



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:08 PM GMT

PDB ID : 4BLB  
Title : Crystal structure of a human Suppressor of fused (SUFU)-GLI1p complex  
Authors : Cherry, A.L.; Finta, C.; Karlstrom, M.; De Sanctis, D.; Toftgard, R.; Jovine, L.  
Deposited on : 2013-05-02  
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](mailto: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.

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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

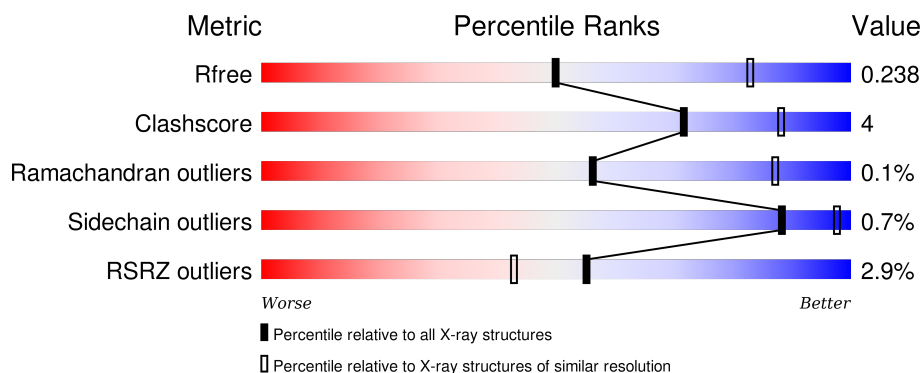
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (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  $\geq 3$ , 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	753	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 89%, yellow 89%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>89%</span> <span>8%</span> <span>••</span> </div> </div>
1	B	753	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 2%, green 2%, green 84%, yellow 84%, yellow 95%, grey 95%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>2%</span> <span>84%</span> <span>11%</span> <span>5%</span> </div> </div>
1	C	753	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 7%, green 7%, green 87%, yellow 87%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>7%</span> <span>87%</span> <span>10%</span> <span>•</span> </div> </div>
1	D	753	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 89%, yellow 89%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>89%</span> <span>8%</span> <span>•</span> </div> </div>
2	E	17	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 0%, green 35%, yellow 35%, yellow 53%, grey 53%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>35%</span> <span>18%</span> <span>47%</span> </div> </div>

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Mol	Chain	Length	Quality of chain
2	F	17	
2	G	17	
2	H	17	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ZN	C	910	-	-	-	X
4	ZN	D	910	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23141 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 MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	734	Total	C	N	O	S	0	0	0
			5757	3681	967	1090	19			
1	B	717	Total	C	N	O	S	0	0	0
			5621	3597	942	1064	18			
1	C	725	Total	C	N	O	S	0	0	0
			5687	3640	957	1072	18			
1	D	730	Total	C	N	O	S	0	0	0
			5727	3666	962	1081	18			

There are 128 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP P0AEX9
A	3	THR	ILE	ENGINEERED MUTATION	UNP P0AEX9
A	360	ALA	GLU	ENGINEERED MUTATION	UNP P0AEX9
A	363	ALA	LYS	ENGINEERED MUTATION	UNP P0AEX9
A	364	ALA	ASP	ENGINEERED MUTATION	UNP P0AEX9
A	368	ASN	ARG	ENGINEERED MUTATION	UNP P0AEX9
A	369	ALA	-	LINKER	UNP P0AEX9
A	370	ALA	-	LINKER	UNP P0AEX9
A	371	ALA	-	LINKER	UNP P0AEX9
A	216	HIS	ALA	ENGINEERED MUTATION	UNP P0AEX9
A	220	HIS	LYS	ENGINEERED MUTATION	UNP P0AEX9
A	401	ASP	TRP	ENGINEERED MUTATION	UNP Q9UMX1
A	402	SER	LEU	ENGINEERED MUTATION	UNP Q9UMX1
A	403	PHE	GLY	ENGINEERED MUTATION	UNP Q9UMX1
A	619	PRO	-	LINKER	UNP Q9UMX1
A	620	SER	-	LINKER	UNP Q9UMX1
A	621	ARG	-	LINKER	UNP Q9UMX1
A	622	GLY	-	LINKER	UNP Q9UMX1
A	625	PRO	-	LINKER	UNP Q9UMX1
A	718	ALA	PRO	ENGINEERED MUTATION	UNP Q9UMX1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	.	-	GLU	DELETION	UNP Q9UMX1
A	.	-	GLU	DELETION	UNP Q9UMX1
A	.	-	PHE	DELETION	UNP Q9UMX1
A	719	ALA	LYS	ENGINEERED MUTATION	UNP Q9UMX1
A	746	VAL	-	EXPRESSION TAG	UNP Q9UMX1
A	747	GLU	-	EXPRESSION TAG	UNP Q9UMX1
A	748	HIS	-	EXPRESSION TAG	UNP Q9UMX1
A	749	HIS	-	EXPRESSION TAG	UNP Q9UMX1
A	750	HIS	-	EXPRESSION TAG	UNP Q9UMX1
A	751	HIS	-	EXPRESSION TAG	UNP Q9UMX1
A	752	HIS	-	EXPRESSION TAG	UNP Q9UMX1
A	753	HIS	-	EXPRESSION TAG	UNP Q9UMX1
B	1	MET	-	EXPRESSION TAG	UNP P0AEX9
B	3	THR	ILE	ENGINEERED MUTATION	UNP P0AEX9
B	360	ALA	GLU	ENGINEERED MUTATION	UNP P0AEX9
B	363	ALA	LYS	ENGINEERED MUTATION	UNP P0AEX9
B	364	ALA	ASP	ENGINEERED MUTATION	UNP P0AEX9
B	368	ASN	ARG	ENGINEERED MUTATION	UNP P0AEX9
B	369	ALA	-	LINKER	UNP P0AEX9
B	370	ALA	-	LINKER	UNP P0AEX9
B	371	ALA	-	LINKER	UNP P0AEX9
B	216	HIS	ALA	ENGINEERED MUTATION	UNP P0AEX9
B	220	HIS	LYS	ENGINEERED MUTATION	UNP P0AEX9
B	401	ASP	TRP	ENGINEERED MUTATION	UNP Q9UMX1
B	402	SER	LEU	ENGINEERED MUTATION	UNP Q9UMX1
B	403	PHE	GLY	ENGINEERED MUTATION	UNP Q9UMX1
B	619	PRO	-	LINKER	UNP Q9UMX1
B	620	SER	-	LINKER	UNP Q9UMX1
B	621	ARG	-	LINKER	UNP Q9UMX1
B	622	GLY	-	LINKER	UNP Q9UMX1
B	625	PRO	-	LINKER	UNP Q9UMX1
B	718	ALA	PRO	ENGINEERED MUTATION	UNP Q9UMX1
B	.	-	GLU	DELETION	UNP Q9UMX1
B	.	-	GLU	DELETION	UNP Q9UMX1
B	.	-	PHE	DELETION	UNP Q9UMX1
B	719	ALA	LYS	ENGINEERED MUTATION	UNP Q9UMX1
B	746	VAL	-	EXPRESSION TAG	UNP Q9UMX1
B	747	GLU	-	EXPRESSION TAG	UNP Q9UMX1
B	748	HIS	-	EXPRESSION TAG	UNP Q9UMX1
B	749	HIS	-	EXPRESSION TAG	UNP Q9UMX1
B	750	HIS	-	EXPRESSION TAG	UNP Q9UMX1
B	751	HIS	-	EXPRESSION TAG	UNP Q9UMX1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	752	HIS	-	EXPRESSION TAG	UNP Q9UMX1
B	753	HIS	-	EXPRESSION TAG	UNP Q9UMX1
C	1	MET	-	EXPRESSION TAG	UNP P0AEX9
C	3	THR	ILE	ENGINEERED MUTATION	UNP P0AEX9
C	360	ALA	GLU	ENGINEERED MUTATION	UNP P0AEX9
C	363	ALA	LYS	ENGINEERED MUTATION	UNP P0AEX9
C	364	ALA	ASP	ENGINEERED MUTATION	UNP P0AEX9
C	368	ASN	ARG	ENGINEERED MUTATION	UNP P0AEX9
C	369	ALA	-	LINKER	UNP P0AEX9
C	370	ALA	-	LINKER	UNP P0AEX9
C	371	ALA	-	LINKER	UNP P0AEX9
C	216	HIS	ALA	ENGINEERED MUTATION	UNP P0AEX9
C	220	HIS	LYS	ENGINEERED MUTATION	UNP P0AEX9
C	401	ASP	TRP	ENGINEERED MUTATION	UNP Q9UMX1
C	402	SER	LEU	ENGINEERED MUTATION	UNP Q9UMX1
C	403	PHE	GLY	ENGINEERED MUTATION	UNP Q9UMX1
C	619	PRO	-	LINKER	UNP Q9UMX1
C	620	SER	-	LINKER	UNP Q9UMX1
C	621	ARG	-	LINKER	UNP Q9UMX1
C	622	GLY	-	LINKER	UNP Q9UMX1
C	625	PRO	-	LINKER	UNP Q9UMX1
C	718	ALA	PRO	ENGINEERED MUTATION	UNP Q9UMX1
C	.	-	GLU	DELETION	UNP Q9UMX1
C	.	-	GLU	DELETION	UNP Q9UMX1
C	.	-	PHE	DELETION	UNP Q9UMX1
C	719	ALA	LYS	ENGINEERED MUTATION	UNP Q9UMX1
C	746	VAL	-	EXPRESSION TAG	UNP Q9UMX1
C	747	GLU	-	EXPRESSION TAG	UNP Q9UMX1
C	748	HIS	-	EXPRESSION TAG	UNP Q9UMX1
C	749	HIS	-	EXPRESSION TAG	UNP Q9UMX1
C	750	HIS	-	EXPRESSION TAG	UNP Q9UMX1
C	751	HIS	-	EXPRESSION TAG	UNP Q9UMX1
C	752	HIS	-	EXPRESSION TAG	UNP Q9UMX1
C	753	HIS	-	EXPRESSION TAG	UNP Q9UMX1
D	1	MET	-	EXPRESSION TAG	UNP P0AEX9
D	3	THR	ILE	ENGINEERED MUTATION	UNP P0AEX9
D	360	ALA	GLU	ENGINEERED MUTATION	UNP P0AEX9
D	363	ALA	LYS	ENGINEERED MUTATION	UNP P0AEX9
D	364	ALA	ASP	ENGINEERED MUTATION	UNP P0AEX9
D	368	ASN	ARG	ENGINEERED MUTATION	UNP P0AEX9
D	369	ALA	-	LINKER	UNP P0AEX9
D	370	ALA	-	LINKER	UNP P0AEX9

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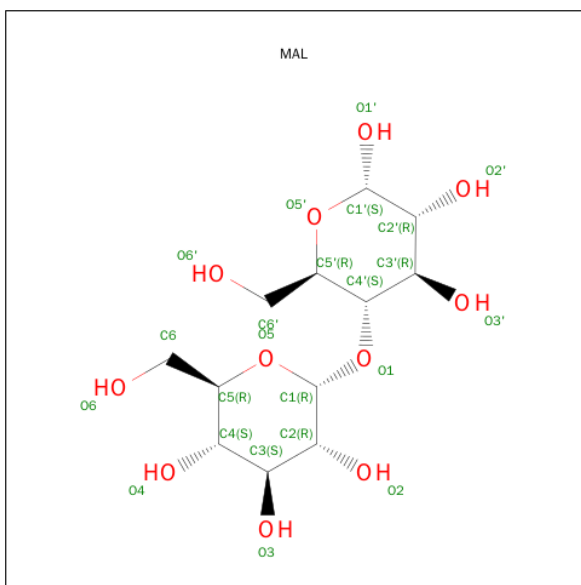
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Chain	Residue	Modelled	Actual	Comment	Reference
D	371	ALA	-	LINKER	UNP P0AEX9
D	216	HIS	ALA	ENGINEERED MUTATION	UNP P0AEX9
D	220	HIS	LYS	ENGINEERED MUTATION	UNP P0AEX9
D	401	ASP	TRP	ENGINEERED MUTATION	UNP Q9UMX1
D	402	SER	LEU	ENGINEERED MUTATION	UNP Q9UMX1
D	403	PHE	GLY	ENGINEERED MUTATION	UNP Q9UMX1
D	619	PRO	-	LINKER	UNP Q9UMX1
D	620	SER	-	LINKER	UNP Q9UMX1
D	621	ARG	-	LINKER	UNP Q9UMX1
D	622	GLY	-	LINKER	UNP Q9UMX1
D	625	PRO	-	LINKER	UNP Q9UMX1
D	718	ALA	PRO	ENGINEERED MUTATION	UNP Q9UMX1
D	.	-	GLU	DELETION	UNP Q9UMX1
D	.	-	GLU	DELETION	UNP Q9UMX1
D	.	-	PHE	DELETION	UNP Q9UMX1
D	719	ALA	LYS	ENGINEERED MUTATION	UNP Q9UMX1
D	746	VAL	-	EXPRESSION TAG	UNP Q9UMX1
D	747	GLU	-	EXPRESSION TAG	UNP Q9UMX1
D	748	HIS	-	EXPRESSION TAG	UNP Q9UMX1
D	749	HIS	-	EXPRESSION TAG	UNP Q9UMX1
D	750	HIS	-	EXPRESSION TAG	UNP Q9UMX1
D	751	HIS	-	EXPRESSION TAG	UNP Q9UMX1
D	752	HIS	-	EXPRESSION TAG	UNP Q9UMX1
D	753	HIS	-	EXPRESSION TAG	UNP Q9UMX1

- Molecule 2 is a protein called ZINC FINGER PROTEIN GLI1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	9	Total	C	N	O	0	0	0
			65	41	11	13			
2	F	9	Total	C	N	O	0	0	0
			65	41	11	13			
2	G	8	Total	C	N	O	0	0	0
			58	37	10	11			
2	H	9	Total	C	N	O	0	0	0
			65	41	11	13			

- Molecule 3 is SUGAR (MALTOSE) (three-letter code: MAL) (formula: C<sub>12</sub>H<sub>22</sub>O<sub>11</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			23	12	11		
3	B	1	Total	C	O	0	0
			23	12	11		
3	C	1	Total	C	O	0	0
			23	12	11		
3	D	1	Total	C	O	0	0
			23	12	11		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

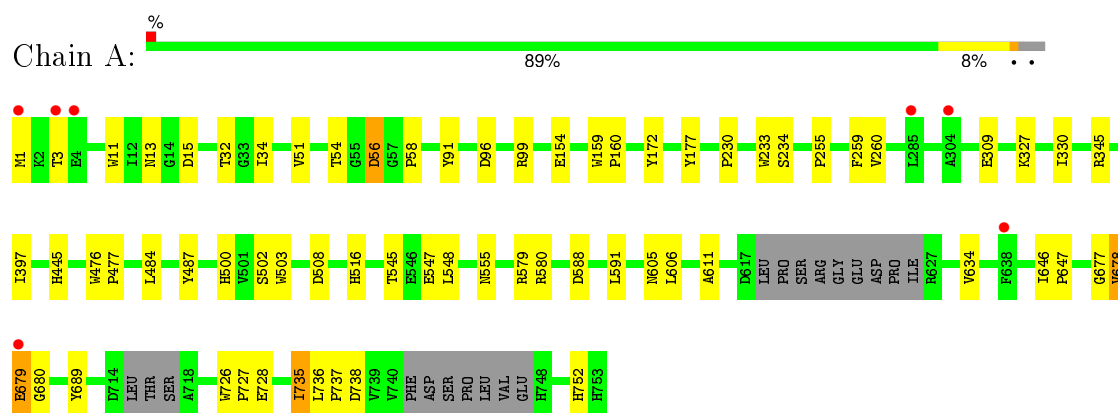
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		
4	A	1	Total	Zn	0	0
			1	1		
4	D	1	Total	Zn	0	0
			1	1		
4	C	1	Total	Zn	0	0
			1	1		



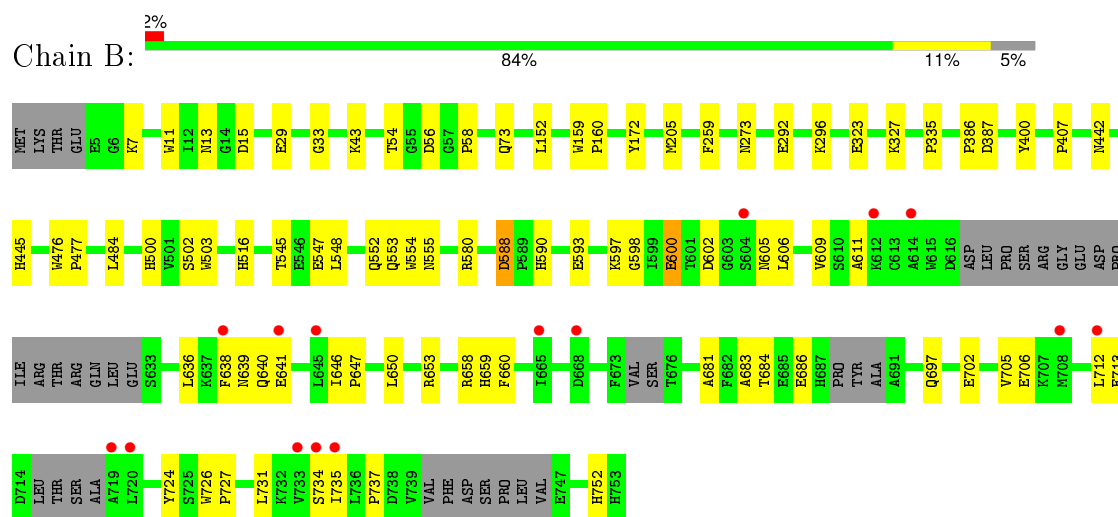
### 3 Residue-property plots [i](#)

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.

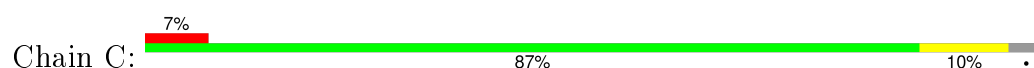
- Molecule 1: MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG

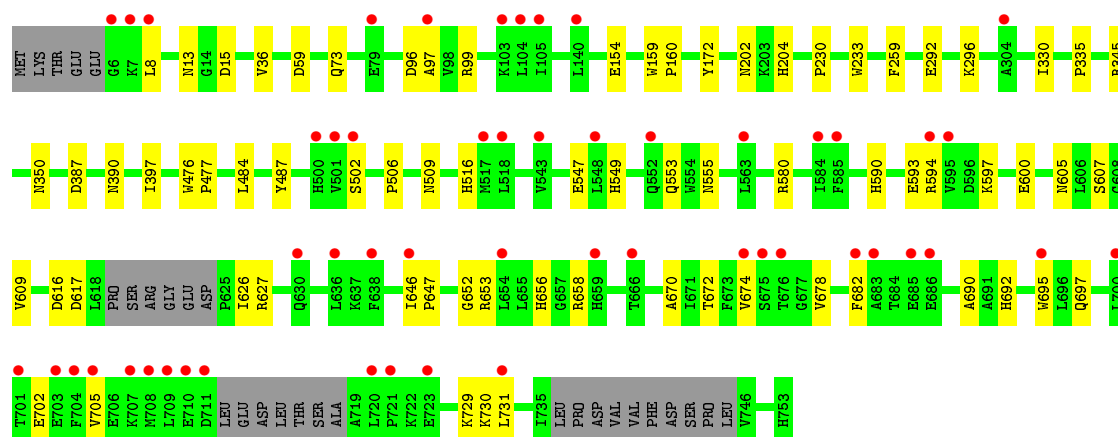


- Molecule 1: MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG

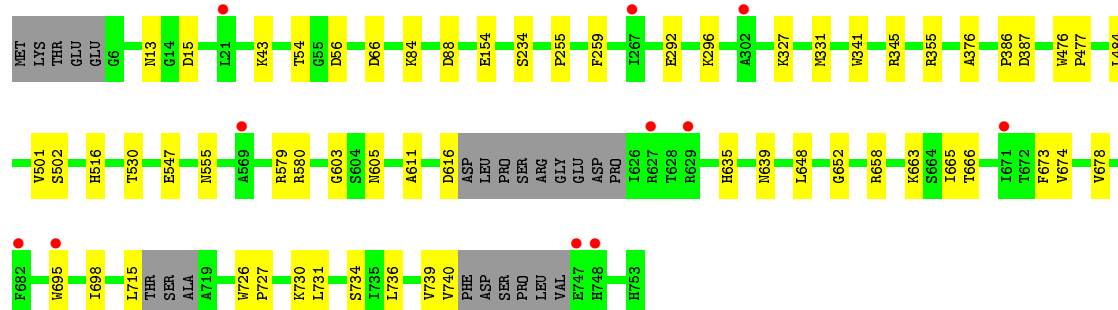
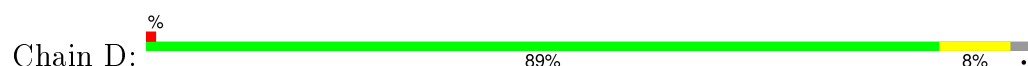


- Molecule 1: MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG

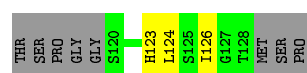
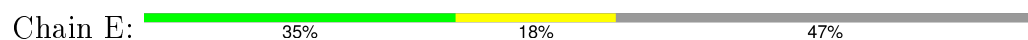




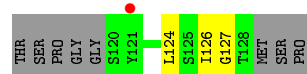
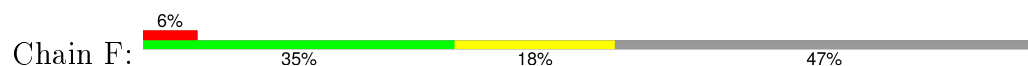
- Molecule 1: MALTOSE-BINDING PERIPLASMIC PROTEIN, SUPPRESSOR OF FUSED HOMOLOG



- Molecule 2: ZINC FINGER PROTEIN GLI1



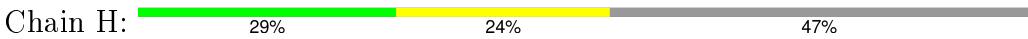
- Molecule 2: ZINC FINGER PROTEIN GLI1



- Molecule 2: ZINC FINGER PROTEIN GLI1



- Molecule 2: ZINC FINGER PROTEIN GLI1



THR	SER	PRO	GLY	GLY	S120	L124	S125	I126	G127	T128	MET	SER	PRO
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	116.30Å 137.60Å 116.54Å 90.00° 105.49° 90.00°	Depositor
Resolution (Å)	19.98 – 2.80 62.73 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.9 (19.98-2.80) 97.9 (62.73-2.80)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.60 (at 2.81Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.197 , 0.234 0.199 , 0.238	Depositor DCC
$R_{free}$ test set	1967 reflections (2.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	80.0	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 46.3	EDS
Estimated twinning fraction	0.019 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 85192 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	23141	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.12% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MAL

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	0.35	0/5913	0.51	1/8040 (0.0%)
1	B	0.38	2/5773 (0.0%)	0.51	2/7847 (0.0%)
1	C	0.36	0/5843	0.50	0/7945
1	D	0.34	0/5883	0.50	0/8001
2	E	0.36	0/66	0.63	0/88
2	F	0.40	0/66	0.59	0/88
2	G	0.32	0/59	0.49	0/78
2	H	0.46	0/66	0.75	0/88
All	All	0.36	2/23669 (0.0%)	0.51	3/32175 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	600	GLU	CG-CD	5.22	1.59	1.51
1	B	600	GLU	CD-OE1	5.18	1.31	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	588	ASP	CB-CG-OD1	-5.50	113.35	118.30
1	B	600	GLU	OE1-CD-OE2	5.14	129.47	123.30
1	A	56	ASP	CB-CG-OD1	-5.10	113.71	118.30

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	5757	0	5587	38	0
1	B	5621	0	5442	63	0
1	C	5687	0	5522	55	0
1	D	5727	0	5562	35	0
2	E	65	0	60	4	0
2	F	65	0	60	7	0
2	G	58	0	53	7	0
2	H	65	0	60	4	0
3	A	23	0	22	0	0
3	B	23	0	22	0	0
3	C	23	0	22	0	0
3	D	23	0	22	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
All	All	23141	0	22434	195	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 4.

All (195) 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:C:652:GLY:O	1:C:658:ARG:NH2	1.97	0.97
1:B:7:LYS:O	1:B:273:ASN:ND2	2.01	0.93
1:B:681:ALA:HB1	1:B:697:GLN:HE22	1.36	0.88
1:B:636:LEU:HB2	1:B:731:LEU:HD13	1.66	0.77
1:B:681:ALA:HB1	1:B:697:GLN:NE2	2.02	0.74
1:C:593:GLU:HG2	1:C:597:LYS:HE3	1.72	0.72
1:B:683:ALA:HB3	1:B:697:GLN:NE2	2.05	0.72
1:B:43:LYS:NZ	1:B:386:PRO:O	2.28	0.66
1:B:597:LYS:HA	1:B:600:GLU:HG2	1.76	0.65
1:B:639:ASN:OD1	1:B:640:GLN:N	2.28	0.65
1:D:663:LYS:HE2	1:D:695:TRP:CH2	2.31	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:547:GLU:OE2	1:A:580:ARG:NH2	2.30	0.65
1:A:54:THR:OG1	1:A:56:ASP:CG	2.36	0.65
1:A:555:ASN:HB3	1:A:605:ASN:HB2	1.79	0.63
1:A:154:GLU:OE1	1:A:345:ARG:NH1	2.32	0.63
1:C:653:ARG:HA	1:C:658:ARG:NH2	2.14	0.62
1:D:13:ASN:ND2	1:D:15:ASP:OD1	2.32	0.62
1:A:678:VAL:O	1:A:680:GLY:N	2.33	0.61
1:C:597:LYS:HA	1:C:600:GLU:OE2	2.01	0.61
1:B:638:PHE:CD2	1:B:731:LEU:HD11	2.35	0.61
1:A:51:VAL:HB	1:A:56:ASP:OD1	2.02	0.60
1:B:54:THR:OG1	1:B:56:ASP:CG	2.42	0.58
1:C:13:ASN:ND2	1:C:15:ASP:OD1	2.36	0.58
1:B:609:VAL:HG21	2:F:124:LEU:HD11	1.85	0.57
2:H:127:GLY:O	2:H:128:THR:OG1	2.15	0.57
1:A:484:LEU:HD23	2:E:126:ILE:HD11	1.86	0.57
1:B:593:GLU:HG2	1:B:597:LYS:HE3	1.86	0.56
1:D:616:ASP:O	1:D:635:HIS:N	2.38	0.56
1:C:547:GLU:OE2	1:C:580:ARG:NH2	2.39	0.56
1:B:638:PHE:HD2	1:B:731:LEU:HD11	1.70	0.55
1:B:683:ALA:HB3	1:B:697:GLN:HE21	1.70	0.55
1:D:43:LYS:NZ	1:D:386:PRO:O	2.39	0.55
1:C:96:ASP:OD1	1:C:99:ARG:NH1	2.40	0.54
1:B:73:GLN:OE1	1:B:335:PRO:HB3	2.07	0.54
1:B:54:THR:OG1	1:B:56:ASP:OD1	2.26	0.54
1:C:609:VAL:HG21	2:G:124:LEU:HD11	1.90	0.54
1:B:555:ASN:HB3	1:B:605:ASN:HB2	1.88	0.54
1:B:552:GLN:O	1:B:606:LEU:HD13	2.07	0.53
1:B:554:TRP:CE2	1:B:598:GLY:HA3	2.42	0.53
1:A:54:THR:OG1	1:A:56:ASP:OD1	2.26	0.53
1:C:656:HIS:CB	1:C:658:ARG:NH1	2.72	0.53
1:B:597:LYS:HA	1:B:600:GLU:CG	2.39	0.53
1:C:674:VAL:HG13	1:C:678:VAL:HG21	1.89	0.53
1:C:154:GLU:OE1	1:C:345:ARG:NH1	2.42	0.53
1:B:547:GLU:OE2	1:B:580:ARG:NH2	2.42	0.53
1:C:8:LEU:HD23	1:C:36:VAL:HG22	1.90	0.53
1:C:670:ALA:HB3	1:C:695:TRP:CE3	2.43	0.53
1:D:555:ASN:HB3	1:D:605:ASN:HB2	1.91	0.52
1:B:724:TYR:HB3	1:B:726:TRP:CH2	2.44	0.52
1:C:484:LEU:HD23	2:G:126:ILE:CD1	2.39	0.52
1:C:653:ARG:HG2	1:C:658:ARG:HE	1.73	0.52
1:B:13:ASN:ND2	1:B:15:ASP:OD1	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:323:GLU:HG2	1:B:327:LYS:NZ	2.24	0.52
1:A:91:TYR:OH	1:A:309:GLU:OE1	2.28	0.51
1:C:506:PRO:HG2	1:C:509:ASN:HA	1.93	0.51
1:D:292:GLU:HG2	1:D:296:LYS:HE3	1.92	0.51
1:B:602:ASP:O	1:B:658:ARG:NH2	2.43	0.51
1:C:590:HIS:CE1	1:C:593:GLU:OE1	2.64	0.51
1:A:736:LEU:N	1:A:737:PRO:HD3	2.26	0.51
1:C:502:SER:HA	1:C:516:HIS:HD2	1.77	0.50
1:C:502:SER:OG	2:G:125:SER:N	2.40	0.50
1:C:682:PHE:CE2	1:C:692:HIS:NE2	2.79	0.50
1:C:555:ASN:HB3	1:C:605:ASN:HB2	1.93	0.49
1:D:255:PRO:HB3	1:D:327:LYS:HD3	1.94	0.49
1:A:445:HIS:HA	1:A:752:HIS:HB3	1.94	0.49
1:C:616:ASP:OD1	1:C:617:ASP:N	2.41	0.49
1:C:656:HIS:CG	1:C:658:ARG:NH1	2.81	0.49
1:B:502:SER:HA	1:B:516:HIS:HD2	1.78	0.48
1:D:674:VAL:CG1	1:D:678:VAL:HG21	2.43	0.48
1:B:484:LEU:HD23	2:F:126:ILE:HD11	1.95	0.48
1:B:735:ILE:HG23	1:B:735:ILE:O	2.13	0.48
1:B:683:ALA:CB	1:B:697:GLN:NE2	2.77	0.48
1:B:684:THR:OG1	1:B:686:GLU:HB2	2.13	0.48
1:C:670:ALA:CB	1:C:695:TRP:CE3	2.96	0.48
1:C:682:PHE:CE1	1:C:690:ALA:HB3	2.49	0.48
1:D:739:VAL:HG23	1:D:740:VAL:HG23	1.96	0.47
1:B:606:LEU:HD12	1:B:606:LEU:N	2.28	0.47
1:B:588:ASP:OD1	1:B:590:HIS:CD2	2.68	0.47
1:A:646:ILE:HB	1:A:647:PRO:HD3	1.96	0.47
1:C:73:GLN:OE1	1:C:335:PRO:CG	2.63	0.47
1:C:626:ILE:HG22	1:C:627:ARG:N	2.30	0.47
1:C:292:GLU:HG2	1:C:296:LYS:HE3	1.96	0.47
1:C:230:PRO:HA	1:C:233:TRP:CE2	2.49	0.47
1:B:29:GLU:O	1:B:33:GLY:N	2.41	0.47
1:A:13:ASN:ND2	1:A:15:ASP:OD1	2.47	0.46
1:B:611:ALA:HB2	2:F:124:LEU:HD12	1.97	0.46
1:D:484:LEU:HD23	2:H:126:ILE:HD11	1.97	0.46
1:D:84:LYS:NZ	1:D:88:ASP:OD1	2.46	0.46
1:C:702:GLU:HA	1:C:705:VAL:HB	1.97	0.46
1:A:11:TRP:CE2	1:A:58:PRO:HG3	2.50	0.46
1:C:292:GLU:O	1:C:296:LYS:HG3	2.16	0.46
1:C:730:LYS:C	1:C:731:LEU:HD12	2.36	0.46
1:D:547:GLU:OE2	1:D:580:ARG:NH2	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:653:ARG:NE	1:B:659:HIS:O	2.40	0.46
1:C:672:THR:HB	1:C:697:GLN:HG2	1.96	0.46
1:B:323:GLU:CG	1:B:327:LYS:NZ	2.79	0.46
1:B:152:LEU:HD11	1:B:205:MET:HE3	1.98	0.46
1:A:735:ILE:HG22	1:A:735:ILE:O	2.16	0.45
1:C:656:HIS:CG	1:C:658:ARG:HH12	2.35	0.45
1:D:476:TRP:N	1:D:477:PRO:CD	2.80	0.45
1:B:597:LYS:CA	1:B:600:GLU:HG2	2.46	0.45
1:C:476:TRP:N	1:C:477:PRO:CD	2.80	0.45
1:C:594:ARG:HA	1:C:597:LYS:HB2	1.98	0.45
1:D:674:VAL:HG13	1:D:678:VAL:HG21	1.97	0.45
1:B:641:GLU:HG2	2:F:124:LEU:HD13	1.99	0.45
1:C:387:ASP:N	1:C:387:ASP:OD1	2.50	0.45
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.52	0.45
1:A:96:ASP:OD1	1:A:99:ARG:NH1	2.49	0.45
1:C:626:ILE:HG22	1:C:627:ARG:HG3	1.98	0.45
1:D:730:LYS:O	1:D:731:LEU:HD12	2.17	0.45
1:A:159:TRP:N	1:A:160:PRO:CD	2.80	0.45
1:B:683:ALA:H	1:B:697:GLN:NE2	2.13	0.44
1:D:736:LEU:O	1:D:739:VAL:HG22	2.18	0.44
1:B:650:LEU:HB3	1:B:705:VAL:HG13	1.98	0.44
1:B:503:TRP:CH2	2:F:127:GLY:HA3	2.52	0.44
1:B:712:LEU:HD12	1:B:713:GLU:N	2.32	0.44
1:C:159:TRP:N	1:C:160:PRO:CD	2.81	0.44
1:D:726:TRP:HB3	1:D:727:PRO:HD2	1.99	0.44
1:D:639:ASN:HA	1:D:734:SER:O	2.17	0.44
1:C:590:HIS:O	1:C:593:GLU:HB3	2.17	0.44
1:B:387:ASP:OD1	1:B:387:ASP:N	2.51	0.44
1:C:653:ARG:HA	1:C:658:ARG:HH21	1.81	0.44
1:D:501:VAL:HG11	2:H:126:ILE:HD11	1.99	0.43
1:D:54:THR:OG1	1:D:56:ASP:OD1	2.34	0.43
1:D:663:LYS:NZ	1:D:695:TRP:HH2	2.16	0.43
1:C:729:LYS:HG2	1:C:731:LEU:CD1	2.48	0.43
1:A:611:ALA:HB2	2:E:124:LEU:HD12	1.99	0.43
1:A:260:VAL:HB	1:A:330:ILE:HA	2.00	0.43
1:D:66:ASP:CB	1:D:331:MET:HE2	2.48	0.43
1:D:387:ASP:N	1:D:387:ASP:OD1	2.49	0.43
1:C:97:ALA:HB2	1:C:330:ILE:HD11	2.01	0.43
1:A:555:ASN:HB3	1:A:605:ASN:CB	2.48	0.43
1:B:588:ASP:OD1	1:B:588:ASP:C	2.57	0.43
1:B:11:TRP:CD2	1:B:58:PRO:HG3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:476:TRP:N	1:A:477:PRO:CD	2.82	0.43
1:D:673:PHE:CE2	1:D:698:ILE:HD11	2.53	0.43
1:D:611:ALA:HB2	2:H:124:LEU:HD12	2.00	0.43
1:A:502:SER:HA	1:A:516:HIS:HD2	1.84	0.43
1:C:202:ASN:HB2	1:C:204:HIS:CD2	2.53	0.43
1:C:607:SER:N	2:G:120:SER:O	2.49	0.43
1:D:331:MET:HE1	1:D:341:TRP:HE1	1.84	0.43
1:C:484:LEU:HD23	2:G:126:ILE:HD11	2.01	0.42
1:D:603:GLY:HA3	1:D:658:ARG:NH2	2.34	0.42
1:B:476:TRP:N	1:B:477:PRO:CD	2.82	0.42
1:B:653:ARG:HD3	1:B:660:PHE:HB2	2.00	0.42
1:B:11:TRP:CE2	1:B:58:PRO:HG3	2.54	0.42
1:D:665:ILE:HG13	1:D:666:THR:HG23	2.01	0.42
1:A:255:PRO:HB3	1:A:327:LYS:HD3	2.01	0.42
1:A:677:GLY:O	1:A:679:GLU:N	2.52	0.42
1:A:503:TRP:H	1:A:516:HIS:HD2	1.67	0.42
1:B:646:ILE:HB	1:B:647:PRO:HD3	2.01	0.42
1:C:476:TRP:CG	1:C:477:PRO:HD3	2.54	0.42
1:B:292:GLU:HG2	1:B:296:LYS:HE3	2.02	0.42
1:A:736:LEU:N	1:A:737:PRO:CD	2.83	0.42
1:A:516:HIS:CE1	1:A:545:THR:HG22	2.55	0.42
1:A:32:THR:HG22	1:A:34:ILE:HD13	2.02	0.42
1:D:502:SER:HA	1:D:516:HIS:HD2	1.84	0.42
1:B:516:HIS:CE1	1:B:545:THR:HG22	2.55	0.42
1:B:500:HIS:HD2	1:B:548:LEU:HD11	1.84	0.42
2:E:123:HIS:HB3	2:E:126:ILE:HG21	2.02	0.41
1:C:350:ASN:OD1	1:C:390:ASN:ND2	2.45	0.41
1:C:487:TYR:CE1	2:G:126:ILE:HG22	2.55	0.41
1:D:154:GLU:OE1	1:D:345:ARG:NH1	2.52	0.41
1:B:641:GLU:OE2	2:F:124:LEU:HD22	2.21	0.41
1:D:665:ILE:HG13	1:D:666:THR:N	2.35	0.41
1:C:397:ILE:H	1:C:397:ILE:HD12	1.84	0.41
1:B:445:HIS:HA	1:B:752:HIS:HB3	2.02	0.41
1:B:554:TRP:NE1	1:B:598:GLY:HA3	2.35	0.41
1:D:579:ARG:NE	1:D:579:ARG:HA	2.35	0.41
1:A:588:ASP:HB3	1:A:591:LEU:CD1	2.50	0.41
1:D:331:MET:HE3	1:D:341:TRP:HZ2	1.86	0.41
1:A:397:ILE:HD12	1:A:397:ILE:H	1.86	0.41
1:B:726:TRP:HB3	1:B:727:PRO:HD2	2.02	0.41
1:B:442:ASN:HA	1:B:445:HIS:O	2.20	0.41
1:A:634:VAL:HG21	1:A:689:TYR:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:GLN:OE1	1:B:335:PRO:CB	2.68	0.41
2:F:126:ILE:HG13	2:F:127:GLY:N	2.36	0.41
1:A:476:TRP:CZ2	1:A:508:ASP:HB3	2.56	0.41
1:A:487:TYR:CD1	2:E:126:ILE:HB	2.56	0.41
1:B:400:TYR:HB2	1:B:407:PRO:HA	2.03	0.41
1:A:500:HIS:HE1	1:A:606:LEU:HD21	1.85	0.41
1:C:8:LEU:HD12	1:C:59:ASP:HB2	2.04	0.40
1:A:516:HIS:CD2	1:A:548:LEU:HD22	2.57	0.40
1:B:734:SER:HB2	1:B:737:PRO:HG3	2.03	0.40
1:C:656:HIS:HB2	1:C:658:ARG:NH1	2.36	0.40
1:C:646:ILE:HB	1:C:647:PRO:HD3	2.04	0.40
1:B:159:TRP:N	1:B:160:PRO:CD	2.84	0.40
2:G:126:ILE:HG13	2:G:127:GLY:N	2.37	0.40
1:C:506:PRO:O	1:C:509:ASN:N	2.48	0.40
1:C:549:HIS:O	1:C:553:GLN:HG2	2.21	0.40
1:D:376:ALA:CB	1:D:530:THR:HG22	2.51	0.40
1:A:172:TYR:HD2	1:A:177:TYR:CZ	2.38	0.40
1:D:648:LEU:O	1:D:652:GLY:N	2.53	0.40
1:B:702:GLU:OE2	1:B:706:GLU:HG3	2.21	0.40
1:A:726:TRP:HB3	1:A:727:PRO:HD2	2.02	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	726/753 (96%)	704 (97%)	18 (2%)	4 (1%)	30	65
1	B	705/753 (94%)	687 (97%)	18 (3%)	0	100	100
1	C	717/753 (95%)	700 (98%)	17 (2%)	0	100	100
1	D	722/753 (96%)	706 (98%)	16 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	7/17 (41%)	7 (100%)	0	0	100	100
2	F	7/17 (41%)	7 (100%)	0	0	100	100
2	G	6/17 (35%)	6 (100%)	0	0	100	100
2	H	7/17 (41%)	7 (100%)	0	0	100	100
All	All	2897/3080 (94%)	2824 (98%)	69 (2%)	4 (0%)	56	87

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	679	GLU
1	A	678	VAL
1	A	738	ASP
1	A	735	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	608/626 (97%)	602 (99%)	6 (1%)	82	96
1	B	593/626 (95%)	590 (100%)	3 (0%)	92	98
1	C	600/626 (96%)	598 (100%)	2 (0%)	94	99
1	D	605/626 (97%)	601 (99%)	4 (1%)	88	97
2	E	7/13 (54%)	7 (100%)	0	100	100
2	F	7/13 (54%)	7 (100%)	0	100	100
2	G	6/13 (46%)	5 (83%)	1 (17%)	3	8
2	H	7/13 (54%)	7 (100%)	0	100	100
All	All	2433/2556 (95%)	2417 (99%)	16 (1%)	88	97

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	3	THR
1	A	234	SER
1	A	259	PHE
1	A	579	ARG
1	A	728	GLU
1	B	172	TYR
1	B	259	PHE
1	B	553	GLN
1	C	172	TYR
1	C	259	PHE
1	D	234	SER
1	D	259	PHE
1	D	355	ARG
1	D	715	LEU
2	G	124	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	500	HIS
1	A	516	HIS
1	B	516	HIS
1	B	590	HIS
1	B	697	GLN
1	C	516	HIS
1	D	500	HIS
1	D	516	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

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z  > 2$
3	MAL	A	900	-	24,24,24	0.59	0	35,35,35	0.61	0
3	MAL	B	900	-	24,24,24	0.60	0	35,35,35	0.59	0
3	MAL	C	900	-	24,24,24	0.57	0	35,35,35	0.73	0
3	MAL	D	900	-	24,24,24	0.54	0	35,35,35	0.87	1 (2%)

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
3	MAL	A	900	-	-	0/8/48/48	0/2/2/2
3	MAL	B	900	-	-	0/8/48/48	0/2/2/2
3	MAL	C	900	-	-	0/8/48/48	0/2/2/2
3	MAL	D	900	-	-	0/8/48/48	0/2/2/2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	900	MAL	C1'-C2'-C3'	2.54	114.20	110.43

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 ⓘ

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	734/753 (97%)	-0.27	7 (0%) 84 77	50, 81, 127, 176	0
1	B	717/753 (95%)	-0.15	15 (2%) 67 56	51, 87, 167, 209	0
1	C	725/753 (96%)	0.11	52 (7%) 18 10	64, 111, 193, 274	0
1	D	730/753 (96%)	-0.24	11 (1%) 76 68	56, 87, 141, 190	0
2	E	9/17 (52%)	-0.11	0 100 100	61, 71, 105, 107	0
2	F	9/17 (52%)	0.33	1 (11%) 7 3	87, 99, 123, 125	0
2	G	8/17 (47%)	0.37	0 100 100	91, 104, 111, 133	0
2	H	9/17 (52%)	0.23	0 100 100	78, 80, 107, 108	0
All	All	2941/3080 (95%)	-0.13	86 (2%) 55 43	50, 90, 167, 274	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	703	GLU	7.6
1	A	1	MET	7.6
1	C	8	LEU	6.1
1	C	710	GLU	6.1
1	C	548	LEU	5.5
1	C	708	MET	5.4
1	C	674	VAL	5.4
1	C	701	THR	5.4
1	B	720	LEU	5.3
1	B	712	LEU	5.2
1	C	595	VAL	5.0
1	C	552	GLN	4.7
1	C	501	VAL	4.5
1	C	630	GLN	4.4
1	C	700	LEU	4.0
1	C	704	PHE	4.0

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Mol	Chain	Res	Type	RSRZ
1	C	103	LYS	3.9
1	C	584	ILE	3.9
1	C	585	PHE	3.9
1	C	105	ILE	3.9
1	C	721	PRO	3.9
1	C	705	VAL	3.8
1	B	708	MET	3.8
1	C	695	TRP	3.7
1	C	518	LEU	3.7
1	A	3	THR	3.6
1	C	711	ASP	3.5
1	D	627	ARG	3.5
1	B	645	LEU	3.4
1	C	104	LEU	3.4
1	C	709	LEU	3.4
1	C	594	ARG	3.3
1	C	682	PHE	3.2
1	C	500	HIS	3.2
1	B	668	ASP	3.1
1	C	686	GLU	3.1
1	C	6	GLY	3.0
1	D	747	GLU	3.0
1	C	659	HIS	3.0
1	C	723	GLU	2.9
1	C	707	LYS	2.9
1	A	4	GLU	2.8
1	C	7	LYS	2.8
1	C	79	GLU	2.8
1	C	676	THR	2.7
1	C	720	LEU	2.7
1	A	285	LEU	2.6
1	D	21	LEU	2.6
1	B	665	ILE	2.5
1	C	731	LEU	2.5
1	C	654	LEU	2.5
1	D	569	ALA	2.4
1	D	695	TRP	2.4
1	C	685	GLU	2.4
1	C	675	SER	2.4
1	B	734	SER	2.4
1	C	563	LEU	2.3
1	A	304	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	140	LEU	2.3
1	B	604	SER	2.3
1	B	719	ALA	2.3
1	D	748	HIS	2.3
1	C	97	ALA	2.3
1	C	517	MET	2.2
1	A	679	GLU	2.2
1	B	614	ALA	2.2
1	C	304	ALA	2.2
1	B	638	PHE	2.2
1	C	683	ALA	2.2
1	C	666	THR	2.2
1	C	646	ILE	2.2
2	F	121	TYR	2.2
1	C	502	SER	2.2
1	B	735	ILE	2.2
1	D	671	ILE	2.2
1	B	612	LYS	2.1
1	D	629	ARG	2.1
1	D	267	ILE	2.1
1	C	636	LEU	2.1
1	D	682	PHE	2.1
1	A	638	PHE	2.1
1	C	543	VAL	2.1
1	C	638	PHE	2.1
1	B	641	GLU	2.0
1	D	302	ALA	2.0
1	B	733	VAL	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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	ZN	D	910	1/1	0.99	0.22	3.52	93,93,93,93	0
4	ZN	C	910	1/1	0.99	0.26	2.34	84,84,84,84	0
4	ZN	B	910	1/1	0.98	0.23	1.15	98,98,98,98	0
4	ZN	A	910	1/1	0.96	0.18	0.85	75,75,75,75	0
3	MAL	C	900	23/23	0.96	0.15	-0.47	63,79,91,98	0
3	MAL	D	900	23/23	0.95	0.15	-0.49	55,64,72,72	0
3	MAL	A	900	23/23	0.96	0.15	-0.83	55,63,73,82	0
3	MAL	B	900	23/23	0.95	0.16	-0.85	47,57,67,71	0

## 6.5 Other polymers

There are no such residues in this entry.