



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:06 AM GMT

PDB ID : 3D4G
Title : ZP-N domain of mammalian sperm receptor ZP3 (crystal form II)
Authors : Jovine, L.; Monne, M.
Deposited on : 2008-05-14
Resolution : 2.30 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 (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

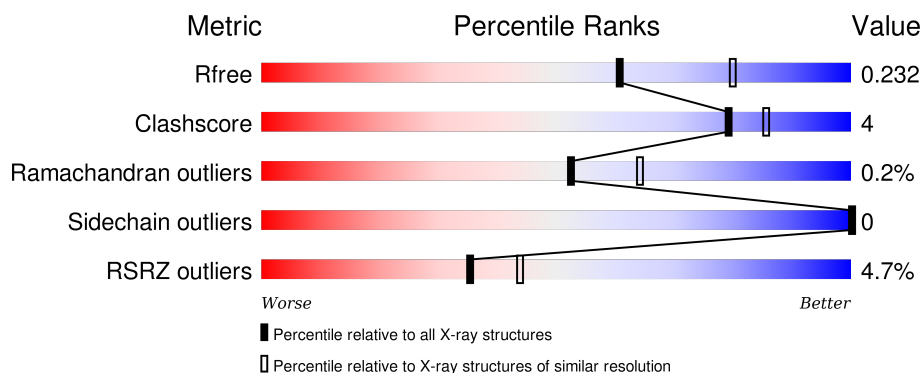
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	481	<div> <div>4%</div> <div>90%</div> <div>7%</div> <div>••</div> </div>
1	B	481	<div> <div>6%</div> <div>88%</div> <div>10%</div> <div>•</div> </div>
1	C	481	<div> <div>4%</div> <div>89%</div> <div>9%</div> <div>•</div> </div>
1	D	481	<div> <div>5%</div> <div>91%</div> <div>7%</div> <div>•</div> </div>
1	E	481	<div> <div>4%</div> <div>91%</div> <div>7%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	481	 4% 90% 8% •
1	G	481	 5% 90% 8% •
1	H	481	 5% 88% 9% •

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 31230 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, LINKER, Zona pellucida protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	471	Total	C	N	O	S	0	0	0
			3637	2325	605	696	11			
1	B	471	Total	C	N	O	S	0	0	0
			3637	2325	605	696	11			
1	C	471	Total	C	N	O	S	0	0	0
			3637	2325	605	696	11			
1	D	469	Total	C	N	O	S	0	0	0
			3621	2316	603	691	11			
1	E	472	Total	C	N	O	S	0	0	0
			3645	2331	606	697	11			
1	F	472	Total	C	N	O	S	0	0	0
			3645	2331	606	697	11			
1	G	472	Total	C	N	O	S	0	0	0
			3645	2331	606	697	11			
1	H	469	Total	C	N	O	S	0	0	0
			3621	2316	603	691	11			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	INITIATING METHIONINE	UNP P0AEX9
A	3	THR	ILE	ENGINEERED	UNP P0AEX9
A	360	ALA	GLU	ENGINEERED	UNP P0AEX9
A	363	ALA	LYS	ENGINEERED	UNP P0AEX9
A	364	ALA	ASP	ENGINEERED	UNP P0AEX9
A	368	ASN	ARG	ENGINEERED	UNP P0AEX9
A	474	LEU	-	EXPRESSION TAG	UNP P10761
A	475	GLU	-	EXPRESSION TAG	UNP P10761
A	476	HIS	-	EXPRESSION TAG	UNP P10761
A	477	HIS	-	EXPRESSION TAG	UNP P10761
A	478	HIS	-	EXPRESSION TAG	UNP P10761
A	479	HIS	-	EXPRESSION TAG	UNP P10761

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Chain	Residue	Modelled	Actual	Comment	Reference
A	480	HIS	-	EXPRESSION TAG	UNP P10761
A	481	HIS	-	EXPRESSION TAG	UNP P10761
B	1	MET	-	INITIATING METHIONINE	UNP P0AEX9
B	3	THR	ILE	ENGINEERED	UNP P0AEX9
B	360	ALA	GLU	ENGINEERED	UNP P0AEX9
B	363	ALA	LYS	ENGINEERED	UNP P0AEX9
B	364	ALA	ASP	ENGINEERED	UNP P0AEX9
B	368	ASN	ARG	ENGINEERED	UNP P0AEX9
B	474	LEU	-	EXPRESSION TAG	UNP P10761
B	475	GLU	-	EXPRESSION TAG	UNP P10761
B	476	HIS	-	EXPRESSION TAG	UNP P10761
B	477	HIS	-	EXPRESSION TAG	UNP P10761
B	478	HIS	-	EXPRESSION TAG	UNP P10761
B	479	HIS	-	EXPRESSION TAG	UNP P10761
B	480	HIS	-	EXPRESSION TAG	UNP P10761
B	481	HIS	-	EXPRESSION TAG	UNP P10761
C	1	MET	-	INITIATING METHIONINE	UNP P0AEX9
C	3	THR	ILE	ENGINEERED	UNP P0AEX9
C	360	ALA	GLU	ENGINEERED	UNP P0AEX9
C	363	ALA	LYS	ENGINEERED	UNP P0AEX9
C	364	ALA	ASP	ENGINEERED	UNP P0AEX9
C	368	ASN	ARG	ENGINEERED	UNP P0AEX9
C	474	LEU	-	EXPRESSION TAG	UNP P10761
C	475	GLU	-	EXPRESSION TAG	UNP P10761
C	476	HIS	-	EXPRESSION TAG	UNP P10761
C	477	HIS	-	EXPRESSION TAG	UNP P10761
C	478	HIS	-	EXPRESSION TAG	UNP P10761
C	479	HIS	-	EXPRESSION TAG	UNP P10761
C	480	HIS	-	EXPRESSION TAG	UNP P10761
C	481	HIS	-	EXPRESSION TAG	UNP P10761
D	1	MET	-	INITIATING METHIONINE	UNP P0AEX9
D	3	THR	ILE	ENGINEERED	UNP P0AEX9
D	360	ALA	GLU	ENGINEERED	UNP P0AEX9
D	363	ALA	LYS	ENGINEERED	UNP P0AEX9
D	364	ALA	ASP	ENGINEERED	UNP P0AEX9
D	368	ASN	ARG	ENGINEERED	UNP P0AEX9
D	474	LEU	-	EXPRESSION TAG	UNP P10761
D	475	GLU	-	EXPRESSION TAG	UNP P10761
D	476	HIS	-	EXPRESSION TAG	UNP P10761
D	477	HIS	-	EXPRESSION TAG	UNP P10761
D	478	HIS	-	EXPRESSION TAG	UNP P10761
D	479	HIS	-	EXPRESSION TAG	UNP P10761

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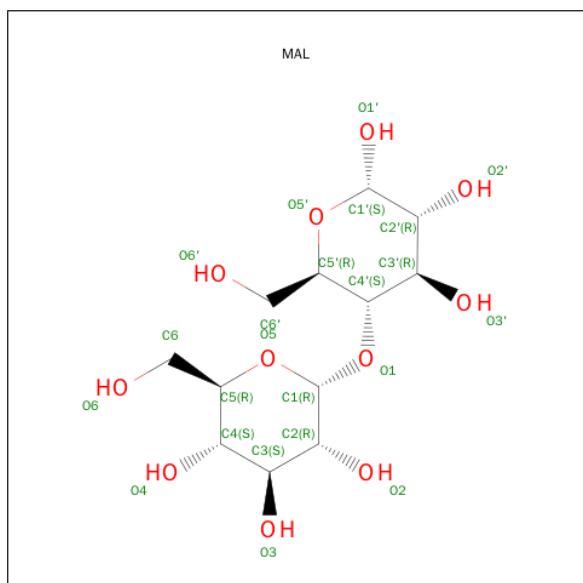
Chain	Residue	Modelled	Actual	Comment	Reference
D	480	HIS	-	EXPRESSION TAG	UNP P10761
D	481	HIS	-	EXPRESSION TAG	UNP P10761
E	1	MET	-	INITIATING METHIONINE	UNP P0AEX9
E	3	THR	ILE	ENGINEERED	UNP P0AEX9
E	360	ALA	GLU	ENGINEERED	UNP P0AEX9
E	363	ALA	LYS	ENGINEERED	UNP P0AEX9
E	364	ALA	ASP	ENGINEERED	UNP P0AEX9
E	368	ASN	ARG	ENGINEERED	UNP P0AEX9
E	474	LEU	-	EXPRESSION TAG	UNP P10761
E	475	GLU	-	EXPRESSION TAG	UNP P10761
E	476	HIS	-	EXPRESSION TAG	UNP P10761
E	477	HIS	-	EXPRESSION TAG	UNP P10761
E	478	HIS	-	EXPRESSION TAG	UNP P10761
E	479	HIS	-	EXPRESSION TAG	UNP P10761
E	480	HIS	-	EXPRESSION TAG	UNP P10761
E	481	HIS	-	EXPRESSION TAG	UNP P10761
F	1	MET	-	INITIATING METHIONINE	UNP P0AEX9
F	3	THR	ILE	ENGINEERED	UNP P0AEX9
F	360	ALA	GLU	ENGINEERED	UNP P0AEX9
F	363	ALA	LYS	ENGINEERED	UNP P0AEX9
F	364	ALA	ASP	ENGINEERED	UNP P0AEX9
F	368	ASN	ARG	ENGINEERED	UNP P0AEX9
F	474	LEU	-	EXPRESSION TAG	UNP P10761
F	475	GLU	-	EXPRESSION TAG	UNP P10761
F	476	HIS	-	EXPRESSION TAG	UNP P10761
F	477	HIS	-	EXPRESSION TAG	UNP P10761
F	478	HIS	-	EXPRESSION TAG	UNP P10761
F	479	HIS	-	EXPRESSION TAG	UNP P10761
F	480	HIS	-	EXPRESSION TAG	UNP P10761
F	481	HIS	-	EXPRESSION TAG	UNP P10761
G	1	MET	-	INITIATING METHIONINE	UNP P0AEX9
G	3	THR	ILE	ENGINEERED	UNP P0AEX9
G	360	ALA	GLU	ENGINEERED	UNP P0AEX9
G	363	ALA	LYS	ENGINEERED	UNP P0AEX9
G	364	ALA	ASP	ENGINEERED	UNP P0AEX9
G	368	ASN	ARG	ENGINEERED	UNP P0AEX9
G	474	LEU	-	EXPRESSION TAG	UNP P10761
G	475	GLU	-	EXPRESSION TAG	UNP P10761
G	476	HIS	-	EXPRESSION TAG	UNP P10761
G	477	HIS	-	EXPRESSION TAG	UNP P10761
G	478	HIS	-	EXPRESSION TAG	UNP P10761
G	479	HIS	-	EXPRESSION TAG	UNP P10761

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Chain	Residue	Modelled	Actual	Comment	Reference
G	480	HIS	-	EXPRESSION TAG	UNP P10761
G	481	HIS	-	EXPRESSION TAG	UNP P10761
H	1	MET	-	INITIATING METHIONINE	UNP P0AEX9
H	3	THR	ILE	ENGINEERED	UNP P0AEX9
H	360	ALA	GLU	ENGINEERED	UNP P0AEX9
H	363	ALA	LYS	ENGINEERED	UNP P0AEX9
H	364	ALA	ASP	ENGINEERED	UNP P0AEX9
H	368	ASN	ARG	ENGINEERED	UNP P0AEX9
H	474	LEU	-	EXPRESSION TAG	UNP P10761
H	475	GLU	-	EXPRESSION TAG	UNP P10761
H	476	HIS	-	EXPRESSION TAG	UNP P10761
H	477	HIS	-	EXPRESSION TAG	UNP P10761
H	478	HIS	-	EXPRESSION TAG	UNP P10761
H	479	HIS	-	EXPRESSION TAG	UNP P10761
H	480	HIS	-	EXPRESSION TAG	UNP P10761
H	481	HIS	-	EXPRESSION TAG	UNP P10761

- Molecule 2 is SUGAR (MALTOSE) (three-letter code: MAL) (formula: C₁₂H₂₂O₁₁).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			23	12	11		
2	E	1	Total	C	O	0	0
			23	12	11		
2	F	1	Total	C	O	0	0
			23	12	11		
2	G	1	Total	C	O	0	0
			23	12	11		
2	H	1	Total	C	O	0	0
			23	12	11		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	3	Total	Ca	0	0
			3	3		
3	D	3	Total	Ca	0	0
			3	3		
3	E	3	Total	Ca	0	0
			3	3		
3	H	3	Total	Ca	0	0
			3	3		
3	B	3	Total	Ca	0	0
			3	3		
3	C	3	Total	Ca	0	0
			3	3		
3	A	3	Total	Ca	0	0
			3	3		
3	F	3	Total	Ca	0	0
			3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	288	Total	O	0	0
			288	288		
4	B	205	Total	O	0	0
			205	205		
4	C	291	Total	O	0	0
			291	291		
4	D	205	Total	O	0	0
			205	205		

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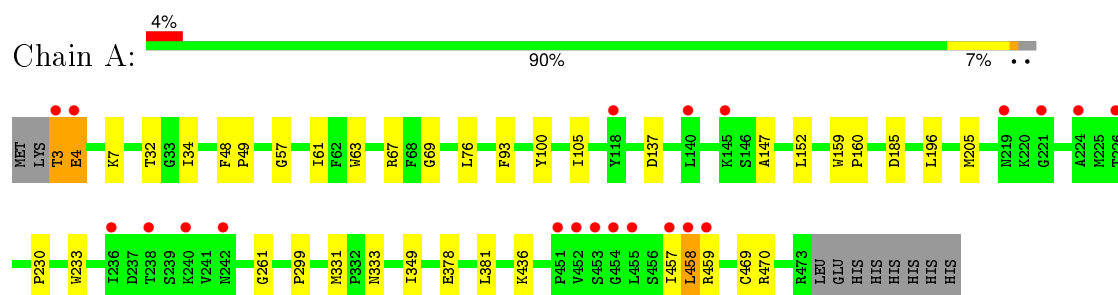
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	282	Total 282	O 282	0	0
4	F	272	Total 272	O 272	0	0
4	G	195	Total 195	O 195	0	0
4	H	196	Total 196	O 196	0	0

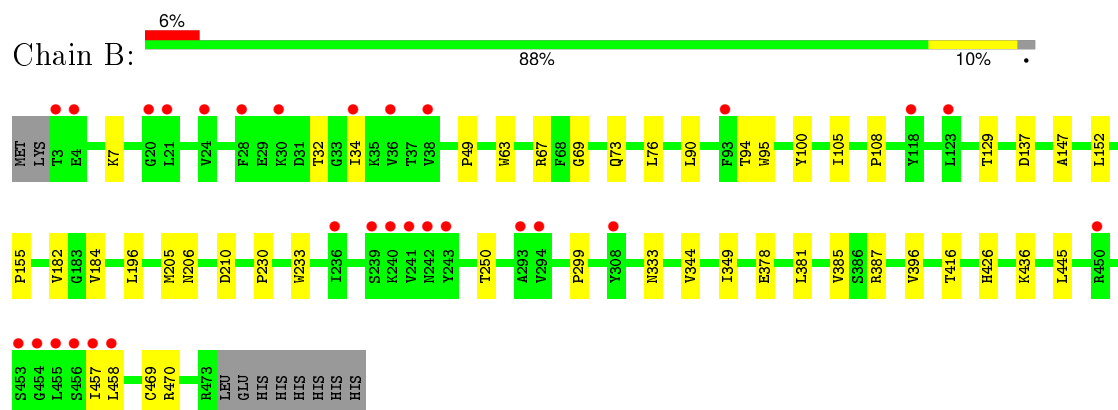
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

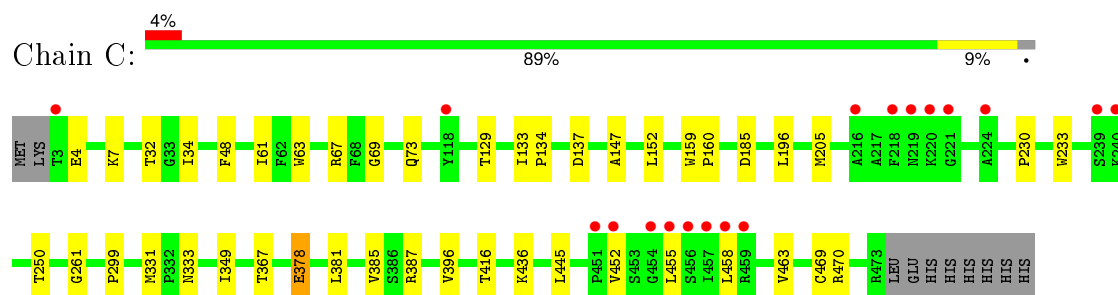
- Molecule 1: Maltose-binding periplasmic protein, LINKER, Zona pellucida protein 3



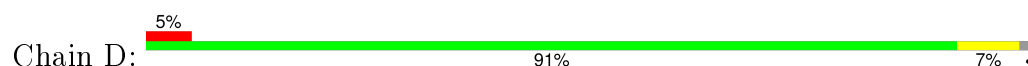
- Molecule 1: Maltose-binding periplasmic protein, LINKER, Zona pellucida protein 3

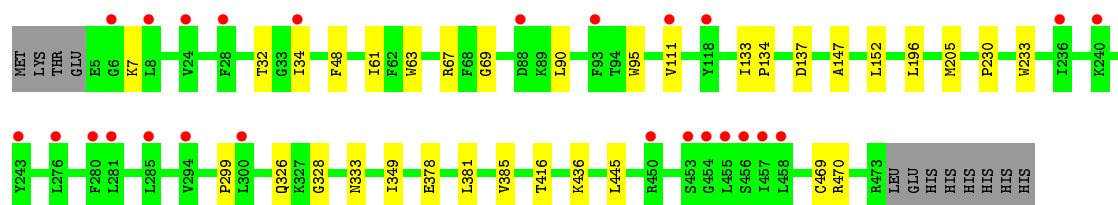


- Molecule 1: Maltose-binding periplasmic protein, LINKER, Zona pellucida protein 3

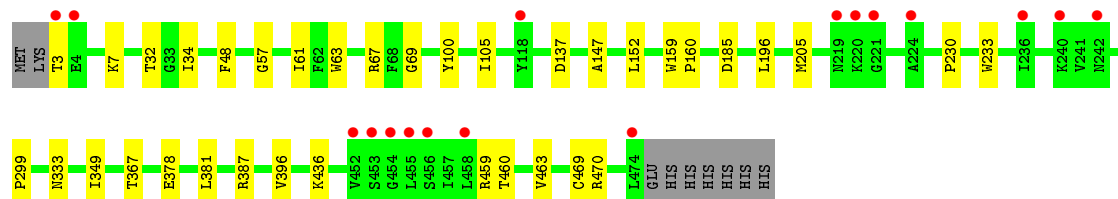
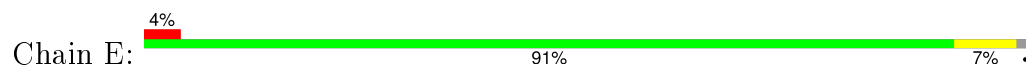


- Molecule 1: Maltose-binding periplasmic protein, LINKER, Zona pellucida protein 3

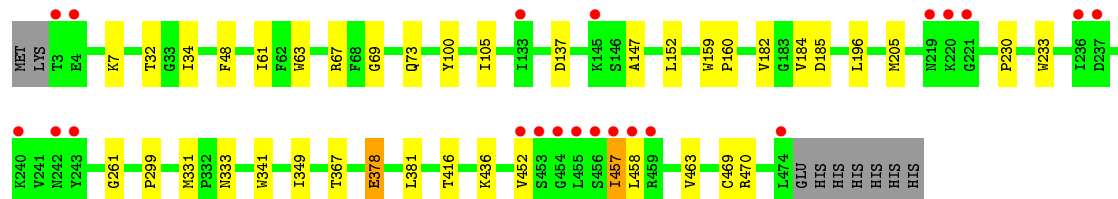
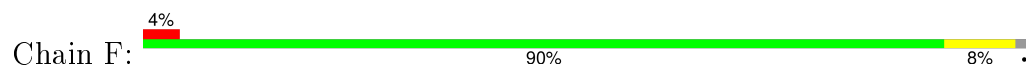




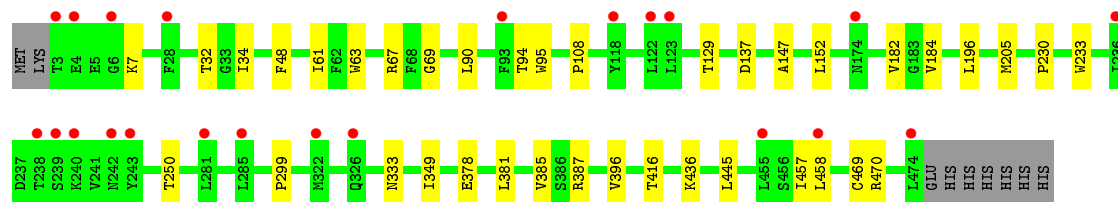
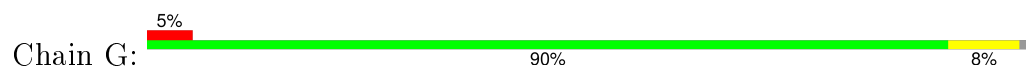
- Molecule 1: Maltose-binding periplasmic protein, LINKER, Zona pellucida protein 3



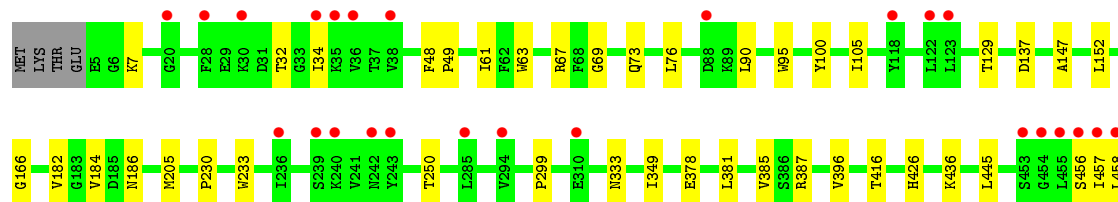
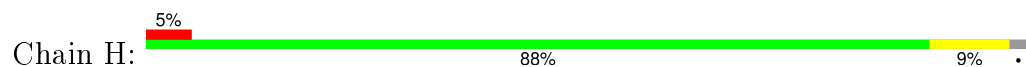
- Molecule 1: Maltose-binding periplasmic protein, LINKER, Zona pellucida protein 3



- Molecule 1: Maltose-binding periplasmic protein, LINKER, Zona pellucida protein 3



- Molecule 1: Maltose-binding periplasmic protein, LINKER, Zona pellucida protein 3



C469	LEU
R470	GLU
R473	HIS
	HIS
	HIS
	HIS
	HIS
	HIS
	HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	91.05Å 91.96Å 140.48Å 89.99° 90.18° 89.98°	Depositor
Resolution (Å)	50.00 – 2.30 47.64 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.5 (50.00-2.30) 95.4 (47.64-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.3.0038	Depositor
R, R_{free}	0.201 , 0.227 0.205 , 0.232	Depositor DCC
R_{free} test set	2139 reflections (1.12%)	DCC
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.185	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.7	EDS
Estimated twinning fraction	0.000 for -k,h,l 0.000 for k,-h,l 0.448 for h,-k,-l 0.459 for -h,k,-l 0.447 for -h,-k,l 0.000 for -k,-h,-l 0.002 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 193225 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	31230	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 57.90 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.2121e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	0.37	0/3719	0.49	0/5055
1	B	0.35	0/3719	0.49	0/5055
1	C	0.36	0/3719	0.49	0/5055
1	D	0.35	0/3703	0.48	0/5033
1	E	0.37	0/3727	0.49	0/5066
1	F	0.36	0/3727	0.49	0/5066
1	G	0.35	0/3727	0.48	0/5066
1	H	0.34	0/3703	0.48	0/5033
All	All	0.36	0/29744	0.49	0/40429

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	F	0	2
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	3	THR	Peptide
1	A	458	LEU	Peptide
1	F	452	VAL	Peptide

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Mol	Chain	Res	Type	Group
1	F	457	ILE	Peptide

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	3637	0	3610	26	0
1	B	3637	0	3610	35	0
1	C	3637	0	3610	30	0
1	D	3621	0	3597	21	0
1	E	3645	0	3621	23	0
1	F	3645	0	3621	26	0
1	G	3645	0	3621	25	0
1	H	3621	0	3597	29	0
2	A	23	0	22	0	0
2	B	23	0	22	0	0
2	C	23	0	22	0	0
2	D	23	0	22	0	0
2	E	23	0	22	0	0
2	F	23	0	22	1	0
2	G	23	0	22	0	0
2	H	23	0	22	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
3	D	3	0	0	0	0
3	E	3	0	0	0	0
3	F	3	0	0	0	0
3	G	3	0	0	0	0
3	H	3	0	0	0	0
4	A	288	0	0	2	0
4	B	205	0	0	2	0
4	C	291	0	0	2	0
4	D	205	0	0	2	0
4	E	282	0	0	2	0
4	F	272	0	0	2	0
4	G	195	0	0	2	0
4	H	196	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	31230	0	29063	204	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 (204) 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:32:THR:HG22	1:D:34:ILE:HD13	1.47	0.97
1:B:32:THR:HG22	1:B:34:ILE:HD13	1.49	0.95
1:H:32:THR:HG22	1:H:34:ILE:HD13	1.48	0.94
1:E:32:THR:HG22	1:E:34:ILE:HD13	1.48	0.93
1:A:32:THR:HG22	1:A:34:ILE:HD13	1.50	0.92
1:C:32:THR:HG22	1:C:34:ILE:HD13	1.50	0.92
1:G:32:THR:HG22	1:G:34:ILE:HD13	1.50	0.90
1:F:32:THR:HG22	1:F:34:ILE:HD13	1.53	0.89
1:F:152:LEU:HD11	1:F:205:MET:HE3	1.56	0.86
1:G:152:LEU:HD11	1:G:205:MET:HE3	1.58	0.85
1:D:152:LEU:HD11	1:D:205:MET:HE3	1.59	0.84
1:C:152:LEU:HD11	1:C:205:MET:HE3	1.61	0.82
1:B:152:LEU:HD11	1:B:205:MET:HE3	1.64	0.79
1:E:152:LEU:HD11	1:E:205:MET:HE3	1.66	0.77
1:A:152:LEU:HD11	1:A:205:MET:HE3	1.67	0.75
1:A:457:ILE:HG22	1:A:458:LEU:HG	1.68	0.75
1:H:152:LEU:HD11	1:H:205:MET:HE3	1.72	0.70
1:E:3:THR:HG23	1:E:57:GLY:O	1.96	0.66
1:B:457:ILE:HG22	1:B:458:LEU:HD13	1.80	0.64
1:B:457:ILE:HG22	1:B:458:LEU:CD1	2.28	0.63
1:G:436:LYS:NZ	4:G:578:HOH:O	2.30	0.62
1:H:436:LYS:NZ	4:H:1824:HOH:O	2.33	0.62
1:E:459:ARG:HG2	1:E:460:THR:N	2.14	0.62
1:B:436:LYS:NZ	4:B:571:HOH:O	2.33	0.62
1:C:458:LEU:HD13	1:D:111:VAL:HG21	1.82	0.60
1:F:48:PHE:CG	1:F:61:ILE:HD12	2.37	0.60
1:C:349:ILE:HD13	1:C:470:ARG:NH2	2.17	0.59
1:F:436:LYS:NZ	4:F:575:HOH:O	2.35	0.58
1:F:349:ILE:HD13	1:F:470:ARG:NH2	2.18	0.57
1:F:416:THR:HG22	1:H:73:GLN:CD	2.23	0.57
1:B:73:GLN:CD	1:C:416:THR:HG22	2.26	0.56
1:A:349:ILE:HD13	1:A:470:ARG:NH2	2.21	0.56
1:G:349:ILE:HD13	1:G:470:ARG:NH2	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:436:LYS:NZ	4:D:576:HOH:O	2.38	0.56
1:H:349:ILE:HD13	1:H:470:ARG:NH2	2.21	0.55
1:B:416:THR:HG22	1:C:73:GLN:CD	2.26	0.55
1:G:7:LYS:HA	1:G:34:ILE:HG23	1.90	0.54
1:G:152:LEU:HD21	1:G:205:MET:HE1	1.90	0.53
1:C:378:GLU:O	1:H:426:HIS:NE2	2.41	0.53
1:B:69:GLY:HA3	1:B:333:ASN:O	2.09	0.53
1:B:90:LEU:HD12	1:B:95:TRP:CZ2	2.43	0.53
1:H:457:ILE:N	1:H:457:ILE:HD12	2.23	0.53
1:G:385:VAL:HG21	1:G:445:LEU:HD21	1.91	0.53
1:A:48:PHE:CG	1:A:61:ILE:HD12	2.44	0.53
1:D:349:ILE:HD13	1:D:470:ARG:NH2	2.24	0.52
1:F:196:LEU:HD11	1:F:205:MET:HE1	1.91	0.52
1:D:7:LYS:HA	1:D:34:ILE:HG23	1.91	0.52
1:F:73:GLN:CD	1:H:416:THR:HG22	2.29	0.52
1:B:7:LYS:HA	1:B:34:ILE:HG23	1.92	0.52
1:D:137:ASP:HA	1:D:147:ALA:HB2	1.92	0.52
1:A:69:GLY:HA3	1:A:333:ASN:O	2.10	0.52
1:A:93:PHE:CD1	1:B:457:ILE:HG23	2.44	0.51
1:D:69:GLY:HA3	1:D:333:ASN:O	2.10	0.51
1:E:349:ILE:HD13	1:E:470:ARG:NH2	2.26	0.51
1:A:436:LYS:NZ	4:A:574:HOH:O	2.43	0.51
1:C:69:GLY:HA3	1:C:333:ASN:O	2.10	0.51
1:H:7:LYS:HA	1:H:34:ILE:HG23	1.93	0.51
1:B:457:ILE:HG22	1:B:458:LEU:N	2.26	0.51
1:G:137:ASP:HA	1:G:147:ALA:HB2	1.92	0.51
1:G:69:GLY:HA3	1:G:333:ASN:O	2.10	0.51
1:F:457:ILE:HG22	1:F:458:LEU:HB2	1.93	0.51
1:F:100:TYR:HB3	1:F:105:ILE:HD13	1.93	0.51
1:B:349:ILE:HD13	1:B:470:ARG:NH2	2.25	0.50
1:E:69:GLY:HA3	1:E:333:ASN:O	2.12	0.50
1:D:90:LEU:HD12	1:D:95:TRP:CZ2	2.46	0.50
1:F:69:GLY:HA3	1:F:333:ASN:O	2.12	0.50
1:B:137:ASP:HA	1:B:147:ALA:HB2	1.94	0.49
1:H:385:VAL:HG21	1:H:445:LEU:HD21	1.95	0.49
1:B:381:LEU:HD13	1:B:469:CYS:SG	2.53	0.49
1:E:48:PHE:CG	1:E:61:ILE:HD12	2.47	0.49
1:E:7:LYS:HA	1:E:34:ILE:HG23	1.94	0.49
1:C:7:LYS:HA	1:C:34:ILE:HG23	1.95	0.49
1:F:7:LYS:HA	1:F:34:ILE:HG23	1.94	0.49
1:B:182:VAL:HG12	1:B:184:VAL:HG13	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:230:PRO:HA	1:E:233:TRP:CE2	2.48	0.48
1:G:90:LEU:HD12	1:G:95:TRP:CZ2	2.48	0.48
1:D:385:VAL:HG21	1:D:445:LEU:HD21	1.96	0.48
1:A:3:THR:O	1:A:4:GLU:HB2	2.13	0.48
1:H:137:ASP:HA	1:H:147:ALA:HB2	1.95	0.48
1:H:90:LEU:HD12	1:H:95:TRP:CZ2	2.48	0.48
1:C:48:PHE:CG	1:C:61:ILE:HD12	2.49	0.47
1:E:63:TRP:CD1	1:E:67:ARG:HG3	2.49	0.47
1:B:426:HIS:NE2	1:F:378:GLU:O	2.47	0.47
1:H:69:GLY:HA3	1:H:333:ASN:O	2.13	0.47
1:A:7:LYS:HA	1:A:34:ILE:HG23	1.96	0.47
1:F:185:ASP:HB2	4:F:678:HOH:O	2.13	0.47
1:F:100:TYR:CB	1:F:105:ILE:HD13	2.44	0.47
1:D:63:TRP:CD1	1:D:67:ARG:HG3	2.50	0.47
1:A:137:ASP:HA	1:A:147:ALA:HB2	1.97	0.47
1:D:233:TRP:HB2	1:D:299:PRO:HG2	1.97	0.47
1:F:367:THR:HG21	1:F:463:VAL:HG11	1.97	0.47
1:F:230:PRO:HA	1:F:233:TRP:CE2	2.50	0.47
1:A:381:LEU:HD13	1:A:469:CYS:SG	2.55	0.47
1:F:381:LEU:HD13	1:F:469:CYS:SG	2.55	0.47
1:D:381:LEU:HD13	1:D:469:CYS:SG	2.56	0.46
1:A:100:TYR:HB3	1:A:105:ILE:HD13	1.97	0.46
1:F:261:GLY:HA2	1:F:331:MET:HE2	1.97	0.46
1:C:230:PRO:HA	1:C:233:TRP:CE2	2.50	0.46
1:B:230:PRO:HA	1:B:233:TRP:CE2	2.51	0.46
1:C:436:LYS:NZ	4:C:586:HOH:O	2.48	0.46
1:E:137:ASP:HA	1:E:147:ALA:HB2	1.97	0.46
1:G:457:ILE:HG22	1:G:458:LEU:N	2.31	0.46
1:D:230:PRO:HA	1:D:233:TRP:CE2	2.51	0.46
1:B:233:TRP:HB2	1:B:299:PRO:HG2	1.98	0.46
1:F:137:ASP:HA	1:F:147:ALA:HB2	1.98	0.46
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.51	0.45
1:B:385:VAL:HG21	1:B:445:LEU:HD21	1.99	0.45
1:E:387:ARG:HA	1:E:396:VAL:HG22	1.98	0.45
1:A:100:TYR:CB	1:A:105:ILE:HD13	2.46	0.45
1:C:133:ILE:N	1:C:134:PRO:CD	2.79	0.45
1:A:63:TRP:CD1	1:A:67:ARG:HG3	2.50	0.45
1:H:100:TYR:HB3	1:H:105:ILE:HD13	1.99	0.45
1:E:233:TRP:HB2	1:E:299:PRO:HG2	1.99	0.45
1:H:182:VAL:HG12	1:H:184:VAL:HG13	1.99	0.45
1:D:133:ILE:N	1:D:134:PRO:CD	2.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:436:LYS:NZ	4:E:1081:HOH:O	2.50	0.45
1:G:182:VAL:HG12	1:G:184:VAL:HG13	1.99	0.45
1:H:233:TRP:HB2	1:H:299:PRO:HG2	1.98	0.45
1:F:233:TRP:HB2	1:F:299:PRO:HG2	1.99	0.45
1:D:48:PHE:CG	1:D:61:ILE:HD12	2.51	0.45
1:C:261:GLY:HA2	1:C:331:MET:HE2	1.98	0.45
1:B:210:ASP:OD2	1:B:436:LYS:NZ	2.46	0.45
1:G:63:TRP:CD1	1:G:67:ARG:HG3	2.51	0.45
1:F:48:PHE:CB	1:F:61:ILE:HD12	2.46	0.44
1:A:458:LEU:O	1:A:459:ARG:C	2.55	0.44
1:G:457:ILE:HG22	1:G:458:LEU:HG	1.98	0.44
1:B:63:TRP:CD1	1:B:67:ARG:HG3	2.52	0.44
1:B:387:ARG:HA	1:B:396:VAL:HG22	1.99	0.44
1:G:230:PRO:HA	1:G:233:TRP:CE2	2.52	0.44
1:F:63:TRP:CD1	1:F:67:ARG:HG3	2.53	0.44
1:C:455:LEU:HB3	1:D:326:GLN:O	2.17	0.44
1:H:457:ILE:HG22	1:H:458:LEU:N	2.32	0.44
1:H:381:LEU:HD13	1:H:469:CYS:SG	2.58	0.44
1:C:452:VAL:O	1:C:452:VAL:HG13	2.17	0.44
1:C:367:THR:HG21	1:C:463:VAL:HG11	1.99	0.44
1:G:233:TRP:HB2	1:G:299:PRO:HG2	1.99	0.44
1:G:381:LEU:HD13	1:G:469:CYS:SG	2.58	0.44
1:C:387:ARG:HA	1:C:396:VAL:HG22	2.00	0.44
1:G:387:ARG:HA	1:G:396:VAL:HG22	2.00	0.43
1:E:152:LEU:HD21	1:E:205:MET:HE1	2.00	0.43
1:A:3:THR:HG21	1:A:57:GLY:O	2.17	0.43
1:G:48:PHE:CG	1:G:61:ILE:HD12	2.53	0.43
1:H:49:PRO:HA	1:H:76:LEU:HD13	1.99	0.43
1:C:152:LEU:HD21	1:C:205:MET:HE1	2.00	0.43
1:C:233:TRP:HB2	1:C:299:PRO:HG2	2.00	0.43
1:C:385:VAL:HG21	1:C:445:LEU:HD21	2.00	0.43
1:C:381:LEU:HD13	1:C:469:CYS:SG	2.59	0.43
1:G:129:THR:HG22	1:G:250:THR:OG1	2.18	0.43
1:C:137:ASP:HA	1:C:147:ALA:HB2	2.01	0.43
1:H:63:TRP:CD1	1:H:67:ARG:HG3	2.52	0.43
1:E:185:ASP:HB2	4:E:1195:HOH:O	2.18	0.43
1:B:94:THR:HB	1:B:108:PRO:HB2	2.00	0.43
1:D:196:LEU:HD11	1:D:205:MET:HE1	2.01	0.43
1:A:233:TRP:HB2	1:A:299:PRO:HG2	2.01	0.43
1:A:261:GLY:HA2	1:A:331:MET:HE2	2.00	0.43
1:E:381:LEU:HD13	1:E:469:CYS:SG	2.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:TRP:CD1	1:C:67:ARG:HG3	2.52	0.43
1:D:416:THR:HG23	4:D:589:HOH:O	2.17	0.42
1:G:94:THR:HB	1:G:108:PRO:HB2	2.00	0.42
1:B:49:PRO:HA	1:B:76:LEU:HD13	2.01	0.42
1:E:159:TRP:N	1:E:160:PRO:CD	2.82	0.42
1:E:367:THR:HG21	1:E:463:VAL:HG11	2.00	0.42
1:H:387:ARG:HA	1:H:396:VAL:HG22	2.01	0.42
1:C:129:THR:HG22	1:C:250:THR:OG1	2.19	0.42
1:F:159:TRP:N	1:F:160:PRO:CD	2.82	0.42
1:A:196:LEU:HD11	1:A:205:MET:HE1	2.00	0.42
1:C:455:LEU:HD22	1:D:328:GLY:O	2.19	0.42
1:C:185:ASP:HB2	4:C:694:HOH:O	2.20	0.42
1:C:4:GLU:HG3	1:C:7:LYS:HE2	2.02	0.42
1:G:457:ILE:HD12	1:G:457:ILE:N	2.35	0.42
1:E:196:LEU:HD11	1:E:205:MET:HE1	2.01	0.42
1:G:152:LEU:HD21	1:G:205:MET:CE	2.50	0.42
1:B:457:ILE:CG2	1:B:458:LEU:HD13	2.49	0.42
1:H:230:PRO:HA	1:H:233:TRP:CE2	2.55	0.42
1:A:49:PRO:HA	1:A:76:LEU:HD13	2.01	0.42
1:E:100:TYR:HB3	1:E:105:ILE:HD13	2.02	0.42
1:B:196:LEU:HD11	1:B:205:MET:HE1	2.02	0.42
1:E:100:TYR:CB	1:E:105:ILE:HD13	2.50	0.42
1:A:48:PHE:CB	1:A:61:ILE:HD12	2.50	0.42
1:G:196:LEU:HD11	1:G:205:MET:HE1	2.01	0.42
1:D:152:LEU:HD21	1:D:205:MET:HE1	2.02	0.42
1:H:100:TYR:CB	1:H:105:ILE:HD13	2.49	0.42
1:G:416:THR:HG23	4:G:587:HOH:O	2.20	0.41
1:B:416:THR:HG23	4:B:583:HOH:O	2.20	0.41
1:B:155:PRO:HB3	1:B:344:VAL:HG12	2.02	0.41
1:B:152:LEU:HD12	1:B:206:ASN:O	2.21	0.41
1:H:48:PHE:CG	1:H:61:ILE:HD12	2.56	0.41
1:C:196:LEU:HD11	1:C:205:MET:HE1	2.01	0.41
1:B:457:ILE:HG22	1:B:458:LEU:HD12	2.00	0.41
1:B:100:TYR:CB	1:B:105:ILE:HD13	2.51	0.41
1:B:129:THR:HG22	1:B:250:THR:OG1	2.21	0.41
1:A:93:PHE:CE1	1:B:457:ILE:HG23	2.55	0.41
1:C:159:TRP:N	1:C:160:PRO:CD	2.84	0.41
1:B:152:LEU:HD21	1:B:205:MET:HE1	2.03	0.40
1:E:349:ILE:HG21	1:E:470:ARG:NH2	2.35	0.40
1:H:129:THR:HG22	1:H:250:THR:OG1	2.21	0.40
1:H:166:GLY:O	1:H:186:ASN:ND2	2.53	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:32:THR:CG2	1:H:34:ILE:HD13	2.34	0.40
1:H:152:LEU:HD21	1:H:205:MET:HE1	2.04	0.40
1:H:456:SER:C	1:H:457:ILE:HD12	2.42	0.40
1:F:341:TRP:CE3	2:F:482:MAL:H61	2.56	0.40
1:A:159:TRP:N	1:A:160:PRO:CD	2.84	0.40
1:A:185:ASP:HB2	4:A:691:HOH:O	2.21	0.40
1:F:182:VAL:HG12	1:F:184:VAL:HG13	2.03	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	469/481 (98%)	457 (97%)	10 (2%)	2 (0%)	39	48
1	B	469/481 (98%)	458 (98%)	10 (2%)	1 (0%)	52	64
1	C	469/481 (98%)	458 (98%)	10 (2%)	1 (0%)	52	64
1	D	467/481 (97%)	457 (98%)	9 (2%)	1 (0%)	52	64
1	E	470/481 (98%)	463 (98%)	6 (1%)	1 (0%)	52	64
1	F	470/481 (98%)	460 (98%)	9 (2%)	1 (0%)	52	64
1	G	470/481 (98%)	460 (98%)	9 (2%)	1 (0%)	52	64
1	H	467/481 (97%)	457 (98%)	9 (2%)	1 (0%)	52	64
All	All	3751/3848 (98%)	3670 (98%)	72 (2%)	9 (0%)	52	64

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	378	GLU
1	F	378	GLU
1	A	378	GLU

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Mol	Chain	Res	Type
1	D	378	GLU
1	E	378	GLU
1	G	378	GLU
1	A	4	GLU
1	B	378	GLU
1	H	378	GLU

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	383/393 (98%)	383 (100%)	0	100	100
1	B	383/393 (98%)	383 (100%)	0	100	100
1	C	383/393 (98%)	383 (100%)	0	100	100
1	D	381/393 (97%)	381 (100%)	0	100	100
1	E	384/393 (98%)	384 (100%)	0	100	100
1	F	384/393 (98%)	384 (100%)	0	100	100
1	G	384/393 (98%)	384 (100%)	0	100	100
1	H	381/393 (97%)	381 (100%)	0	100	100
All	All	3063/3144 (97%)	3063 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

Of 32 ligands modelled in this entry, 24 are monoatomic - leaving 8 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$
2	MAL	A	482	-	24,24,24	0.77	1 (4%)	35,35,35	0.86	1 (2%)
2	MAL	B	482	-	24,24,24	0.90	0	35,35,35	1.00	3 (8%)
2	MAL	C	482	-	24,24,24	0.73	0	35,35,35	0.90	1 (2%)
2	MAL	D	482	-	24,24,24	0.87	0	35,35,35	0.90	2 (5%)
2	MAL	E	482	-	24,24,24	0.72	0	35,35,35	0.86	2 (5%)
2	MAL	F	482	-	24,24,24	0.73	0	35,35,35	0.78	1 (2%)
2	MAL	G	482	-	24,24,24	0.85	0	35,35,35	0.88	2 (5%)
2	MAL	H	482	-	24,24,24	0.91	0	35,35,35	0.93	2 (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	MAL	A	482	-	-	0/8/48/48	0/2/2/2
2	MAL	B	482	-	-	0/8/48/48	0/2/2/2
2	MAL	C	482	-	-	0/8/48/48	0/2/2/2
2	MAL	D	482	-	-	0/8/48/48	0/2/2/2
2	MAL	E	482	-	-	0/8/48/48	0/2/2/2
2	MAL	F	482	-	-	0/8/48/48	0/2/2/2
2	MAL	G	482	-	-	0/8/48/48	0/2/2/2
2	MAL	H	482	-	-	0/8/48/48	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	482	MAL	O5'-C1'	2.14	1.47	1.43

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	482	MAL	C6-C5-C4	-2.52	106.81	113.02
2	B	482	MAL	C6-C5-C4	-2.48	106.90	113.02
2	H	482	MAL	C6-C5-C4	-2.34	107.25	113.02
2	B	482	MAL	C1-O1-C4'	-2.33	111.93	118.01
2	G	482	MAL	C6-C5-C4	-2.30	107.35	113.02
2	D	482	MAL	C1-O1-C4'	-2.13	112.43	118.01
2	E	482	MAL	C6-C5-C4	-2.12	107.77	113.02
2	G	482	MAL	C1-O1-C4'	-2.04	112.67	118.01
2	F	482	MAL	C1-O1-C4'	-2.02	112.73	118.01
2	A	482	MAL	O1-C1-C2	2.00	112.98	108.10
2	H	482	MAL	O5'-C1'-C2'	2.08	113.12	109.80
2	B	482	MAL	O5'-C1'-C2'	2.15	113.23	109.80
2	E	482	MAL	O5'-C1'-C2'	2.43	113.67	109.80
2	C	482	MAL	O5'-C1'-C2'	2.52	113.82	109.80

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	482	MAL	1	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

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	471/481 (97%)	0.20	21 (4%) 37 46	28, 52, 97, 152	0
1	B	471/481 (97%)	0.41	29 (6%) 24 32	29, 64, 111, 134	0
1	C	471/481 (97%)	0.19	18 (3%) 44 53	28, 52, 98, 161	0
1	D	469/481 (97%)	0.44	25 (5%) 30 39	30, 66, 117, 143	0
1	E	472/481 (98%)	0.24	17 (3%) 46 55	28, 52, 97, 162	0
1	F	472/481 (98%)	0.25	21 (4%) 38 47	28, 53, 96, 166	0
1	G	472/481 (98%)	0.39	22 (4%) 35 44	30, 65, 114, 139	0
1	H	469/481 (97%)	0.39	25 (5%) 30 39	30, 66, 113, 133	0
All	All	3767/3848 (97%)	0.31	178 (4%) 35 44	28, 58, 109, 166	0

All (178) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	453	SER	7.9
1	F	453	SER	7.0
1	G	3	THR	6.4
1	F	454	GLY	5.9
1	F	3	THR	5.9
1	F	455	LEU	5.7
1	E	455	LEU	5.4
1	E	458	LEU	5.0
1	F	458	LEU	4.9
1	D	455	LEU	4.9
1	F	456	SER	4.8
1	E	3	THR	4.8
1	C	454	GLY	4.6
1	E	456	SER	4.4
1	E	454	GLY	4.3
1	C	455	LEU	4.3

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Mol	Chain	Res	Type	RSRZ
1	F	474	LEU	4.3
1	B	458	LEU	4.2
1	F	452	VAL	4.1
1	G	4	GLU	4.1
1	A	459	ARG	4.1
1	A	454	GLY	4.1
1	D	281	LEU	4.0
1	D	6	GLY	3.9
1	A	457	ILE	3.9
1	D	34	ILE	3.9
1	G	123	LEU	3.8
1	H	458	LEU	3.7
1	C	221	GLY	3.7
1	G	236	ILE	3.6
1	D	456	SER	3.6
1	H	36	VAL	3.6
1	D	280	PHE	3.6
1	D	236	ILE	3.5
1	G	285	LEU	3.5
1	B	3	THR	3.5
1	A	140	LEU	3.5
1	A	455	LEU	3.5
1	D	458	LEU	3.5
1	A	452	VAL	3.4
1	C	457	ILE	3.4
1	D	28	PHE	3.4
1	H	455	LEU	3.4
1	B	240	LYS	3.4
1	D	285	LEU	3.4
1	E	236	ILE	3.3
1	B	24	VAL	3.3
1	C	456	SER	3.3
1	D	453	SER	3.3
1	F	236	ILE	3.2
1	G	239	SER	3.2
1	H	240	LYS	3.1
1	G	243	TYR	3.1
1	H	35	LYS	3.1
1	A	4	GLU	3.1
1	H	236	ILE	3.0
1	G	458	LEU	3.0
1	B	36	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	458	LEU	3.0
1	H	20	GLY	2.9
1	C	452	VAL	2.9
1	F	237	ASP	2.9
1	A	3	THR	2.9
1	D	457	ILE	2.9
1	D	454	GLY	2.9
1	A	145	LYS	2.9
1	E	118	TYR	2.9
1	B	123	LEU	2.9
1	B	455	LEU	2.9
1	F	221	GLY	2.9
1	G	455	LEU	2.9
1	H	285	LEU	2.9
1	B	243	TYR	2.8
1	E	224	ALA	2.8
1	C	3	THR	2.8
1	B	34	ILE	2.8
1	B	454	GLY	2.8
1	E	221	GLY	2.8
1	G	281	LEU	2.8
1	B	453	SER	2.8
1	C	459	ARG	2.8
1	F	242	ASN	2.8
1	H	457	ILE	2.7
1	D	240	LYS	2.7
1	D	93	PHE	2.7
1	E	4	GLU	2.7
1	B	456	SER	2.7
1	H	239	SER	2.7
1	C	220	LYS	2.7
1	B	236	ILE	2.7
1	H	28	PHE	2.7
1	H	242	ASN	2.7
1	A	219	ASN	2.6
1	C	219	ASN	2.6
1	D	8	LEU	2.6
1	A	236	ILE	2.6
1	B	457	ILE	2.6
1	E	219	ASN	2.6
1	G	118	TYR	2.6
1	B	38	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	242	ASN	2.6
1	E	474	LEU	2.6
1	H	453	SER	2.5
1	E	452	VAL	2.5
1	F	4	GLU	2.5
1	H	243	TYR	2.5
1	D	24	VAL	2.5
1	H	30	LYS	2.5
1	G	6	GLY	2.5
1	E	240	LYS	2.4
1	D	243	TYR	2.4
1	A	240	LYS	2.4
1	A	242	ASN	2.4
1	C	458	LEU	2.4
1	F	243	TYR	2.4
1	F	220	LYS	2.4
1	C	451	PRO	2.4
1	F	459	ARG	2.4
1	H	88	ASP	2.4
1	G	28	PHE	2.3
1	A	221	GLY	2.3
1	B	20	GLY	2.3
1	H	122	LEU	2.3
1	C	240	LYS	2.3
1	G	326	GLN	2.3
1	B	93	PHE	2.3
1	H	34	ILE	2.3
1	B	4	GLU	2.3
1	B	293	ALA	2.3
1	B	294	VAL	2.3
1	G	122	LEU	2.3
1	C	224	ALA	2.2
1	D	294	VAL	2.2
1	F	219	ASN	2.2
1	B	450	ARG	2.2
1	E	220	LYS	2.2
1	H	454	GLY	2.2
1	H	310	GLU	2.2
1	D	276	LEU	2.2
1	A	226	THR	2.2
1	G	174	ASN	2.2
1	A	451	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	239	SER	2.2
1	C	239	SER	2.2
1	F	133	ILE	2.2
1	F	457	ILE	2.2
1	B	118	TYR	2.2
1	B	28	PHE	2.2
1	H	456	SER	2.2
1	G	474	LEU	2.2
1	D	450	ARG	2.1
1	H	294	VAL	2.1
1	C	218	PHE	2.1
1	H	123	LEU	2.1
1	A	224	ALA	2.1
1	D	111	VAL	2.1
1	B	308	TYR	2.1
1	G	240	LYS	2.1
1	A	453	SER	2.1
1	D	88	ASP	2.1
1	H	38	VAL	2.1
1	G	238	THR	2.1
1	B	30	LYS	2.1
1	E	242	ASN	2.1
1	G	242	ASN	2.1
1	B	241	VAL	2.1
1	A	118	TYR	2.1
1	B	21	LEU	2.1
1	A	238	THR	2.0
1	G	322	MET	2.0
1	D	118	TYR	2.0
1	F	240	LYS	2.0
1	F	145	LYS	2.0
1	G	93	PHE	2.0
1	C	216	ALA	2.0
1	H	118	TYR	2.0
1	D	300	LEU	2.0
1	C	118	TYR	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(Å ²)	Q<0.9
2	MAL	E	482	23/23	0.96	0.14	1.72	28,34,39,41	0
2	MAL	C	482	23/23	0.97	0.13	0.88	29,35,41,41	0
2	MAL	F	482	23/23	0.96	0.14	0.76	30,36,42,42	0
3	CA	E	484	1/1	0.83	0.22	0.54	58,58,58,58	0
2	MAL	A	482	23/23	0.97	0.13	0.50	27,33,39,40	0
2	MAL	B	482	23/23	0.95	0.13	0.42	39,44,50,52	0
2	MAL	D	482	23/23	0.96	0.14	0.11	40,46,55,57	0
2	MAL	G	482	23/23	0.97	0.12	0.06	39,43,50,50	0
2	MAL	H	482	23/23	0.95	0.13	-0.04	39,43,51,52	0
3	CA	C	484	1/1	0.87	0.15	-0.05	57,57,57,57	0
3	CA	A	484	1/1	0.84	0.14	-0.45	57,57,57,57	0
3	CA	F	484	1/1	0.94	0.12	-0.70	57,57,57,57	0
3	CA	E	483	1/1	0.92	0.07	-2.33	67,67,67,67	0
3	CA	G	484	1/1	0.96	0.05	-3.26	54,54,54,54	0
3	CA	D	484	1/1	0.97	0.04	-3.67	55,55,55,55	0
3	CA	C	483	1/1	0.92	0.04	-3.73	64,64,64,64	0
3	CA	B	484	1/1	0.94	0.05	-4.71	54,54,54,54	0
3	CA	H	484	1/1	0.95	0.04	-5.17	54,54,54,54	0
3	CA	G	485	1/1	0.82	0.16	-	59,59,59,59	0
3	CA	H	483	1/1	0.93	0.07	-	89,89,89,89	0
3	CA	A	485	1/1	0.86	0.10	-	61,61,61,61	0
3	CA	D	485	1/1	0.90	0.18	-	58,58,58,58	0
3	CA	B	483	1/1	0.83	0.07	-	93,93,93,93	0
3	CA	E	485	1/1	0.95	0.14	-	61,61,61,61	0
3	CA	F	485	1/1	0.93	0.10	-	60,60,60,60	0
3	CA	G	483	1/1	0.86	0.05	-	85,85,85,85	0
3	CA	A	483	1/1	0.78	0.06	-	70,70,70,70	0
3	CA	H	485	1/1	0.97	0.11	-	59,59,59,59	0
3	CA	B	485	1/1	0.96	0.21	-	59,59,59,59	0
3	CA	D	483	1/1	0.89	0.16	-	95,95,95,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	CA	F	483	1/1	0.83	0.07	-	71,71,71,71	0
3	CA	C	485	1/1	0.87	0.13	-	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.