



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 11:54 pm GMT

PDB ID : 1VRB
Title : Crystal structure of Putative asparaginyl hydroxylase (2636534) from *Bacillus subtilis* at 2.60 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2005-02-18
Resolution : 2.60 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 : trunk28620
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) : recalc28949

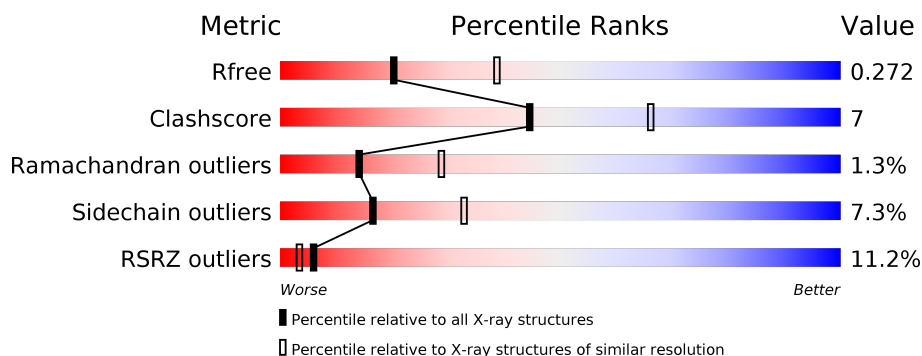
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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	342	<div> <div>11%</div> <div> <div></div> <div>71%</div> <div>17%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	342	<div> <div>9%</div> <div> <div></div> <div>72%</div> <div>17%</div> <div>•</div> <div>9%</div> </div> </div>
1	C	342	<div> <div>9%</div> <div> <div></div> <div>68%</div> <div>18%</div> <div>•</div> <div>13%</div> </div> </div>
1	D	342	<div> <div>11%</div> <div> <div></div> <div>70%</div> <div>19%</div> <div>•</div> <div>8%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9036 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 Putative asparaginyl hydroxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	Se	0	0	0
			2223	1456	367	395	5			
1	B	311	Total	C	N	O	Se	0	0	0
			2283	1491	378	409	5			
1	C	298	Total	C	N	O	Se	0	0	0
			2210	1446	367	392	5			
1	D	313	Total	C	N	O	Se	0	0	0
			2294	1498	385	406	5			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MSE	-	LEADER SEQUENCE	UNP P46327
A	-10	GLY	-	LEADER SEQUENCE	UNP P46327
A	-9	SER	-	LEADER SEQUENCE	UNP P46327
A	-8	ASP	-	LEADER SEQUENCE	UNP P46327
A	-7	LYS	-	LEADER SEQUENCE	UNP P46327
A	-6	ILE	-	LEADER SEQUENCE	UNP P46327
A	-5	HIS	-	LEADER SEQUENCE	UNP P46327
A	-4	HIS	-	LEADER SEQUENCE	UNP P46327
A	-3	HIS	-	LEADER SEQUENCE	UNP P46327
A	-2	HIS	-	LEADER SEQUENCE	UNP P46327
A	-1	HIS	-	LEADER SEQUENCE	UNP P46327
A	0	HIS	-	LEADER SEQUENCE	UNP P46327
A	1	MSE	MET	MODIFIED RESIDUE	UNP P46327
A	18	MSE	MET	MODIFIED RESIDUE	UNP P46327
A	64	MSE	MET	MODIFIED RESIDUE	UNP P46327
A	174	MSE	MET	MODIFIED RESIDUE	UNP P46327
A	215	MSE	MET	MODIFIED RESIDUE	UNP P46327
A	248	MSE	MET	MODIFIED RESIDUE	UNP P46327
B	-11	MSE	-	LEADER SEQUENCE	UNP P46327
B	-10	GLY	-	LEADER SEQUENCE	UNP P46327
B	-9	SER	-	LEADER SEQUENCE	UNP P46327

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	ASP	-	LEADER SEQUENCE	UNP P46327
B	-7	LYS	-	LEADER SEQUENCE	UNP P46327
B	-6	ILE	-	LEADER SEQUENCE	UNP P46327
B	-5	HIS	-	LEADER SEQUENCE	UNP P46327
B	-4	HIS	-	LEADER SEQUENCE	UNP P46327
B	-3	HIS	-	LEADER SEQUENCE	UNP P46327
B	-2	HIS	-	LEADER SEQUENCE	UNP P46327
B	-1	HIS	-	LEADER SEQUENCE	UNP P46327
B	0	HIS	-	LEADER SEQUENCE	UNP P46327
B	1	MSE	MET	MODIFIED RESIDUE	UNP P46327
B	18	MSE	MET	MODIFIED RESIDUE	UNP P46327
B	64	MSE	MET	MODIFIED RESIDUE	UNP P46327
B	174	MSE	MET	MODIFIED RESIDUE	UNP P46327
B	215	MSE	MET	MODIFIED RESIDUE	UNP P46327
B	248	MSE	MET	MODIFIED RESIDUE	UNP P46327
C	-11	MSE	-	LEADER SEQUENCE	UNP P46327
C	-10	GLY	-	LEADER SEQUENCE	UNP P46327
C	-9	SER	-	LEADER SEQUENCE	UNP P46327
C	-8	ASP	-	LEADER SEQUENCE	UNP P46327
C	-7	LYS	-	LEADER SEQUENCE	UNP P46327
C	-6	ILE	-	LEADER SEQUENCE	UNP P46327
C	-5	HIS	-	LEADER SEQUENCE	UNP P46327
C	-4	HIS	-	LEADER SEQUENCE	UNP P46327
C	-3	HIS	-	LEADER SEQUENCE	UNP P46327
C	-2	HIS	-	LEADER SEQUENCE	UNP P46327
C	-1	HIS	-	LEADER SEQUENCE	UNP P46327
C	0	HIS	-	LEADER SEQUENCE	UNP P46327
C	1	MSE	MET	MODIFIED RESIDUE	UNP P46327
C	18	MSE	MET	MODIFIED RESIDUE	UNP P46327
C	64	MSE	MET	MODIFIED RESIDUE	UNP P46327
C	174	MSE	MET	MODIFIED RESIDUE	UNP P46327
C	215	MSE	MET	MODIFIED RESIDUE	UNP P46327
C	248	MSE	MET	MODIFIED RESIDUE	UNP P46327
D	-11	MSE	-	LEADER SEQUENCE	UNP P46327
D	-10	GLY	-	LEADER SEQUENCE	UNP P46327
D	-9	SER	-	LEADER SEQUENCE	UNP P46327
D	-8	ASP	-	LEADER SEQUENCE	UNP P46327
D	-7	LYS	-	LEADER SEQUENCE	UNP P46327
D	-6	ILE	-	LEADER SEQUENCE	UNP P46327
D	-5	HIS	-	LEADER SEQUENCE	UNP P46327
D	-4	HIS	-	LEADER SEQUENCE	UNP P46327
D	-3	HIS	-	LEADER SEQUENCE	UNP P46327

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	HIS	-	LEADER SEQUENCE	UNP P46327
D	-1	HIS	-	LEADER SEQUENCE	UNP P46327
D	0	HIS	-	LEADER SEQUENCE	UNP P46327
D	1	MSE	MET	MODIFIED RESIDUE	UNP P46327
D	18	MSE	MET	MODIFIED RESIDUE	UNP P46327
D	64	MSE	MET	MODIFIED RESIDUE	UNP P46327
D	174	MSE	MET	MODIFIED RESIDUE	UNP P46327
D	215	MSE	MET	MODIFIED RESIDUE	UNP P46327
D	248	MSE	MET	MODIFIED RESIDUE	UNP P46327

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

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

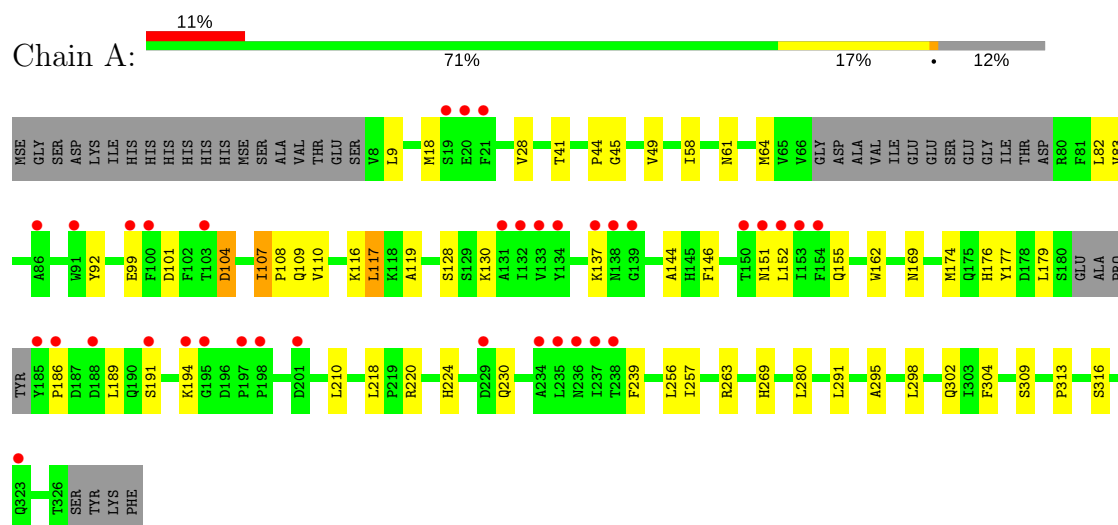
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	6	Total O 6 6	0	0
3	B	3	Total O 3 3	0	0
3	C	5	Total O 5 5	0	0
3	D	8	Total O 8 8	0	0

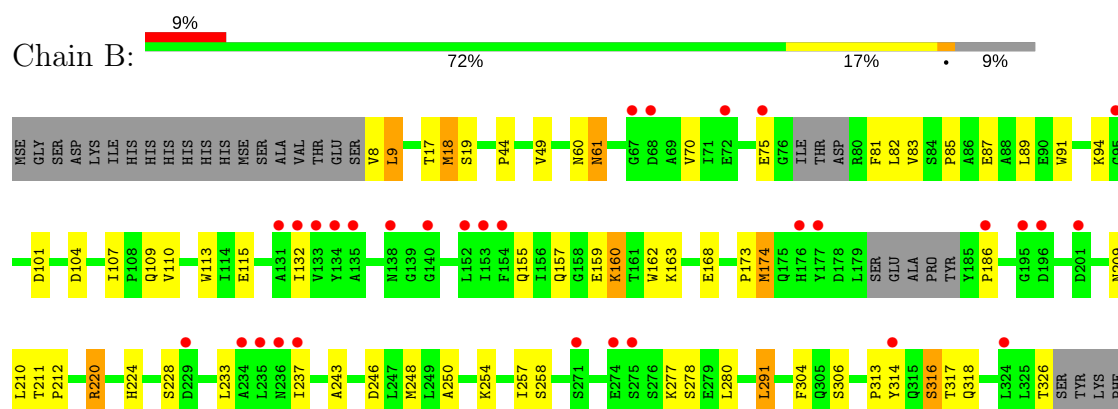
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

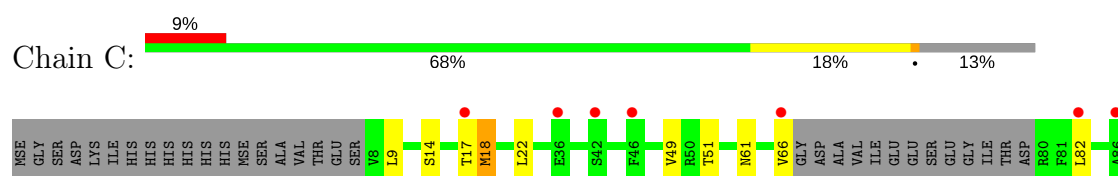
• Molecule 1: Putative asparaginyl hydroxylase

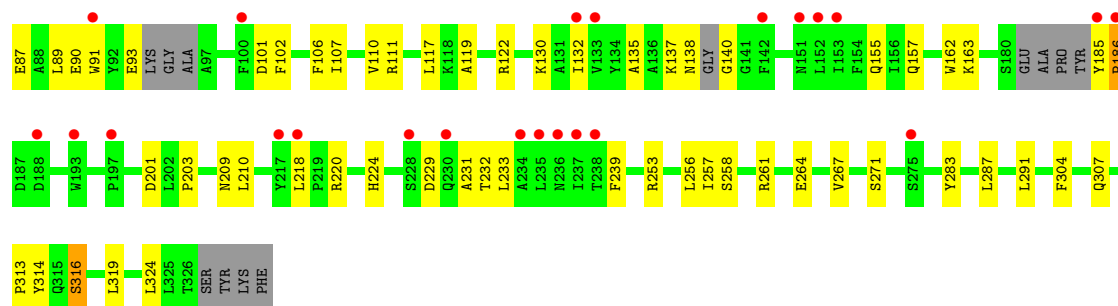


• Molecule 1: Putative asparaginyl hydroxylase

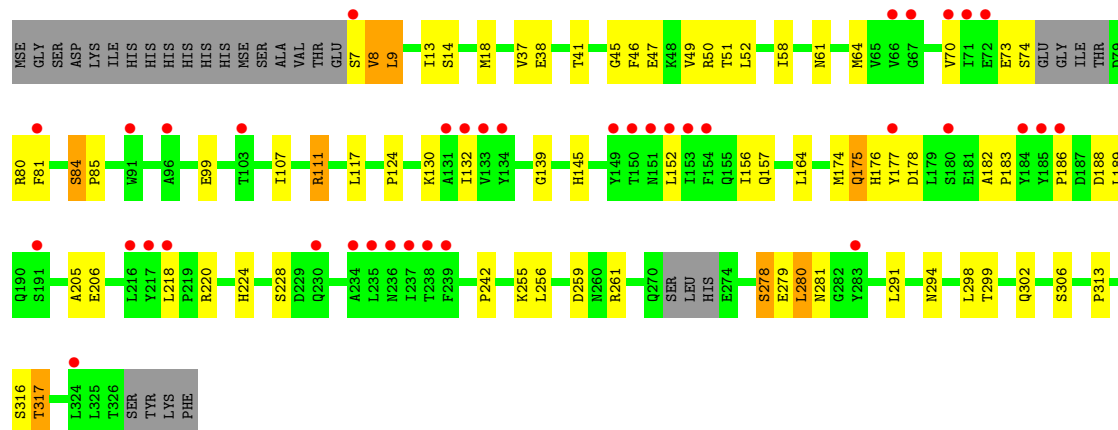


• Molecule 1: Putative asparaginyl hydroxylase





• Molecule 1: Putative asparaginyl hydroxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.30Å 104.66Å 286.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.66 – 2.60 28.66 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.3 (28.66-2.60) 99.4 (28.66-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.221 , 0.272 0.224 , 0.272	Depositor DCC
R_{free} test set	2243 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	69.9	Xtriage
Anisotropy	0.432	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 84.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9036	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 3.24% 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

5.1 Standard geometry

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	0.75	0/2275	0.79	0/3116
1	B	0.74	0/2336	0.76	1/3200 (0.0%)
1	C	0.68	1/2260 (0.0%)	0.72	0/3093
1	D	0.81	3/2347 (0.1%)	0.80	1/3214 (0.0%)
All	All	0.75	4/9218 (0.0%)	0.77	2/12623 (0.0%)

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	D	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	93	GLU	C-O	10.32	1.43	1.23
1	D	74	SER	CA-CB	9.21	1.66	1.52
1	D	70	VAL	C-O	6.41	1.35	1.23
1	D	73	GLU	CA-C	6.13	1.68	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	9	LEU	CA-CB-CG	5.36	127.62	115.30
1	D	111	ARG	NE-CZ-NH1	5.11	122.86	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	182	ALA	Peptide
1	D	7	SER	Peptide
1	D	80	ARG	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	2223	0	2053	31	0
1	B	2283	0	2106	39	0
1	C	2210	0	2037	36	0
1	D	2294	0	2135	36	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
3	A	6	0	0	0	0
3	B	3	0	0	0	0
3	C	5	0	0	0	0
3	D	8	0	0	0	0
All	All	9036	0	8331	130	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 7.

All (130) 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:18:MSE:CE	1:D:18:MSE:SE	2.15	1.44
1:B:18:MSE:CE	1:B:18:MSE:SE	2.15	1.43
1:B:155:GLN:HG3	1:B:210:LEU:HB3	1.60	0.82
1:D:256:LEU:HD11	1:D:291:LEU:HD12	1.73	0.71
1:C:49:VAL:HG12	1:C:49:VAL:O	1.90	0.71
1:C:256:LEU:HD11	1:C:291:LEU:HD12	1.73	0.70
1:C:138:ASN:C	1:C:140:GLY:N	2.46	0.69
1:D:313:PRO:O	1:D:317:THR:HG23	1.93	0.69
1:B:159:GLU:HG2	1:B:211:THR:HG22	1.73	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:TRP:O	1:B:94:LYS:O	2.11	0.67
1:D:9:LEU:HD22	1:D:117:LEU:HD11	1.76	0.66
1:A:174:MSE:HG3	1:B:257:ILE:HD11	1.77	0.66
1:A:257:ILE:HD11	1:B:174:MSE:HG3	1.78	0.65
1:D:145:HIS:HB2	1:D:177:TYR:O	1.97	0.64
1:A:155:GLN:HG3	1:A:210:LEU:HB3	1.80	0.62
1:A:144:ALA:HB3	1:A:179:LEU:HD12	1.81	0.62
1:C:314:TYR:HA	1:D:317:THR:HG21	1.82	0.62
1:D:278:SER:O	1:D:280:LEU:N	2.33	0.61
1:C:313:PRO:HA	1:C:316:SER:HB2	1.82	0.60
1:B:82:LEU:O	1:B:83:VAL:HG23	2.01	0.60
1:C:256:LEU:CD1	1:C:291:LEU:HD12	2.35	0.56
1:D:313:PRO:HA	1:D:316:SER:OG	2.05	0.56
1:C:261:ARG:NE	1:C:261:ARG:HA	2.20	0.56
1:A:45:GLY:HA3	1:A:107:ILE:HG21	1.88	0.56
1:C:162:TRP:NE1	1:C:210:LEU:HD22	2.20	0.56
1:B:254:LYS:O	1:B:257:ILE:HG22	2.07	0.55
1:D:176:HIS:NE2	1:D:178:ASP:CG	2.60	0.55
1:B:314:TYR:CE1	1:B:318:GLN:NE2	2.77	0.53
1:C:130:LYS:HE3	1:C:132:ILE:HD11	1.89	0.53
1:D:45:GLY:HA3	1:D:107:ILE:HG21	1.91	0.53
1:B:87:GLU:O	1:B:91:TRP:HD1	1.92	0.53
1:C:155:GLN:HG3	1:C:210:LEU:HB3	1.91	0.53
1:A:44:PRO:O	1:A:110:VAL:HG12	2.08	0.53
1:C:117:LEU:HD23	1:C:239:PHE:HZ	1.74	0.53
1:C:257:ILE:HD11	1:D:174:MSE:HG3	1.91	0.53
1:C:324:LEU:HD13	1:D:64:MSE:HE1	1.91	0.52
1:B:313:PRO:HA	1:B:316:SER:HB2	1.91	0.52
1:C:218:LEU:HD21	1:C:224:HIS:CE1	2.45	0.52
1:D:130:LYS:HE3	1:D:132:ILE:HD11	1.90	0.52
1:A:9:LEU:HD23	1:A:18:MSE:HE1	1.91	0.52
1:A:256:LEU:HD11	1:A:291:LEU:CD1	2.40	0.51
1:A:104:ASP:N	1:A:104:ASP:OD1	2.44	0.51
1:A:49:VAL:O	1:A:49:VAL:CG1	2.59	0.51
1:B:243:ALA:O	1:B:246:ASP:HB2	2.10	0.51
1:A:291:LEU:HD22	1:B:248:MSE:HE2	1.93	0.50
1:A:176:HIS:CG	1:A:177:TYR:N	2.79	0.50
1:D:9:LEU:HD23	1:D:18:MSE:HE1	1.93	0.50
1:C:17:THR:O	1:C:18:MSE:HB2	2.12	0.50
1:C:117:LEU:HD23	1:C:239:PHE:CZ	2.47	0.50
1:A:92:TYR:CE1	1:A:137:LYS:HB2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:LYS:O	1:A:119:ALA:HB3	2.12	0.49
1:A:218:LEU:HD21	1:A:224:HIS:ND1	2.27	0.49
1:A:295:ALA:HB1	1:B:291:LEU:HD23	1.95	0.48
1:B:101:ASP:HA	1:B:132:ILE:HG23	1.95	0.48
1:C:163:LYS:HE2	1:C:201:ASP:O	2.13	0.48
1:D:139:GLY:HA2	1:D:228:SER:O	2.13	0.48
1:D:255:LYS:HE2	1:D:294:ASN:O	2.14	0.48
1:C:137:LYS:HA	1:C:231:ALA:HA	1.95	0.48
1:B:49:VAL:O	1:B:49:VAL:HG12	2.14	0.47
1:A:298:LEU:HD12	1:A:302:GLN:OE1	2.14	0.47
1:D:117:LEU:HD21	1:D:152:LEU:CD1	2.44	0.47
1:D:256:LEU:HD11	1:D:291:LEU:CD1	2.44	0.47
1:C:261:ARG:HA	1:C:261:ARG:HE	1.79	0.47
1:D:164:LEU:O	1:D:205:ALA:HB1	2.15	0.47
1:C:102:PHE:CD1	1:C:319:LEU:HD13	2.49	0.47
1:D:84:SER:OG	1:D:85:PRO:HD2	2.14	0.47
1:A:280:LEU:HD23	1:A:280:LEU:HA	1.76	0.46
1:B:44:PRO:HB2	1:B:109:GLN:HB3	1.97	0.46
1:C:119:ALA:O	1:C:122:ARG:N	2.43	0.46
1:A:101:ASP:OD1	1:A:130:LYS:HD2	2.15	0.46
1:C:89:LEU:C	1:C:91:TRP:H	2.19	0.46
1:D:175:GLN:NE2	1:D:188:ASP:HB2	2.31	0.46
1:A:313:PRO:HA	1:A:316:SER:OG	2.15	0.46
1:A:82:LEU:O	1:A:83:VAL:HG23	2.16	0.46
1:A:162:TRP:NE1	1:A:210:LEU:HD22	2.31	0.46
1:B:87:GLU:O	1:B:91:TRP:CD1	2.69	0.46
1:B:173:PRO:HG2	1:B:220:ARG:HD3	1.97	0.45
1:B:60:ASN:ND2	1:B:326:THR:C	2.70	0.45
1:B:160:LYS:HG3	1:B:228:SER:HB2	1.97	0.45
1:D:9:LEU:HD22	1:D:117:LEU:CD1	2.44	0.45
1:B:104:ASP:N	1:B:104:ASP:OD1	2.49	0.45
1:A:218:LEU:HD21	1:A:224:HIS:CE1	2.52	0.45
1:C:135:ALA:HB2	1:C:233:LEU:HD12	1.99	0.45
1:B:17:THR:O	1:B:18:MSE:HB2	2.18	0.44
1:C:106:PHE:C	1:C:107:ILE:HG13	2.36	0.44
1:C:185:TYR:O	1:C:186:PRO:C	2.56	0.44
1:B:157:GLN:HB2	1:B:233:LEU:HB3	2.01	0.43
1:B:162:TRP:NE1	1:B:210:LEU:HD22	2.33	0.43
1:B:8:VAL:HG12	1:B:9:LEU:N	2.34	0.43
1:C:49:VAL:CG1	1:C:49:VAL:O	2.62	0.43
1:A:117:LEU:HD22	1:A:239:PHE:CZ	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:46:PHE:O	1:D:47:GLU:C	2.56	0.43
1:B:277:LYS:O	1:B:280:LEU:N	2.42	0.43
1:B:61:ASN:O	1:B:85:PRO:HG3	2.18	0.43
1:A:64:MSE:HB3	1:A:99:GLU:HB3	1.99	0.43
1:A:269:HIS:CE1	1:B:304:PHE:HB3	2.54	0.43
1:D:299:THR:HG23	1:D:302:GLN:NE2	2.33	0.43
1:C:101:ASP:CG	1:C:130:LYS:HZ2	2.22	0.42
1:B:209:ASN:HB2	1:C:203:PRO:HB2	2.01	0.42
1:D:259:ASP:OD1	1:D:261:ARG:HD3	2.20	0.42
1:D:38:GLU:O	1:D:41:THR:OG1	2.28	0.42
1:D:218:LEU:HD21	1:D:224:HIS:CE1	2.55	0.42
1:B:107:ILE:O	1:B:110:VAL:HG12	2.20	0.42
1:B:250:ALA:HB3	1:B:306:SER:OG	2.20	0.41
1:B:60:ASN:HD21	1:B:326:THR:C	2.23	0.41
1:D:175:GLN:HE21	1:D:188:ASP:HB2	1.84	0.41
1:D:298:LEU:HA	1:D:298:LEU:HD12	1.90	0.41
1:B:113:TRP:HZ3	1:B:237:ILE:HD12	1.84	0.41
1:B:163:LYS:O	1:B:224:HIS:HA	2.20	0.41
1:C:313:PRO:HA	1:C:316:SER:CB	2.50	0.41
1:C:304:PHE:O	1:C:307:GLN:HG2	2.20	0.41
1:C:264:GLU:OE1	1:C:283:TYR:OH	2.35	0.41
1:D:13:ILE:O	1:D:14:SER:C	2.59	0.41
1:B:277:LYS:O	1:B:278:SER:C	2.57	0.41
1:C:87:GLU:O	1:C:91:TRP:CD1	2.74	0.41
1:A:108:PRO:O	1:A:109:GLN:C	2.60	0.41
1:A:146:PHE:CD1	1:A:146:PHE:C	2.94	0.41
1:B:211:THR:HB	1:B:212:PRO:CD	2.51	0.41
1:C:314:TYR:HD1	1:D:317:THR:HG22	1.85	0.41
1:A:169:ASN:HA	1:A:194:LYS:O	2.20	0.41
1:C:66:VAL:O	1:C:66:VAL:HG12	2.21	0.41
1:A:263:ARG:HG2	1:B:174:MSE:HE2	2.03	0.41
1:A:304:PHE:CZ	1:B:280:LEU:HD23	2.55	0.41
1:D:124:PRO:HG2	1:D:242:PRO:HG2	2.03	0.41
1:D:278:SER:O	1:D:281:ASN:N	2.53	0.40
1:C:17:THR:OG1	1:C:18:MSE:N	2.54	0.40
1:D:37:VAL:O	1:D:38:GLU:C	2.60	0.40
1:C:82:LEU:HA	1:C:82:LEU:HD12	1.96	0.40
1:D:50:ARG:HA	1:D:156:ILE:HG21	2.03	0.40
1:D:64:MSE:HB3	1:D:99:GLU:HB3	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	296/342 (86%)	275 (93%)	18 (6%)	3 (1%)	18	37
1	B	305/342 (89%)	280 (92%)	21 (7%)	4 (1%)	14	29
1	C	288/342 (84%)	266 (92%)	18 (6%)	4 (1%)	13	26
1	D	307/342 (90%)	281 (92%)	21 (7%)	5 (2%)	11	23
All	All	1196/1368 (87%)	1102 (92%)	78 (6%)	16 (1%)	14	29

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	186	PRO
1	B	18	MSE
1	B	75	GLU
1	C	14	SER
1	C	18	MSE
1	D	183	PRO
1	A	309	SER
1	C	90	GLU
1	D	278	SER
1	D	279	GLU
1	D	186	PRO
1	A	41	THR
1	C	186	PRO
1	D	8	VAL
1	B	186	PRO
1	B	70	VAL

5.3.2 Protein sidechains [i](#)

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	205/294 (70%)	192 (94%)	13 (6%)	21	42
1	B	212/294 (72%)	199 (94%)	13 (6%)	22	43
1	C	206/294 (70%)	189 (92%)	17 (8%)	13	25
1	D	213/294 (72%)	195 (92%)	18 (8%)	12	24
All	All	836/1176 (71%)	775 (93%)	61 (7%)	16	33

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	VAL
1	A	58	ILE
1	A	61	ASN
1	A	104	ASP
1	A	107	ILE
1	A	117	LEU
1	A	128	SER
1	A	151	ASN
1	A	152	LEU
1	A	189	LEU
1	A	191	SER
1	A	220	ARG
1	A	230	GLN
1	B	19	SER
1	B	61	ASN
1	B	81	PHE
1	B	89	LEU
1	B	115	GLU
1	B	160	LYS
1	B	168	GLU
1	B	174	MSE
1	B	220	ARG
1	B	258	SER
1	B	291	LEU
1	B	316	SER
1	B	317	THR
1	C	9	LEU
1	C	22	LEU
1	C	51	THR
1	C	61	ASN

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Mol	Chain	Res	Type
1	C	110	VAL
1	C	111	ARG
1	C	157	GLN
1	C	209	ASN
1	C	220	ARG
1	C	229	ASP
1	C	232	THR
1	C	253	ARG
1	C	258	SER
1	C	267	VAL
1	C	271	SER
1	C	287	LEU
1	C	316	SER
1	D	8	VAL
1	D	9	LEU
1	D	49	VAL
1	D	51	THR
1	D	52	LEU
1	D	58	ILE
1	D	61	ASN
1	D	81	PHE
1	D	84	SER
1	D	111	ARG
1	D	157	GLN
1	D	175	GLN
1	D	189	LEU
1	D	206	GLU
1	D	220	ARG
1	D	280	LEU
1	D	306	SER
1	D	317	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	157	GLN
1	B	60	ASN
1	B	138	ASN
1	B	155	GLN
1	B	318	GLN
1	C	155	GLN

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Mol	Chain	Res	Type
1	C	270	GLN
1	C	307	GLN
1	D	268	ASN

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 4 ligands modelled in this entry, 4 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	297/342 (86%)	0.52	36 (12%) 5 3	59, 67, 76, 83	0
1	B	306/342 (89%)	0.35	31 (10%) 8 5	59, 67, 75, 87	0
1	C	293/342 (85%)	0.36	30 (10%) 7 4	60, 67, 76, 84	0
1	D	308/342 (90%)	0.38	38 (12%) 5 2	58, 67, 76, 91	0
All	All	1204/1368 (88%)	0.40	135 (11%) 6 3	58, 67, 76, 91	0

All (135) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	67	GLY	7.2
1	A	236	ASN	6.2
1	A	153	ILE	5.4
1	C	236	ASN	5.3
1	C	132	ILE	5.1
1	D	185	TYR	4.9
1	D	153	ILE	4.9
1	A	132	ILE	4.8
1	D	152	LEU	4.8
1	D	236	ASN	4.7
1	A	151	ASN	4.7
1	D	151	ASN	4.6
1	A	235	LEU	4.5
1	B	235	LEU	4.4
1	C	237	ILE	4.3
1	B	133	VAL	4.3
1	D	72	GLU	4.2
1	A	133	VAL	4.1
1	D	234	ALA	4.1
1	A	134	TYR	4.0
1	A	234	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	237	ILE	3.9
1	D	235	LEU	3.8
1	A	139	GLY	3.8
1	C	151	ASN	3.8
1	D	66	VAL	3.7
1	D	184	TYR	3.7
1	C	185	TYR	3.6
1	C	238	THR	3.6
1	A	237	ILE	3.5
1	A	152	LEU	3.5
1	A	198	PRO	3.5
1	C	153	ILE	3.4
1	D	71	ILE	3.4
1	B	131	ALA	3.3
1	A	185	TYR	3.3
1	A	86	ALA	3.3
1	B	195	GLY	3.2
1	B	234	ALA	3.2
1	C	91	TRP	3.2
1	D	7	SER	3.2
1	D	186	PRO	3.2
1	D	238	THR	3.1
1	C	46	PHE	3.0
1	A	99	GLU	3.0
1	B	132	ILE	3.0
1	B	95	GLY	3.0
1	D	230	GLN	3.0
1	C	133	VAL	2.9
1	B	176	HIS	2.9
1	B	67	GLY	2.9
1	D	133	VAL	2.9
1	A	138	ASN	2.9
1	B	271	SER	2.9
1	A	154	PHE	2.9
1	A	195	GLY	2.8
1	C	193	TRP	2.8
1	B	229	ASP	2.8
1	A	131	ALA	2.8
1	B	274	GLU	2.8
1	D	218	LEU	2.8
1	B	134	TYR	2.7
1	A	100	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	42	SER	2.7
1	D	131	ALA	2.7
1	B	140	GLY	2.7
1	B	314	TYR	2.7
1	B	237	ILE	2.7
1	B	196	ASP	2.7
1	B	135	ALA	2.6
1	D	134	TYR	2.6
1	C	152	LEU	2.6
1	A	191	SER	2.6
1	C	197	PRO	2.5
1	A	188	ASP	2.5
1	B	177	TYR	2.5
1	B	201	ASP	2.5
1	C	234	ALA	2.5
1	B	138	ASN	2.5
1	C	217	TYR	2.5
1	A	229	ASP	2.5
1	C	235	LEU	2.5
1	C	82	LEU	2.4
1	C	228	SER	2.4
1	C	100	PHE	2.4
1	B	72	GLU	2.4
1	D	150	THR	2.4
1	D	96	ALA	2.4
1	D	239	PHE	2.4
1	A	103	THR	2.4
1	B	153	ILE	2.4
1	D	216	LEU	2.4
1	A	20	GLU	2.4
1	C	17	THR	2.3
1	D	132	ILE	2.3
1	C	142	PHE	2.3
1	D	180	SER	2.3
1	A	201	ASP	2.3
1	B	75	GLU	2.3
1	D	191	SER	2.3
1	D	103	THR	2.3
1	C	36	GLU	2.3
1	C	188	ASP	2.3
1	A	137	LYS	2.3
1	B	68	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	323	GLN	2.2
1	A	194	LYS	2.2
1	C	230	GLN	2.2
1	C	218	LEU	2.2
1	D	154	PHE	2.2
1	C	275	SER	2.2
1	C	186	PRO	2.2
1	B	236	ASN	2.2
1	B	186	PRO	2.2
1	B	154	PHE	2.2
1	D	217	TYR	2.2
1	A	197	PRO	2.2
1	A	186	PRO	2.2
1	C	86	ALA	2.2
1	B	152	LEU	2.1
1	B	324	LEU	2.1
1	D	70	VAL	2.1
1	D	149	TYR	2.1
1	D	177	TYR	2.1
1	D	324	LEU	2.1
1	A	91	TRP	2.1
1	A	150	THR	2.1
1	A	19	SER	2.1
1	D	283	TYR	2.1
1	D	91	TRP	2.1
1	C	66	VAL	2.0
1	B	275	SER	2.0
1	A	238	THR	2.0
1	A	21	PHE	2.0
1	D	81	PHE	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 ⓘ

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	C	555	1/1	0.75	0.14	-	104,104,104,104	0
2	FE	B	555	1/1	0.84	0.18	-	104,104,104,104	0
2	FE	D	555	1/1	0.95	0.04	-	85,85,85,85	0
2	FE	A	555	1/1	0.91	0.24	-	104,104,104,104	0

6.5 Other polymers [i](#)

There are no such residues in this entry.