



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 29, 2017 – 03:24 PM EDT

PDB ID : 3FPV
Title : Crystal Structure of HbpS
Authors : Ortiz de Orue Lucana, D.; Bogel, G.; Zou, P.; Groves, M.R.
Deposited on : unknown
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

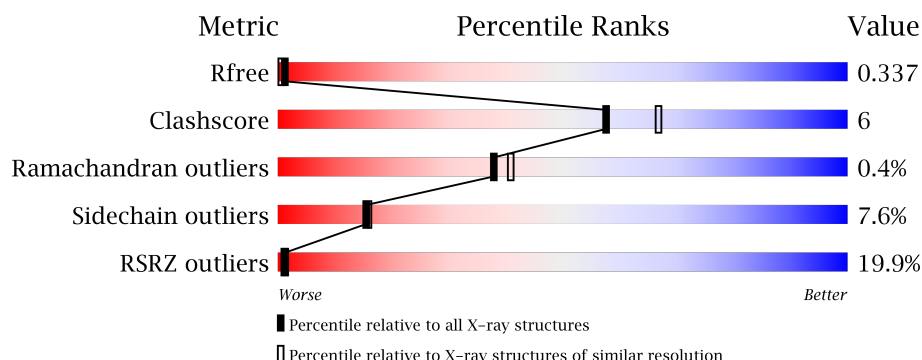
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



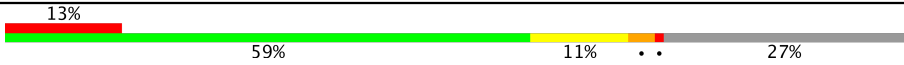


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	192	<div> <div>11%</div> <div> <div>60%</div> <div>11%</div> <div>27%</div> </div> </div>
1	B	192	<div> <div>18%</div> <div> <div>64%</div> <div>7%</div> <div>27%</div> </div> </div>
1	C	192	<div> <div>14%</div> <div> <div>61%</div> <div>10%</div> <div>27%</div> </div> </div>
1	D	192	<div> <div>14%</div> <div> <div>60%</div> <div>11%</div> <div>27%</div> </div> </div>
1	E	192	<div> <div>21%</div> <div> <div>61%</div> <div>10%</div> <div>27%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	192	
1	G	192	
1	H	192	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8497 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 Extracellular haem-binding protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	0	0	0
			997	614	185	198			
1	B	141	Total	C	N	O	0	0	0
			997	614	185	198			
1	C	141	Total	C	N	O	0	0	0
			997	614	185	198			
1	D	141	Total	C	N	O	0	0	0
			997	614	185	198			
1	E	141	Total	C	N	O	0	0	0
			997	614	185	198			
1	F	141	Total	C	N	O	0	0	0
			997	614	185	198			
1	G	141	Total	C	N	O	0	0	0
			997	614	185	198			
1	H	141	Total	C	N	O	0	0	0
			997	614	185	198			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-35	GLY	-	EXPRESSION TAG	UNP Q9RIM2
A	-34	ALA	-	EXPRESSION TAG	UNP Q9RIM2
A	-33	GLY	-	EXPRESSION TAG	UNP Q9RIM2
A	-32	ALA	-	EXPRESSION TAG	UNP Q9RIM2
B	-35	GLY	-	EXPRESSION TAG	UNP Q9RIM2
B	-34	ALA	-	EXPRESSION TAG	UNP Q9RIM2
B	-33	GLY	-	EXPRESSION TAG	UNP Q9RIM2
B	-32	ALA	-	EXPRESSION TAG	UNP Q9RIM2
C	-35	GLY	-	EXPRESSION TAG	UNP Q9RIM2
C	-34	ALA	-	EXPRESSION TAG	UNP Q9RIM2
C	-33	GLY	-	EXPRESSION TAG	UNP Q9RIM2
C	-32	ALA	-	EXPRESSION TAG	UNP Q9RIM2
D	-35	GLY	-	EXPRESSION TAG	UNP Q9RIM2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-34	ALA	-	EXPRESSION TAG	UNP Q9RIM2
D	-33	GLY	-	EXPRESSION TAG	UNP Q9RIM2
D	-32	ALA	-	EXPRESSION TAG	UNP Q9RIM2
E	-35	GLY	-	EXPRESSION TAG	UNP Q9RIM2
E	-34	ALA	-	EXPRESSION TAG	UNP Q9RIM2
E	-33	GLY	-	EXPRESSION TAG	UNP Q9RIM2
E	-32	ALA	-	EXPRESSION TAG	UNP Q9RIM2
F	-35	GLY	-	EXPRESSION TAG	UNP Q9RIM2
F	-34	ALA	-	EXPRESSION TAG	UNP Q9RIM2
F	-33	GLY	-	EXPRESSION TAG	UNP Q9RIM2
F	-32	ALA	-	EXPRESSION TAG	UNP Q9RIM2
G	-35	GLY	-	EXPRESSION TAG	UNP Q9RIM2
G	-34	ALA	-	EXPRESSION TAG	UNP Q9RIM2
G	-33	GLY	-	EXPRESSION TAG	UNP Q9RIM2
G	-32	ALA	-	EXPRESSION TAG	UNP Q9RIM2
H	-35	GLY	-	EXPRESSION TAG	UNP Q9RIM2
H	-34	ALA	-	EXPRESSION TAG	UNP Q9RIM2
H	-33	GLY	-	EXPRESSION TAG	UNP Q9RIM2
H	-32	ALA	-	EXPRESSION TAG	UNP Q9RIM2

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total Fe 1 1	0	0
2	E	2	Total Fe 2 2	0	0
2	H	1	Total Fe 1 1	0	0
2	B	1	Total Fe 1 1	0	0
2	C	1	Total Fe 1 1	0	0
2	A	1	Total Fe 1 1	0	0
2	F	1	Total Fe 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	81	Total O 81 81	0	0

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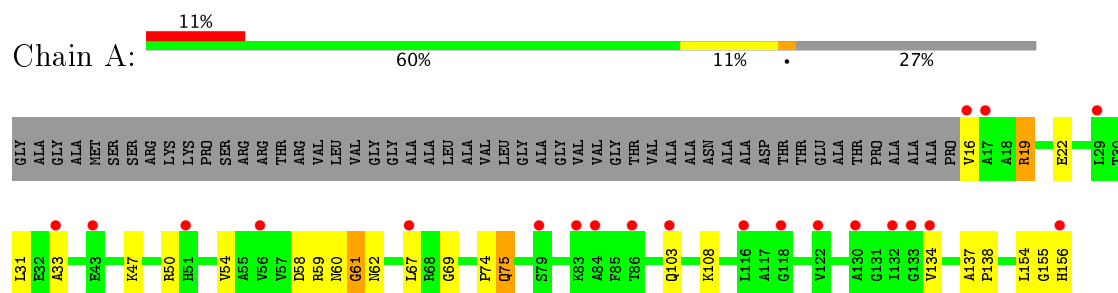
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	101	Total	O	0	0
			101	101		
3	C	57	Total	O	0	0
			57	57		
3	D	50	Total	O	0	0
			50	50		
3	E	35	Total	O	0	0
			35	35		
3	F	84	Total	O	0	0
			84	84		
3	G	46	Total	O	0	0
			46	46		
3	H	59	Total	O	0	0
			59	59		

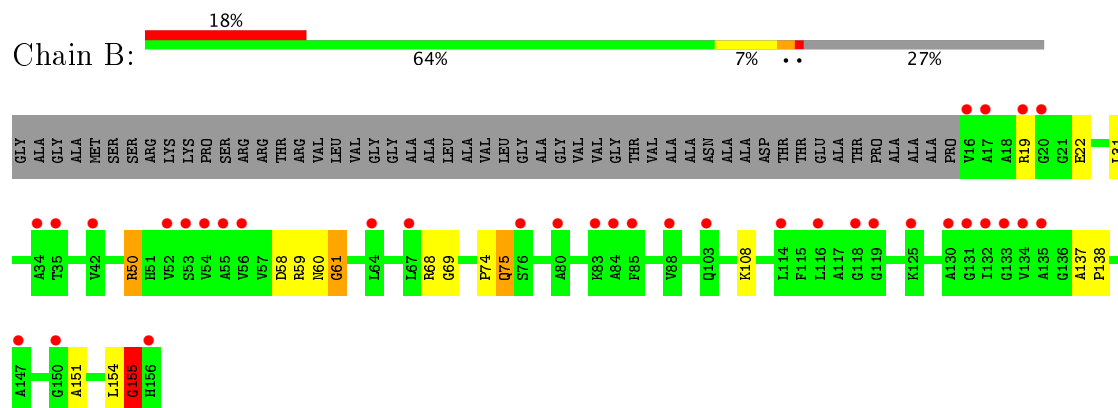
3 Residue-property plots [i](#)

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

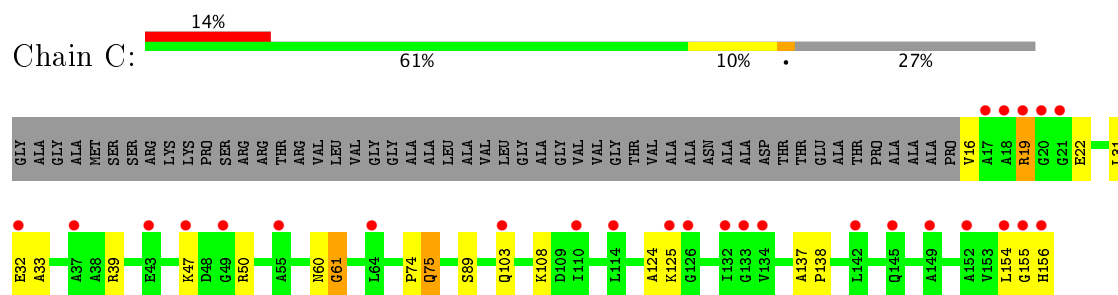
- Molecule 1: Extracellular haem-binding protein



- Molecule 1: Extracellular haem-binding protein

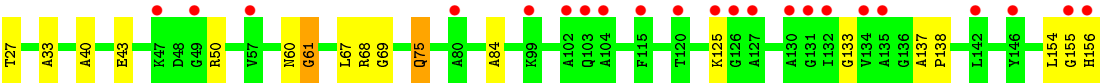


- Molecule 1: Extracellular haem-binding protein



- Molecule 1: Extracellular haem-binding protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	152.52Å 152.52Å 152.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.86 – 2.20 19.86 – 2.21	Depositor EDS
% Data completeness (in resolution range)	99.1 (19.86-2.20) 99.1 (19.86-2.21)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.21Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.201 , 0.264 0.295 , 0.337	Depositor DCC
R_{free} test set	2914 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.033 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8497	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.84% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.21	0/1012	0.36	0/1376
1	B	0.21	0/1012	0.36	0/1376
1	C	0.20	0/1012	0.35	0/1376
1	D	0.20	0/1012	0.36	0/1376
1	E	0.21	0/1012	0.36	0/1376
1	F	0.21	0/1012	0.36	0/1376
1	G	0.20	0/1012	0.35	0/1376
1	H	0.20	0/1012	0.35	0/1376
All	All	0.21	0/8096	0.36	0/11008

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	4
1	B	0	5
1	C	0	6
1	D	0	8
1	E	0	10
1	F	0	8
1	G	0	4
1	H	0	6
All	All	0	51

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (51) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	154	LEU	Peptide
1	A	19	ARG	Peptide
1	A	60	ASN	Peptide
1	A	61	GLY	Peptide
1	B	154	LEU	Peptide
1	B	155	GLY	Peptide
1	B	19	ARG	Peptide
1	B	60	ASN	Peptide
1	B	61	GLY	Peptide
1	C	124	ALA	Peptide
1	C	154	LEU	Peptide
1	C	155	GLY	Peptide
1	C	19	ARG	Peptide
1	C	60	ASN	Peptide
1	C	61	GLY	Peptide
1	D	124	ALA	Peptide
1	D	135	ALA	Peptide
1	D	136	GLY	Peptide
1	D	154	LEU	Peptide
1	D	155	GLY	Peptide
1	D	19	ARG	Peptide
1	D	60	ASN	Peptide
1	D	61	GLY	Peptide
1	E	124	ALA	Peptide
1	E	135	ALA	Peptide
1	E	154	LEU	Peptide
1	E	16	VAL	Peptide
1	E	18	ALA	Peptide
1	E	19	ARG	Peptide
1	E	22	GLU	Peptide
1	E	50	ARG	Peptide
1	E	60	ASN	Peptide
1	E	61	GLY	Peptide
1	F	135	ALA	Peptide
1	F	136	GLY	Peptide
1	F	154	LEU	Peptide
1	F	155	GLY	Peptide
1	F	16	VAL	Peptide
1	F	19	ARG	Peptide
1	F	60	ASN	Peptide
1	F	61	GLY	Peptide
1	G	155	GLY	Peptide
1	G	19	ARG	Peptide

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Mol	Chain	Res	Type	Group
1	G	60	ASN	Peptide
1	G	61	GLY	Peptide
1	H	154	LEU	Peptide
1	H	155	GLY	Peptide
1	H	16	VAL	Peptide
1	H	19	ARG	Peptide
1	H	60	ASN	Peptide
1	H	61	GLY	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	997	0	996	20	0
1	B	997	0	996	17	0
1	C	997	0	996	11	0
1	D	997	0	996	14	0
1	E	997	0	996	15	0
1	F	997	0	996	22	0
1	G	997	0	996	12	0
1	H	997	0	996	16	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	2	0	0	0	0
2	F	1	0	0	0	0
2	H	1	0	0	0	0
3	A	81	0	0	0	0
3	B	101	0	0	1	0
3	C	57	0	0	0	0
3	D	50	0	0	0	0
3	E	35	0	0	0	0
3	F	84	0	0	2	0
3	G	46	0	0	0	0
3	H	59	0	0	1	0
All	All	8497	0	7968	90	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 6.

All (90) 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:22:GLU:HB2	1:H:33:ALA:HB2	1.56	0.85
1:A:155:GLY:HA2	1:F:22:GLU:OE1	1.80	0.82
1:A:33:ALA:HB2	1:F:22:GLU:HB3	1.65	0.77
1:C:33:ALA:HB2	1:H:22:GLU:HB2	1.69	0.75
1:A:155:GLY:N	1:A:156:HIS:HA	2.02	0.74
1:D:22:GLU:HB2	1:E:33:ALA:HB2	1.69	0.74
1:B:155:GLY:HA2	1:G:22:GLU:OE1	1.91	0.70
1:A:22:GLU:HB3	1:F:33:ALA:HB2	1.74	0.68
1:H:19:ARG:HG3	1:H:20:GLY:H	1.60	0.66
1:H:19:ARG:HG3	1:H:20:GLY:N	2.13	0.63
1:D:33:ALA:HB2	1:E:22:GLU:HG3	1.79	0.62
1:B:137:ALA:HB1	1:B:138:PRO:HD2	1.81	0.62
1:A:33:ALA:CB	1:F:22:GLU:HB3	2.30	0.60
1:B:61:GLY:H	1:D:74:PRO:HB3	1.65	0.60
1:C:31:LEU:HD11	1:E:67:LEU:HB3	1.83	0.60
1:E:61:GLY:H	1:G:74:PRO:HB3	1.67	0.59
1:A:75:GLN:NE2	1:A:75:GLN:H	2.01	0.59
1:D:75:GLN:NE2	1:D:75:GLN:H	1.99	0.58
1:D:33:ALA:CB	1:E:22:GLU:HG3	2.32	0.58
1:B:75:GLN:H	1:B:75:GLN:HE21	1.53	0.57
1:B:75:GLN:H	1:B:75:GLN:NE2	2.02	0.57
1:B:58:ASP:O	1:B:61:GLY:HA2	2.05	0.57
1:A:75:GLN:HE21	1:A:75:GLN:H	1.53	0.55
1:A:22:GLU:OE1	1:F:155:GLY:HA2	2.08	0.54
1:F:31:LEU:HD11	1:H:67:LEU:HB3	1.89	0.54
1:F:75:GLN:NE2	1:F:75:GLN:H	2.06	0.54
1:F:17:ALA:HB1	3:F:1210:HOH:O	2.07	0.53
1:A:22:GLU:HB3	1:F:33:ALA:CB	2.38	0.53
1:H:16:VAL:N	3:H:1176:HOH:O	2.42	0.53
1:E:48:ASP:N	1:E:48:ASP:OD1	2.42	0.53
1:A:69:GLY:HA2	1:B:31:LEU:HD22	1.91	0.53
1:A:74:PRO:HB3	1:G:61:GLY:H	1.74	0.52
1:H:137:ALA:HB1	1:H:138:PRO:HD2	1.90	0.52
1:B:151:ALA:HB1	1:G:16:VAL:HG21	1.90	0.52
1:C:75:GLN:H	1:C:75:GLN:NE2	2.07	0.52
1:D:75:GLN:HE21	1:D:75:GLN:H	1.57	0.51
1:B:137:ALA:HB1	1:B:138:PRO:CD	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:74:PRO:HB3	1:H:61:GLY:H	1.75	0.51
1:G:58:ASP:O	1:G:61:GLY:HA2	2.11	0.51
1:H:75:GLN:H	1:H:75:GLN:HE21	1.58	0.51
1:C:74:PRO:HB3	1:D:61:GLY:H	1.76	0.51
1:H:75:GLN:NE2	1:H:75:GLN:H	2.08	0.50
1:E:75:GLN:HE21	1:E:75:GLN:H	1.59	0.50
1:F:75:GLN:HE21	1:F:75:GLN:H	1.58	0.50
1:G:75:GLN:H	1:G:75:GLN:NE2	2.08	0.50
1:D:54:VAL:HG22	1:D:134:VAL:HG22	1.93	0.50
1:F:31:LEU:HD22	1:H:69:GLY:HA2	1.94	0.49
1:B:59:ARG:O	1:B:59:ARG:HG3	2.12	0.49
1:D:39:ARG:NH2	1:G:32:GLU:OE1	2.44	0.49
1:G:137:ALA:HB1	1:G:138:PRO:HD2	1.94	0.49
1:A:62:ASN:HB3	1:B:68:ARG:CZ	2.42	0.49
1:C:137:ALA:HB1	1:C:138:PRO:HD2	1.95	0.49
1:A:137:ALA:HB1	1:A:138:PRO:HD2	1.95	0.49
1:E:155:GLY:HA3	1:E:156:HIS:HA	1.53	0.48
1:F:58:ASP:O	1:F:61:GLY:CA	2.61	0.48
1:B:58:ASP:O	1:B:61:GLY:CA	2.61	0.48
1:F:54:VAL:HG22	1:F:134:VAL:HG22	1.94	0.48
1:C:75:GLN:NE2	1:D:89:SER:HB3	2.29	0.47
1:A:58:ASP:O	1:A:61:GLY:HA2	2.15	0.47
1:C:39:ARG:HH11	1:E:39:ARG:HH11	1.63	0.47
1:F:62:ASN:HB3	1:H:68:ARG:CZ	2.45	0.47
1:D:33:ALA:HB2	1:E:22:GLU:CG	2.44	0.46
1:C:33:ALA:CB	1:H:22:GLU:HB2	2.44	0.46
1:H:137:ALA:HB1	1:H:138:PRO:CD	2.45	0.46
1:F:58:ASP:O	1:F:61:GLY:HA2	2.16	0.45
1:A:54:VAL:HG22	1:A:134:VAL:HG22	1.98	0.45
1:A:58:ASP:O	1:A:61:GLY:CA	2.64	0.45
1:E:152:ALA:O	1:E:156:HIS:NE2	2.50	0.45
1:A:31:LEU:HD22	1:B:69:GLY:HA2	1.98	0.45
1:G:125:LYS:HD2	1:G:125:LYS:HA	1.79	0.45
1:A:67:LEU:HB3	1:B:31:LEU:HD11	1.99	0.44
1:B:50:ARG:HD3	3:B:717:HOH:O	2.16	0.44
1:D:33:ALA:HB2	1:E:22:GLU:HB2	2.00	0.44
1:A:137:ALA:HB1	1:A:138:PRO:CD	2.48	0.44
1:B:74:PRO:HB3	1:F:61:GLY:H	1.81	0.44
1:C:61:GLY:H	1:F:74:PRO:HB3	1.81	0.44
1:B:22:GLU:OE1	1:G:155:GLY:HA2	2.18	0.44
1:E:137:ALA:HB1	1:E:138:PRO:CD	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:50:ARG:HG2	1:E:138:PRO:HD3	2.00	0.43
1:G:58:ASP:HB2	1:G:64:LEU:HD21	2.01	0.43
1:F:155:GLY:C	1:F:156:HIS:CG	2.92	0.43
1:C:89:SER:HB3	1:F:75:GLN:NE2	2.34	0.42
1:A:59:ARG:O	1:A:59:ARG:HG3	2.18	0.42
1:F:105:PRO:O	1:F:108:LYS:HG2	2.20	0.42
1:F:25:GLN:HG2	3:F:1209:HOH:O	2.19	0.42
1:D:35:THR:O	1:D:39:ARG:HG3	2.20	0.42
1:F:150:GLY:O	1:F:153:VAL:HG22	2.19	0.42
1:H:84:ALA:HB2	1:H:133:GLY:N	2.36	0.41
1:H:40:ALA:HA	1:H:43:GLU:HB3	2.03	0.41
1:D:68:ARG:CZ	1:G:62:ASN:HB3	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	139/192 (72%)	136 (98%)	3 (2%)	0	100	100
1	B	139/192 (72%)	137 (99%)	1 (1%)	1 (1%)	25	24
1	C	139/192 (72%)	134 (96%)	4 (3%)	1 (1%)	25	24
1	D	139/192 (72%)	132 (95%)	6 (4%)	1 (1%)	25	24
1	E	139/192 (72%)	134 (96%)	4 (3%)	1 (1%)	25	24
1	F	139/192 (72%)	136 (98%)	2 (1%)	1 (1%)	25	24
1	G	139/192 (72%)	138 (99%)	1 (1%)	0	100	100
1	H	139/192 (72%)	135 (97%)	4 (3%)	0	100	100
All	All	1112/1536 (72%)	1082 (97%)	25 (2%)	5 (0%)	38	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	125	LYS
1	D	125	LYS
1	E	125	LYS
1	F	155	GLY
1	B	155	GLY

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	95/125 (76%)	88 (93%)	7 (7%)	16	17
1	B	95/125 (76%)	92 (97%)	3 (3%)	44	56
1	C	95/125 (76%)	86 (90%)	9 (10%)	10	9
1	D	95/125 (76%)	86 (90%)	9 (10%)	10	9
1	E	95/125 (76%)	89 (94%)	6 (6%)	21	23
1	F	95/125 (76%)	87 (92%)	8 (8%)	13	12
1	G	95/125 (76%)	86 (90%)	9 (10%)	10	9
1	H	95/125 (76%)	88 (93%)	7 (7%)	16	17
All	All	760/1000 (76%)	702 (92%)	58 (8%)	15	16

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	VAL
1	A	19	ARG
1	A	47	LYS
1	A	50	ARG
1	A	75	GLN
1	A	103	GLN
1	A	108	LYS
1	B	50	ARG
1	B	75	GLN
1	B	108	LYS
1	C	16	VAL

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Mol	Chain	Res	Type
1	C	19	ARG
1	C	32	GLU
1	C	47	LYS
1	C	50	ARG
1	C	75	GLN
1	C	103	GLN
1	C	108	LYS
1	C	156	HIS
1	D	19	ARG
1	D	22	GLU
1	D	32	GLU
1	D	47	LYS
1	D	50	ARG
1	D	75	GLN
1	D	108	LYS
1	D	132	ILE
1	D	141	ASP
1	E	22	GLU
1	E	48	ASP
1	E	50	ARG
1	E	75	GLN
1	E	96	GLU
1	E	108	LYS
1	F	16	VAL
1	F	19	ARG
1	F	47	LYS
1	F	50	ARG
1	F	75	GLN
1	F	103	GLN
1	F	120	THR
1	F	156	HIS
1	G	16	VAL
1	G	43	GLU
1	G	47	LYS
1	G	50	ARG
1	G	75	GLN
1	G	79	SER
1	G	103	GLN
1	G	154	LEU
1	G	156	HIS
1	H	16	VAL
1	H	22	GLU

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Mol	Chain	Res	Type
1	H	27	THR
1	H	50	ARG
1	H	75	GLN
1	H	125	LYS
1	H	156	HIS

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

Mol	Chain	Res	Type
1	A	25	GLN
1	A	75	GLN
1	A	156	HIS
1	B	75	GLN
1	C	25	GLN
1	C	75	GLN
1	D	75	GLN
1	D	145	GLN
1	E	25	GLN
1	E	75	GLN
1	F	25	GLN
1	F	75	GLN
1	G	25	GLN
1	G	75	GLN
1	H	75	GLN
1	H	156	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 8 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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	141/192 (73%)	0.95	21 (14%) 3 2	64, 67, 77, 80	0
1	B	141/192 (73%)	1.21	35 (24%) 1 1	63, 68, 77, 82	0
1	C	141/192 (73%)	1.19	27 (19%) 1 1	64, 67, 76, 79	0
1	D	141/192 (73%)	1.14	27 (19%) 1 1	64, 67, 76, 78	0
1	E	141/192 (73%)	1.73	41 (29%) 1 0	64, 70, 74, 77	0
1	F	141/192 (73%)	1.02	24 (17%) 2 2	63, 68, 77, 82	0
1	G	141/192 (73%)	0.95	22 (15%) 2 2	64, 67, 76, 82	0
1	H	141/192 (73%)	1.25	28 (19%) 1 1	62, 66, 75, 79	0
All	All	1128/1536 (73%)	1.18	225 (19%) 1 1	62, 68, 76, 82	0

All (225) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	152	ALA	7.7
1	H	156	HIS	7.4
1	E	50	ARG	7.2
1	D	19	ARG	7.0
1	E	20	GLY	6.4
1	E	156	HIS	6.3
1	E	19	ARG	6.1
1	D	20	GLY	6.0
1	E	126	GLY	5.9
1	F	156	HIS	5.8
1	D	152	ALA	5.8
1	B	134	VAL	5.4
1	C	126	GLY	5.3
1	F	19	ARG	5.0
1	H	47	LYS	4.9
1	D	126	GLY	4.8

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Mol	Chain	Res	Type	RSRZ
1	C	125	LYS	4.7
1	A	156	HIS	4.7
1	A	134	VAL	4.7
1	D	156	HIS	4.6
1	H	19	ARG	4.6
1	C	19	ARG	4.6
1	B	156	HIS	4.5
1	G	50	ARG	4.5
1	E	51	HIS	4.5
1	E	127	ALA	4.5
1	G	156	HIS	4.4
1	D	17	ALA	4.4
1	B	132	ILE	4.4
1	E	155	GLY	4.4
1	E	145	GLN	4.3
1	B	133	GLY	4.3
1	B	53	SER	4.3
1	F	155	GLY	4.3
1	B	135	ALA	4.2
1	G	19	ARG	4.1
1	E	124	ALA	4.1
1	F	16	VAL	3.9
1	E	125	LYS	3.8
1	H	155	GLY	3.8
1	E	47	LYS	3.8
1	A	132	ILE	3.8
1	D	155	GLY	3.8
1	H	21	GLY	3.8
1	E	23	LEU	3.7
1	B	116	LEU	3.7
1	C	49	GLY	3.7
1	E	21	GLY	3.7
1	D	21	GLY	3.6
1	C	32	GLU	3.6
1	C	155	GLY	3.6
1	C	43	GLU	3.6
1	H	80	ALA	3.6
1	D	25	GLN	3.6
1	E	134	VAL	3.5
1	B	55	ALA	3.5
1	B	54	VAL	3.5
1	G	155	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	125	LYS	3.5
1	D	114	LEU	3.4
1	E	146	TYR	3.4
1	C	18	ALA	3.4
1	B	84	ALA	3.4
1	F	114	LEU	3.4
1	B	131	GLY	3.4
1	E	46	GLU	3.3
1	E	25	GLN	3.3
1	G	51	HIS	3.3
1	E	141	ASP	3.3
1	E	103	GLN	3.3
1	E	148	ARG	3.3
1	B	125	LYS	3.3
1	B	19	ARG	3.3
1	C	156	HIS	3.2
1	H	24	THR	3.2
1	D	116	LEU	3.2
1	D	134	VAL	3.1
1	H	102	ALA	3.1
1	A	84	ALA	3.1
1	B	80	ALA	3.1
1	H	134	VAL	3.1
1	D	23	LEU	3.1
1	H	126	GLY	3.1
1	E	138	PRO	3.0
1	E	135	ALA	3.0
1	H	125	LYS	3.0
1	C	47	LYS	3.0
1	A	133	GLY	3.0
1	C	20	GLY	3.0
1	F	119	GLY	3.0
1	B	42	VAL	3.0
1	B	56	VAL	3.0
1	H	120	THR	3.0
1	H	132	ILE	2.9
1	H	49	GLY	2.9
1	E	70	ASP	2.9
1	H	99	LYS	2.9
1	G	142	LEU	2.9
1	C	17	ALA	2.9
1	C	55	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	E	43	GLU	2.9
1	H	127	ALA	2.9
1	G	134	VAL	2.8
1	G	146	TYR	2.8
1	C	154	LEU	2.8
1	A	130	ALA	2.8
1	G	56	VAL	2.8
1	A	116	LEU	2.8
1	B	20	GLY	2.8
1	A	83	LYS	2.7
1	F	86	THR	2.7
1	B	119	GLY	2.7
1	F	133	GLY	2.7
1	F	56	VAL	2.7
1	A	103	GLN	2.7
1	D	18	ALA	2.7
1	E	137	ALA	2.7
1	C	132	ILE	2.7
1	D	38	ALA	2.7
1	C	110	ILE	2.6
1	D	24	THR	2.6
1	D	133	GLY	2.6
1	H	57	VAL	2.6
1	G	103	GLN	2.6
1	E	17	ALA	2.6
1	G	84	ALA	2.6
1	C	133	GLY	2.6
1	E	147	ALA	2.6
1	A	43	GLU	2.6
1	A	122	VAL	2.6
1	B	118	GLY	2.6
1	F	126	GLY	2.5
1	H	20	GLY	2.5
1	B	34	ALA	2.5
1	C	145	GLN	2.5
1	C	37	ALA	2.5
1	F	94	THR	2.5
1	B	88	VAL	2.5
1	D	103	GLN	2.5
1	C	21	GLY	2.5
1	A	16	VAL	2.5
1	B	16	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	130	ALA	2.5
1	F	132	ILE	2.5
1	C	103	GLN	2.5
1	E	27	THR	2.4
1	E	116	LEU	2.4
1	F	134	VAL	2.4
1	H	16	VAL	2.4
1	E	99	LYS	2.4
1	F	135	ALA	2.4
1	H	115	PHE	2.4
1	D	132	ILE	2.4
1	B	114	LEU	2.4
1	G	46	GLU	2.4
1	D	49	GLY	2.4
1	E	52	VAL	2.4
1	G	86	THR	2.4
1	F	20	GLY	2.4
1	A	86	THR	2.3
1	E	102	ALA	2.3
1	E	142	LEU	2.3
1	B	35	THR	2.3
1	G	138	PRO	2.3
1	H	104	ALA	2.3
1	H	142	LEU	2.3
1	B	83	LYS	2.3
1	A	17	ALA	2.3
1	B	17	ALA	2.3
1	F	130	ALA	2.3
1	E	118	GLY	2.3
1	B	76	SER	2.3
1	F	37	ALA	2.3
1	A	67	LEU	2.2
1	A	118	GLY	2.2
1	C	134	VAL	2.2
1	D	50	ARG	2.2
1	E	117	ALA	2.2
1	F	113	THR	2.2
1	D	153	VAL	2.2
1	F	57	VAL	2.2
1	E	36	LYS	2.2
1	G	125	LYS	2.2
1	G	70	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	114	LEU	2.2
1	A	56	VAL	2.2
1	D	79	SER	2.2
1	E	96	GLU	2.2
1	C	152	ALA	2.2
1	F	84	ALA	2.2
1	H	103	GLN	2.2
1	C	64	LEU	2.2
1	G	29	LEU	2.2
1	H	23	LEU	2.2
1	B	147	ALA	2.2
1	F	83	LYS	2.2
1	H	130	ALA	2.2
1	B	67	LEU	2.1
1	E	56	VAL	2.1
1	G	47	LYS	2.1
1	B	103	GLN	2.1
1	F	110	ILE	2.1
1	G	145	GLN	2.1
1	B	64	LEU	2.1
1	F	116	LEU	2.1
1	D	135	ALA	2.1
1	B	150	GLY	2.1
1	H	131	GLY	2.1
1	G	65	VAL	2.1
1	H	146	TYR	2.1
1	A	29	LEU	2.1
1	A	51	HIS	2.1
1	D	32	GLU	2.1
1	B	52	VAL	2.1
1	B	85	PHE	2.1
1	C	142	LEU	2.0
1	E	151	ALA	2.0
1	H	135	ALA	2.0
1	F	49	GLY	2.0
1	G	119	GLY	2.0
1	A	33	ALA	2.0
1	A	79	SER	2.0
1	C	149	ALA	2.0
1	G	48	ASP	2.0
1	D	56	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, 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	FE	E	157	1/1	0.50	0.32	2.00	132,132,132,132	0
2	FE	E	158	1/1	0.87	0.26	1.43	99,99,99,99	0
2	FE	B	157	1/1	0.99	0.15	-2.40	55,55,55,55	0
2	FE	D	157	1/1	0.95	0.08	-2.97	72,72,72,72	0
2	FE	A	157	1/1	0.98	0.07	-3.16	72,72,72,72	0
2	FE	F	157	1/1	0.93	0.09	-3.22	69,69,69,69	0
2	FE	C	157	1/1	0.91	0.05	-3.59	77,77,77,77	0
2	FE	H	157	1/1	0.89	0.07	-3.90	80,80,80,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.