2	A	320	8PP	C1-C2	3.27	1.44	1.40
2	B	321	8PP	C4-C3	3.34	1.45	1.38
2	A	320	8PP	C5-C4	3.59	1.46	1.38
2	D	323	8PP	C1-C2	3.80	1.45	1.40
2	D	323	8PP	C5-C4	3.81	1.47	1.38
2	A	320	8PP	C4-C3	3.91	1.46	1.38
2	C	322	8PP	C5-C4	4.06	1.47	1.38
2	D	323	8PP	C4-C3	4.13	1.46	1.38
2	C	322	8PP	C4-C3	4.31	1.47	1.38
2	C	322	8PP	C6-C1	4.58	1.47	1.39
2	D	323	8PP	C6-C1	5.02	1.48	1.39
2	A	320	8PP	C6-C1	5.08	1.48	1.39
2	B	321	8PP	C6-C1	5.12	1.48	1.39
2	A	320	8PP	C8-C9	8.08	1.53	1.33

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	322	8PP	C7-C8-C9	-6.17	117.64	127.43
2	D	323	8PP	C7-C8-C9	-5.73	118.34	127.43
2	A	320	8PP	C7-C8-C9	-5.56	118.61	127.43
2	B	321	8PP	C41-C39-C40	-5.50	105.74	115.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	320	8PP	C41-C39-C40	-5.40	105.92	115.29
2	C	322	8PP	C11-C9-C8	-5.28	110.30	121.10
2	C	322	8PP	C41-C39-C40	-5.26	106.16	115.29
2	C	322	8PP	C32-C33-C34	-5.09	114.90	127.68
2	A	320	8PP	C16-C14-C13	-5.03	110.81	121.10
2	D	323	8PP	C41-C39-C40	-4.97	106.67	115.29
2	B	321	8PP	C32-C33-C34	-4.94	115.26	127.68
2	C	322	8PP	C12-C13-C14	-4.91	115.34	127.68
2	C	322	8PP	C17-C18-C19	-4.82	115.57	127.68
2	A	320	8PP	C32-C33-C34	-4.79	115.66	127.68
2	D	323	8PP	C32-C33-C34	-4.76	115.73	127.68
2	A	320	8PP	C36-C34-C33	-4.72	111.44	121.10
2	B	321	8PP	C12-C13-C14	-4.69	115.91	127.68
2	B	321	8PP	C36-C34-C33	-4.53	111.84	121.10
2	D	323	8PP	C36-C34-C33	-4.52	111.85	121.10
2	B	321	8PP	C30-C29-C31	-4.44	107.58	115.29
2	B	321	8PP	C7-C8-C9	-4.40	120.45	127.43
2	C	322	8PP	C36-C34-C33	-4.40	112.10	121.10
2	D	323	8PP	C22-C23-C24	-4.26	116.98	127.68
2	B	321	8PP	C15-C14-C13	-4.25	112.34	123.69
2	A	320	8PP	C22-C23-C24	-4.22	117.08	127.68
2	B	321	8PP	C11-C9-C8	-4.10	112.71	121.10
2	A	320	8PP	C11-C9-C8	-4.09	112.72	121.10
2	C	322	8PP	C22-C23-C24	-4.03	117.55	127.68
2	B	321	8PP	C22-C23-C24	-3.99	117.66	127.68
2	D	323	8PP	C15-C14-C13	-3.73	113.73	123.69
2	D	323	8PP	C11-C9-C8	-3.68	113.58	121.10
2	C	322	8PP	C37-C38-C39	-3.55	118.76	127.68
2	C	322	8PP	C15-C14-C13	-3.49	114.37	123.69
2	D	323	8PP	C37-C38-C39	-3.39	119.17	127.68
2	A	320	8PP	C37-C38-C39	-3.28	119.44	127.68
2	B	321	8PP	C37-C38-C39	-3.26	119.49	127.68
2	C	322	8PP	C16-C14-C13	-3.12	114.72	121.10
2	D	323	8PP	C16-C14-C13	-2.93	115.09	121.10
2	B	321	8PP	C17-C18-C19	-2.89	120.41	127.68
2	D	323	8PP	C17-C18-C19	-2.83	120.58	127.68
2	C	322	8PP	C30-C29-C31	-2.77	110.48	115.29
2	A	320	8PP	C17-C18-C19	-2.73	120.83	127.68
2	D	323	8PP	C30-C29-C31	-2.50	110.94	115.29
2	C	322	8PP	C2-C7-C8	-2.25	107.93	112.48
2	B	321	8PP	C4-C5-C6	-2.14	117.26	120.21
2	A	320	8PP	C15-C14-C13	-2.08	118.13	123.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	322	8PP	C20-C19-C18	-2.01	118.32	123.69
2	B	321	8PP	C30-C29-C28	2.10	129.31	123.69
2	A	320	8PP	C16-C17-C18	2.38	120.14	111.97
2	C	322	8PP	C26-C24-C23	2.82	126.87	121.10
2	A	320	8PP	C26-C24-C23	2.92	127.08	121.10
2	B	321	8PP	C12-C11-C9	2.95	122.92	112.93
2	B	321	8PP	C26-C24-C23	2.97	127.19	121.10
2	D	323	8PP	C26-C24-C23	3.05	127.35	121.10
2	D	323	8PP	C41-C39-C38	3.39	132.74	123.69
2	D	323	8PP	C10-C9-C8	3.45	132.91	123.69
2	C	322	8PP	C41-C39-C38	3.48	132.99	123.69
2	B	321	8PP	C41-C39-C38	3.53	133.11	123.69
2	B	321	8PP	C2-C7-C8	3.60	119.77	112.48
2	A	320	8PP	C17-C16-C14	3.66	125.30	112.93
2	A	320	8PP	C41-C39-C38	3.68	133.52	123.69
2	A	320	8PP	C10-C9-C8	3.72	133.63	123.69
2	D	323	8PP	C35-C34-C36	3.79	121.87	115.29
2	B	321	8PP	C35-C34-C36	3.81	121.91	115.29
2	C	322	8PP	C35-C34-C36	3.87	122.01	115.29
2	A	320	8PP	C35-C34-C36	4.06	122.33	115.29
2	B	321	8PP	C10-C9-C11	4.86	123.73	115.29
2	C	322	8PP	C10-C9-C11	6.43	126.45	115.29
2	B	321	8PP	C15-C14-C16	7.44	128.19	115.29
2	C	322	8PP	C15-C14-C16	9.00	130.91	115.29
2	D	323	8PP	C15-C14-C16	9.03	130.95	115.29
2	A	320	8PP	C15-C14-C16	9.08	131.05	115.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 32 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	320	8PP	8	0
2	B	321	8PP	8	0
2	C	322	8PP	6	0
2	D	323	8PP	10	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	169/191 (88%)	-0.11	7 (4%) 38 36	17, 32, 64, 75	0
1	B	169/191 (88%)	-0.19	3 (1%) 69 66	18, 30, 50, 71	0
1	C	169/191 (88%)	-0.28	2 (1%) 79 77	16, 27, 44, 68	0
1	D	169/191 (88%)	-0.06	5 (2%) 51 48	15, 33, 56, 69	0
All	All	676/764 (88%)	-0.16	17 (2%) 58 55	15, 31, 54, 75	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	117	ASP	5.2
1	D	148	TRP	4.6
1	A	191	LYS	3.9
1	D	117	ASP	3.6
1	B	191	LYS	3.3
1	A	115	ASP	3.3
1	A	67	ALA	3.1
1	D	191	LYS	2.7
1	D	115	ASP	2.6
1	B	115	ASP	2.5
1	A	66	PRO	2.4
1	C	116	GLY	2.4
1	C	191	LYS	2.3
1	A	64	LYS	2.2
1	B	30	GLU	2.1
1	A	113	LYS	2.0
1	D	116	GLY	2.0

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 [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(\AA^2)	Q<0.9
2	8PP	D	323	47/47	0.83	0.26	3.88	27,36,67,70	0
2	8PP	C	322	47/47	0.86	0.24	2.88	28,35,64,65	0
2	8PP	B	321	47/47	0.90	0.22	2.64	19,30,43,75	0
2	8PP	A	320	47/47	0.86	0.25	2.42	21,33,53,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.