



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 01:32 PM GMT

PDB ID : 3TYH
Title : Crystal structure of oxo-copper clusters binding to ferric binding protein from *Neisseria gonorrhoeae*
Authors : Chen, W.J.; Wang, H.F.; Zhou, C.J.; Ye, D.R.; Huang, J.; Tan, X.S.; Zhong, W.Q.
Deposited on : 2011-09-26
Resolution : 2.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

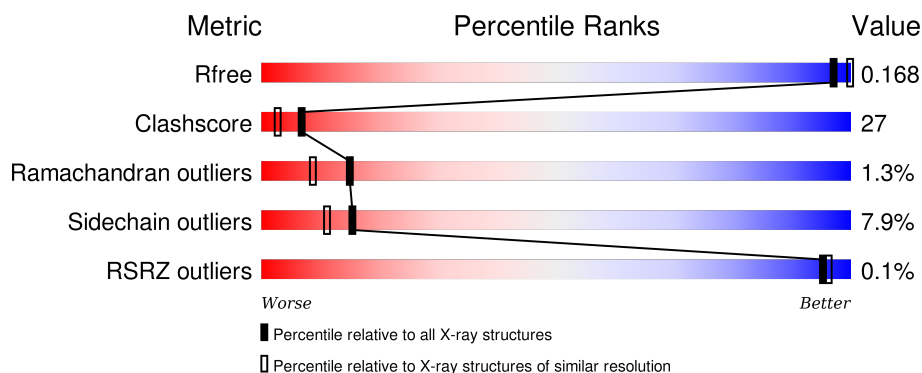
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	309	 61% 36% •
1	B	309	 57% 38% 5%
1	C	309	 62% 33% •
1	D	309	 55% 37% 6% •
1	E	309	 65% 29% 5%

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Mol	Chain	Length	Quality of chain
1	F	309	<div><div></div><div>61%34%5% •</div></div>
1	G	309	<div><div></div><div>60%35%5%</div></div>
1	H	309	<div><div>%<div></div></div><div>48%47%5%</div></div>
1	I	309	<div><div></div><div>47%47%6%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 22889 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 FbpA protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			
1	B	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			
1	C	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			
1	D	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			
1	E	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			
1	F	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			
1	G	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			
1	H	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			
1	I	309	Total	C	N	O	S	0	0	0
			2378	1508	423	446	1			

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	2	Total	Cu	0	0
			2	2		
2	D	2	Total	Cu	0	0
			2	2		
2	E	2	Total	Cu	0	0
			2	2		
2	H	2	Total	Cu	0	0
			2	2		
2	B	2	Total	Cu	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	I	2	Total 2	Cu 2	0	0
2	C	2	Total 2	Cu 2	0	0
2	A	2	Total 2	Cu 2	0	0
2	F	2	Total 2	Cu 2	0	0

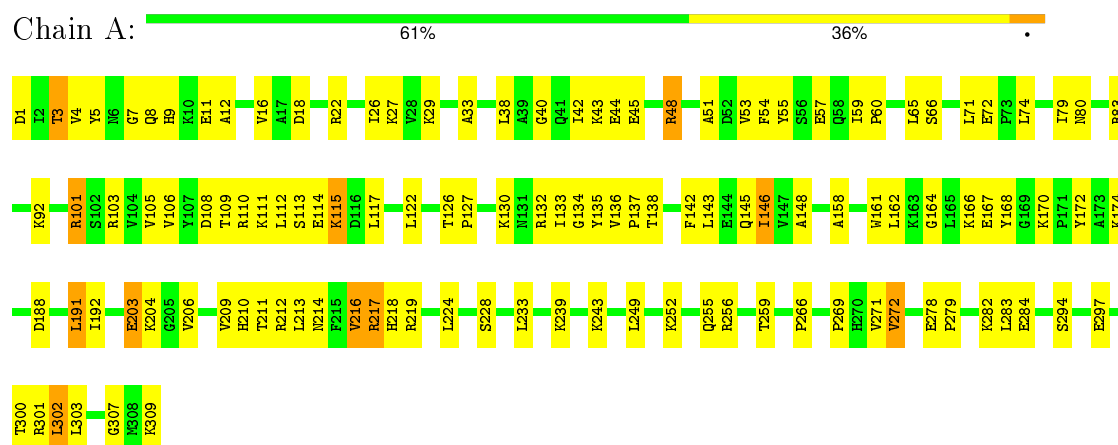
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	217	Total 217	O 217	0	0
3	B	166	Total 166	O 166	0	0
3	C	188	Total 188	O 188	0	0
3	D	177	Total 177	O 177	0	0
3	E	178	Total 178	O 178	0	0
3	F	181	Total 181	O 181	0	0
3	G	123	Total 123	O 123	0	0
3	H	120	Total 120	O 120	0	0
3	I	119	Total 119	O 119	0	0

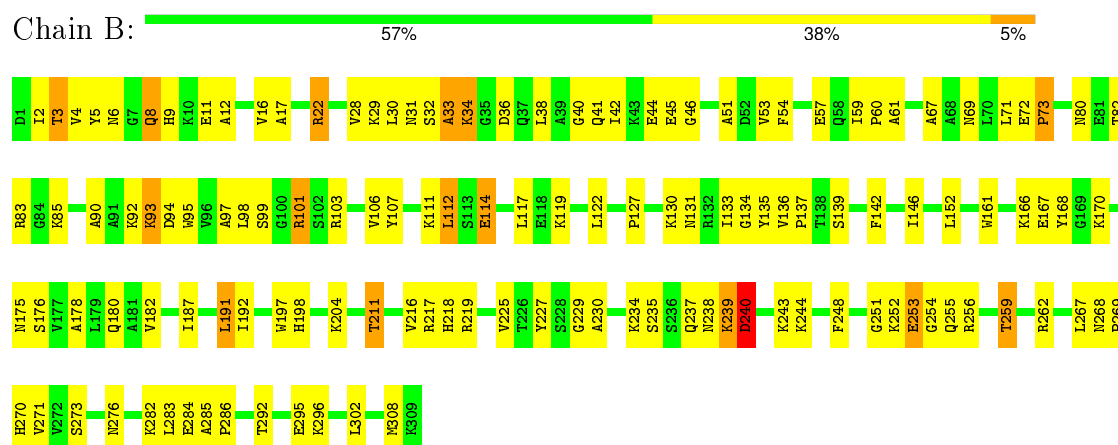
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

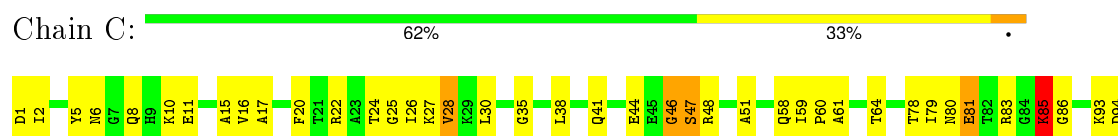
• Molecule 1: FbpA protein

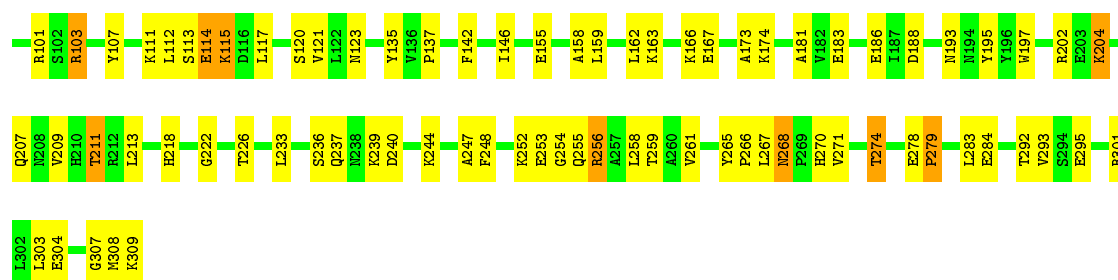


• Molecule 1: FbpA protein



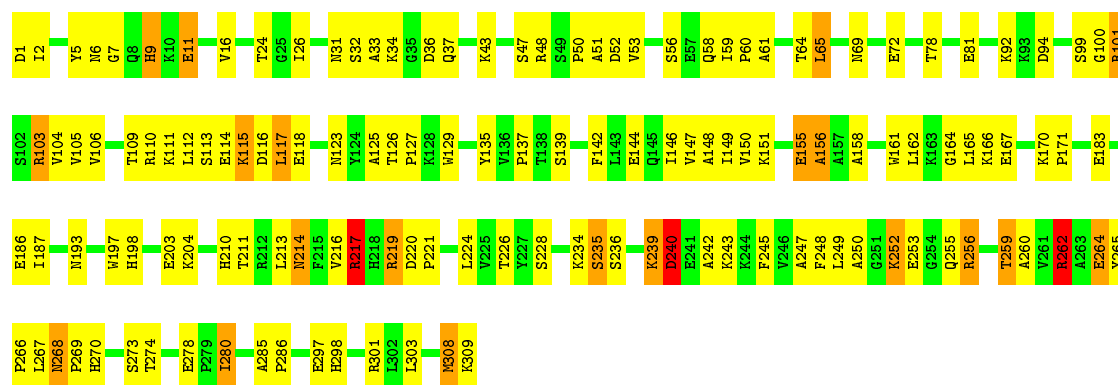
• Molecule 1: FbpA protein





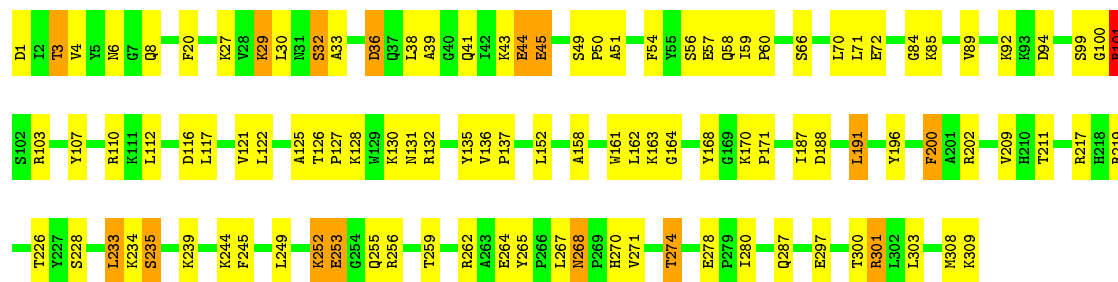
- Molecule 1: FbpA protein

Chain D: 55% 37% 6% .



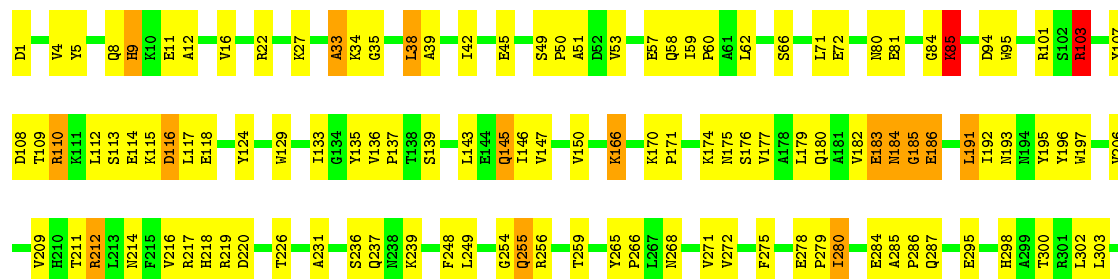
- Molecule 1: FbpA protein

Chain E: 65% 29% 5%



- Molecule 1: FbpA protein

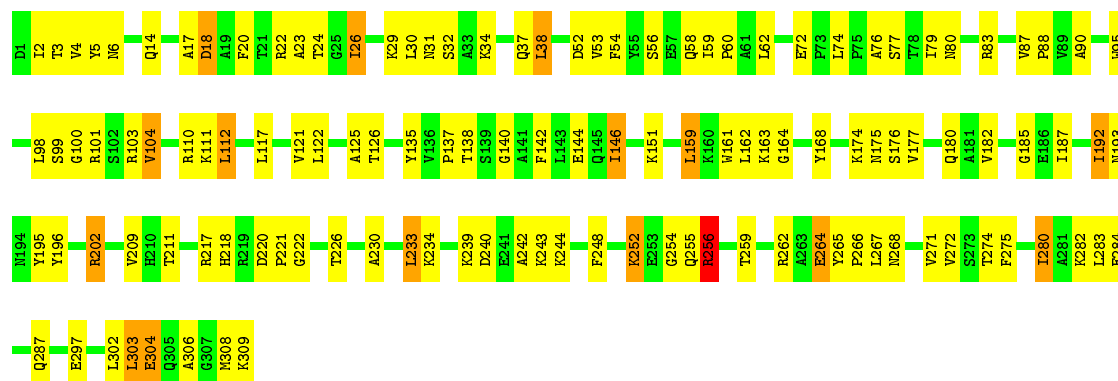
Chain F: 61% 34% 5% .



G307
K308
K309

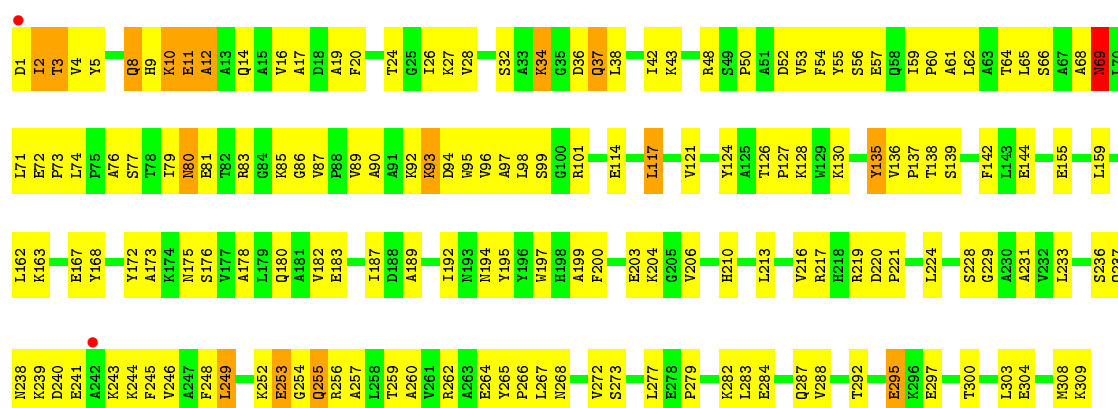
• Molecule 1: FbpA protein

Chain G: 



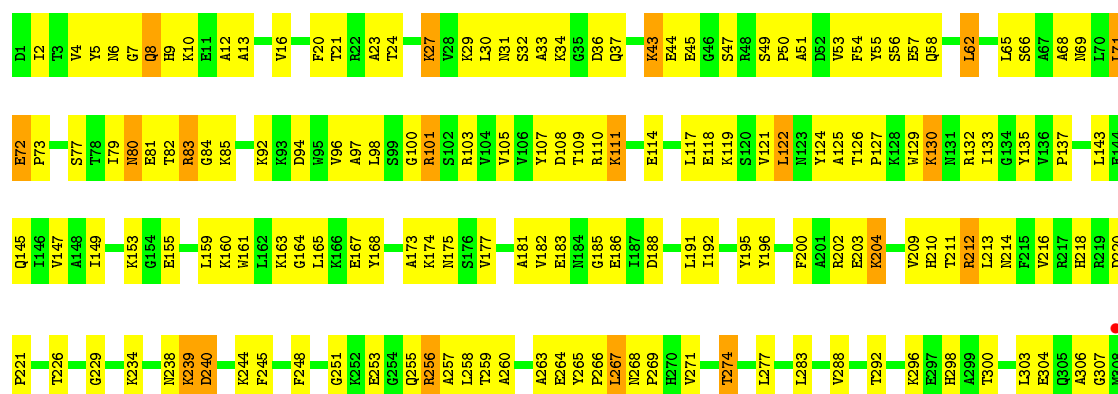
• Molecule 1: FbpA protein

Chain H: 



• Molecule 1: FbpA protein

Chain I: 



K309

4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	146.51Å 146.51Å 114.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.96 – 2.10 47.96 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.96-2.10) 100.0 (47.96-2.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.30 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.143 , 0.169 0.147 , 0.168	Depositor DCC
R_{free} test set	8174 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.001	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 41.7	EDS
Estimated twinning fraction	0.541 for H, K, L 0.459 for H+K, -K, -L 0.004 for -h,-k,l 0.018 for h,-h-k,-l 0.045 for -k,-h,-l	Xtriage
Reported twinning fraction	0.541 for H, K, L 0.459 for H+K, -K, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 160455 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	22889	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.75% 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: CU

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	0.69	0/2423	0.93	3/3280 (0.1%)
1	B	0.64	0/2423	0.97	3/3280 (0.1%)
1	C	0.65	0/2423	0.95	2/3280 (0.1%)
1	D	0.60	0/2423	0.92	3/3280 (0.1%)
1	E	0.63	0/2423	0.94	5/3280 (0.2%)
1	F	0.73	1/2423 (0.0%)	1.01	9/3280 (0.3%)
1	G	0.55	0/2423	0.90	2/3280 (0.1%)
1	H	0.48	0/2423	0.87	1/3280 (0.0%)
1	I	0.48	0/2423	0.88	2/3280 (0.1%)
All	All	0.61	1/21807 (0.0%)	0.93	30/29520 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	85	LYS	C-N	-6.11	1.22	1.33

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	103	ARG	NE-CZ-NH2	-12.65	113.97	120.30
1	F	85	LYS	C-N-CA	-9.06	103.27	122.30
1	E	101	ARG	NE-CZ-NH1	-8.52	116.04	120.30
1	A	101	ARG	NE-CZ-NH2	-7.37	116.61	120.30
1	E	103	ARG	NE-CZ-NH1	7.10	123.85	120.30

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	2378	0	2424	124	0
1	B	2378	0	2424	118	0
1	C	2378	0	2424	111	0
1	D	2378	0	2424	154	0
1	E	2378	0	2424	88	0
1	F	2378	0	2423	133	0
1	G	2378	0	2425	112	0
1	H	2378	0	2424	172	0
1	I	2378	0	2424	155	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
2	I	2	0	0	0	0
3	A	217	0	0	17	0
3	B	166	0	0	14	0
3	C	188	0	0	11	0
3	D	177	0	0	19	0
3	E	178	0	0	3	0
3	F	181	0	0	14	0
3	G	123	0	0	8	0
3	H	120	0	0	30	0
3	I	119	0	0	18	0
All	All	22889	0	21816	1160	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 27.

The worst 5 of 1160 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:252:LYS:HE2	1:B:256:ARG:NH2	1.45	1.30
1:C:202:ARG:HH22	1:C:274:THR:HG23	1.07	1.12
1:A:284:GLU:HB3	3:A:1406:HOH:O	1.49	1.11
1:I:202:ARG:HH22	1:I:274:THR:HG23	1.14	1.11
1:F:192:ILE:HG22	1:F:193:ASN:H	1.13	1.08

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	307/309 (99%)	297 (97%)	10 (3%)	0	100	100
1	B	307/309 (99%)	280 (91%)	21 (7%)	6 (2%)	9	4
1	C	307/309 (99%)	284 (92%)	18 (6%)	5 (2%)	12	6
1	D	307/309 (99%)	286 (93%)	16 (5%)	5 (2%)	12	6
1	E	307/309 (99%)	296 (96%)	9 (3%)	2 (1%)	26	21
1	F	307/309 (99%)	284 (92%)	18 (6%)	5 (2%)	12	6
1	G	307/309 (99%)	284 (92%)	19 (6%)	4 (1%)	15	9
1	H	307/309 (99%)	262 (85%)	39 (13%)	6 (2%)	9	4
1	I	307/309 (99%)	274 (89%)	30 (10%)	3 (1%)	19	13
All	All	2763/2781 (99%)	2547 (92%)	180 (6%)	36 (1%)	15	9

5 of 36 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	85	LYS
1	H	8	GLN
1	H	81	GLU
1	B	33	ALA
1	B	85	LYS

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	245/245 (100%)	230 (94%)	15 (6%)	23	19
1	B	245/245 (100%)	228 (93%)	17 (7%)	19	15
1	C	245/245 (100%)	227 (93%)	18 (7%)	17	13
1	D	245/245 (100%)	223 (91%)	22 (9%)	12	8
1	E	245/245 (100%)	224 (91%)	21 (9%)	13	9
1	F	245/245 (100%)	232 (95%)	13 (5%)	28	25
1	G	245/245 (100%)	228 (93%)	17 (7%)	19	15
1	H	245/245 (100%)	223 (91%)	22 (9%)	12	8
1	I	245/245 (100%)	215 (88%)	30 (12%)	6	3
All	All	2205/2205 (100%)	2030 (92%)	175 (8%)	15	11

5 of 175 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	163	LYS
1	F	145	GLN
1	I	122	LEU
1	E	191	LEU
1	E	268	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 70 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	255	GLN
1	F	255	GLN
1	I	69	ASN
1	E	268	ASN
1	F	198	HIS

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 18 ligands modelled in this entry, 18 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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	309/309 (100%)	-0.77	0 100 100	12, 20, 31, 41	0
1	B	309/309 (100%)	-0.66	0 100 100	11, 23, 45, 62	0
1	C	309/309 (100%)	-0.74	0 100 100	12, 23, 37, 55	0
1	D	309/309 (100%)	-0.72	0 100 100	15, 26, 38, 51	0
1	E	309/309 (100%)	-0.79	0 100 100	13, 23, 34, 46	0
1	F	309/309 (100%)	-0.78	0 100 100	13, 22, 36, 48	0
1	G	309/309 (100%)	-0.71	0 100 100	18, 29, 41, 50	0
1	H	309/309 (100%)	-0.37	2 (0%) 90 92	23, 38, 59, 65	0
1	I	309/309 (100%)	-0.55	1 (0%) 94 95	22, 36, 49, 67	0
All	All	2781/2781 (100%)	-0.68	3 (0%) 95 96	11, 27, 48, 67	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	308	MET	3.0
1	H	242	ALA	2.2
1	H	1	ASP	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	CU	B	310	1/1	0.98	0.09	-0.10	41,41,41,41	0
2	CU	F	311	1/1	0.92	0.09	-	61,61,61,61	0
2	CU	D	311	1/1	0.95	0.10	-	50,50,50,50	0
2	CU	H	311	1/1	0.94	0.06	-	71,71,71,71	0
2	CU	C	311	1/1	0.99	0.09	-	37,37,37,37	0
2	CU	A	311	1/1	1.00	0.09	-	37,37,37,37	0
2	CU	G	311	1/1	0.96	0.07	-	55,55,55,55	0
2	CU	E	311	1/1	0.99	0.09	-	45,45,45,45	0
2	CU	I	310	1/1	0.99	0.05	-	46,46,46,46	0
2	CU	I	311	1/1	0.95	0.06	-	64,64,64,64	0
2	CU	C	310	1/1	0.99	0.07	-	49,49,49,49	0
2	CU	F	310	1/1	0.91	0.08	-	53,53,53,53	0
2	CU	A	310	1/1	0.99	0.09	-	38,38,38,38	0
2	CU	D	310	1/1	0.97	0.12	-	45,45,45,45	0
2	CU	G	310	1/1	0.98	0.05	-	53,53,53,53	0
2	CU	E	310	1/1	0.99	0.08	-	46,46,46,46	0
2	CU	H	310	1/1	0.96	0.06	-	56,56,56,56	0
2	CU	B	311	1/1	0.98	0.09	-	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.