



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 18, 2017 – 02:48 PM EDT

PDB ID : 6ALX
Title : Structure of F. tularensis MglA-SspA solved in the presence of polyP
Authors : Schumacher, M.A.
Deposited on : unknown
Resolution : 3.30 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
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
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-20029824

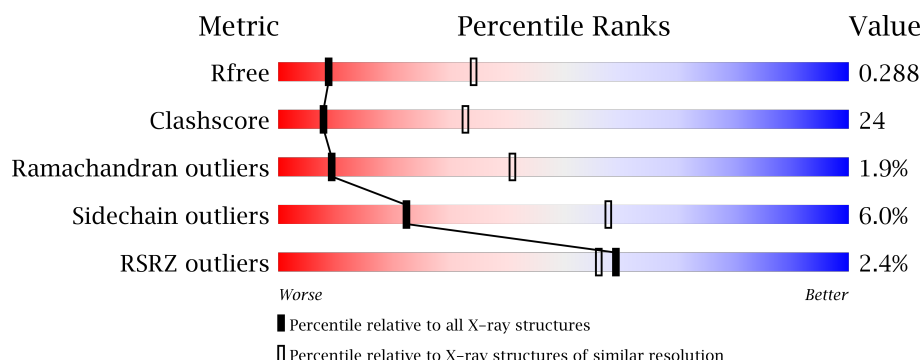
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	100719	1034 (3.36-3.24)
Clashscore	112137	1100 (3.36-3.24)
Ramachandran outliers	110173	1081 (3.36-3.24)
Sidechain outliers	110143	1080 (3.36-3.24)
RSRZ outliers	101464	1039 (3.36-3.24)

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	B	211	<div> <div>5%</div> <div>49%</div> <div>38%</div> <div>5%</div> <div>8%</div> </div>
1	C	211	<div> <div>%</div> <div>47%</div> <div>40%</div> <div>5%</div> <div>8%</div> </div>
2	A	204	<div> <div>%</div> <div>56%</div> <div>42%</div> <div>.</div> </div>
2	D	204	<div> <div>%</div> <div>55%</div> <div>43%</div> <div>.</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6230 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 Stringent starvation protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	195	Total	C	N	O	S	0	0	0
			1526	980	249	287	10			
1	B	195	Total	C	N	O	S	0	0	0
			1523	978	249	287	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP A0A0E2ZL39
C	1	ASN	-	expression tag	UNP A0A0E2ZL39
C	2	ALA	-	expression tag	UNP A0A0E2ZL39
C	3	MET	-	expression tag	UNP A0A0E2ZL39
B	0	SER	-	expression tag	UNP A0A0E2ZL39
B	1	ASN	-	expression tag	UNP A0A0E2ZL39
B	2	ALA	-	expression tag	UNP A0A0E2ZL39
B	3	MET	-	expression tag	UNP A0A0E2ZL39

- Molecule 2 is a protein called Macrophage growth locus A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	203	Total	C	N	O	S	0	0	0
			1595	1037	256	297	5			
2	D	203	Total	C	N	O	S	0	0	0
			1586	1033	253	295	5			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	SER	-	expression tag	UNP A0A0E2ZLH6
A	3	ASN	-	expression tag	UNP A0A0E2ZLH6
A	4	ALA	-	expression tag	UNP A0A0E2ZLH6
D	2	SER	-	expression tag	UNP A0A0E2ZLH6

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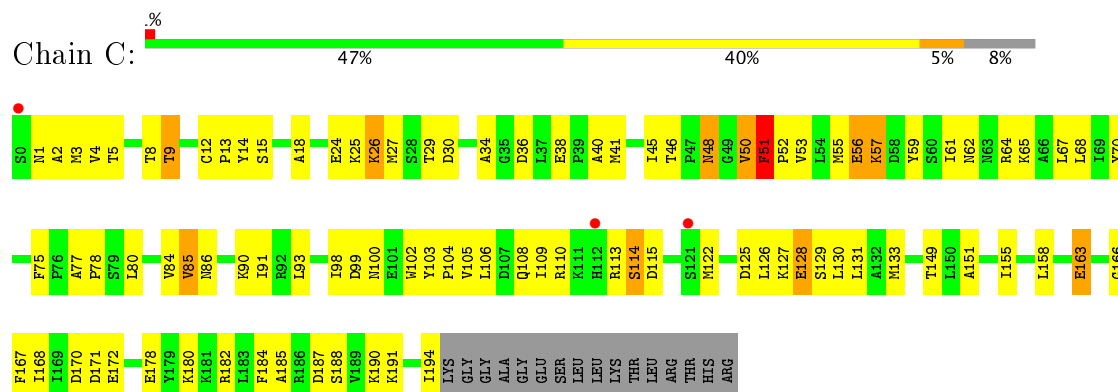
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Chain	Residue	Modelled	Actual	Comment	Reference
D	3	ASN	-	expression tag	UNP A0A0E2ZLH6
D	4	ALA	-	expression tag	UNP A0A0E2ZLH6

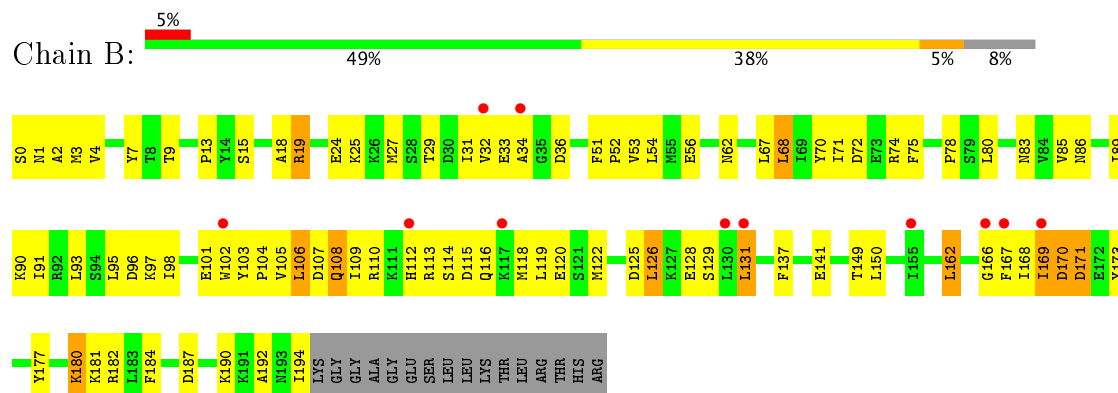
3 Residue-property plots [i](#)

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.

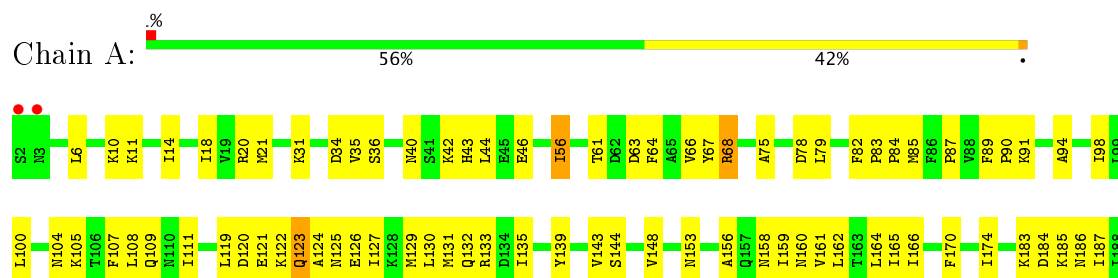
• Molecule 1: Stringent starvation protein A



• Molecule 1: Stringent starvation protein A

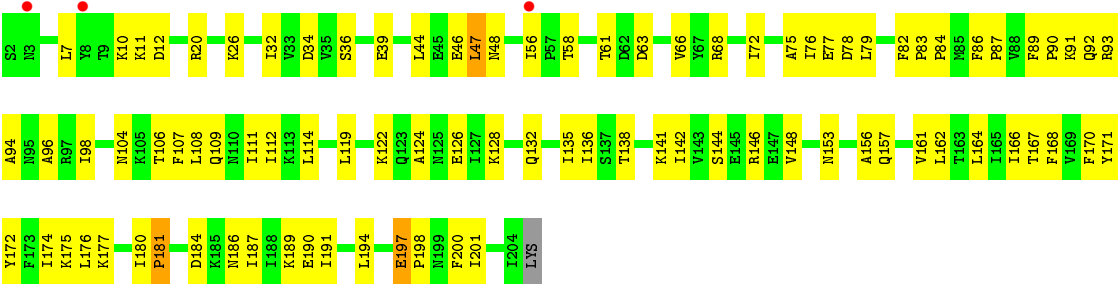


• Molecule 2: Macrophage growth locus A





● Molecule 2: Macrophage growth locus A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.25Å 114.30Å 141.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.75 – 3.30 60.75 – 3.30	Depositor EDS
% Data completeness (in resolution range)	94.2 (60.75-3.30) 93.0 (60.75-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.6.4 _486	Depositor
R, R_{free}	0.209 , 0.294 0.204 , 0.288	Depositor DCC
R_{free} test set	1597 reflections (10.06%)	DCC
Wilson B-factor (Å ²)	80.9	Xtriage
Anisotropy	0.352	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 73.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6230	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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](#)

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	B	0.32	0/1551	0.51	0/2100
1	C	0.39	0/1554	0.56	0/2103
2	A	0.34	0/1623	0.55	0/2210
2	D	0.33	0/1614	0.58	0/2199
All	All	0.35	0/6342	0.55	0/8612

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	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	51	PHE	Peptide

5.2 Too-close contacts [i](#)

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	B	1523	0	1503	81	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1526	0	1510	72	0
2	A	1595	0	1612	77	0
2	D	1586	0	1599	76	0
All	All	6230	0	6224	297	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 24.

All (297) 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)
2:A:119:LEU:HB3	2:A:123:GLN:HE21	1.21	1.02
2:A:186:ASN:HA	2:A:189:LYS:HG2	1.41	1.02
2:A:132:GLN:HE22	2:D:153:ASN:HD21	1.15	0.94
1:B:187:ASP:HA	1:B:190:LYS:HB2	1.49	0.93
2:A:119:LEU:HG	2:A:123:GLN:HG3	1.48	0.92
1:C:65:LYS:NZ	1:C:100:ASN:HD21	1.67	0.90
1:C:102:TRP:O	1:C:105:VAL:HG12	1.73	0.89
1:B:15:SER:HA	1:B:52:PRO:HG3	1.53	0.88
1:B:181:LYS:HA	1:B:184:PHE:CZ	2.09	0.87
2:A:61:THR:HG23	2:A:63:ASP:H	1.42	0.85
2:A:191:ILE:O	2:A:195:LEU:HD12	1.77	0.84
1:B:113:ARG:HA	1:B:119:LEU:HD21	1.60	0.83
1:B:141:GLU:HA	1:B:182:ARG:NH2	1.93	0.83
1:C:65:LYS:HZ1	1:C:100:ASN:HD21	1.25	0.83
2:D:119:LEU:HD23	2:D:124:ALA:HB2	1.60	0.83
2:D:122:LYS:O	2:D:126:GLU:HG2	1.79	0.82
1:B:169:ILE:HG23	1:B:180:LYS:HZ2	1.45	0.81
1:B:126:LEU:HD11	1:B:167:PHE:CZ	2.16	0.80
1:B:169:ILE:O	1:B:170:ASP:HB2	1.81	0.80
1:C:127:LYS:O	1:C:131:LEU:HG	1.82	0.79
1:C:5:THR:HG22	1:C:30:ASP:HB2	1.64	0.79
1:C:125:ASP:O	1:C:128:GLU:HG3	1.84	0.78
2:A:119:LEU:CB	2:A:123:GLN:HE21	1.97	0.78
2:A:119:LEU:HB3	2:A:123:GLN:NE2	1.98	0.78
1:B:102:TRP:HH2	1:B:129:SER:HB3	1.48	0.78
2:D:11:LYS:HD2	2:D:12:ASP:H	1.49	0.77
1:C:50:VAL:HG21	1:C:62:ASN:CG	2.04	0.77
1:C:25:LYS:O	1:C:26:LYS:HD2	1.85	0.77
2:D:142:ILE:HD12	2:D:166:ILE:HD12	1.68	0.76
1:B:169:ILE:HG23	1:B:180:LYS:NZ	2.02	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:153:ASN:ND2	2:A:156:ALA:HB2	2.02	0.74
2:A:100:LEU:O	2:A:100:LEU:HD23	1.88	0.73
2:D:61:THR:HG23	2:D:63:ASP:H	1.54	0.72
2:A:14:ILE:O	2:A:18:ILE:HG13	1.90	0.72
2:A:197:GLU:HG2	2:A:198:PRO:HD2	1.72	0.72
2:A:94:ALA:O	2:A:98:ILE:HG13	1.90	0.71
2:A:123:GLN:OE1	2:A:124:ALA:N	2.21	0.71
2:A:189:LYS:N	2:A:189:LYS:HD3	2.05	0.71
2:A:119:LEU:HG	2:A:123:GLN:CG	2.19	0.70
1:B:19:ARG:HE	1:B:31:ILE:HD13	1.56	0.70
1:C:126:LEU:HB3	1:C:167:PHE:CE1	2.30	0.67
1:B:101:GLU:O	1:B:104:PRO:HD2	1.95	0.67
1:B:25:LYS:HA	1:B:78:PRO:HD2	1.76	0.67
1:B:53:VAL:HG22	1:B:62:ASN:HB3	1.77	0.67
1:C:65:LYS:NZ	1:C:100:ASN:ND2	2.41	0.66
2:D:170:PHE:O	2:D:174:ILE:HG13	1.94	0.66
1:B:19:ARG:NH1	1:B:192:ALA:HA	2.11	0.66
2:A:184:ASP:HB3	2:A:187:ILE:HD12	1.79	0.64
1:C:191:LYS:HD2	1:C:194:ILE:HD11	1.78	0.64
2:A:85:MET:HE2	2:A:85:MET:HA	1.80	0.63
1:B:141:GLU:HA	1:B:182:ARG:HH22	1.63	0.63
1:B:125:ASP:HA	1:B:128:GLU:HG3	1.81	0.62
1:B:115:ASP:OD2	1:B:118:MET:HG3	2.00	0.62
1:C:46:THR:HB	1:C:53:VAL:HG21	1.82	0.61
1:C:13:PRO:HG2	1:C:14:TYR:CD2	2.36	0.61
2:D:11:LYS:HD2	2:D:12:ASP:N	2.15	0.61
2:A:135:ILE:HD11	2:A:170:PHE:CE1	2.36	0.60
2:D:135:ILE:HD11	2:D:170:PHE:CE1	2.36	0.60
1:B:169:ILE:HG22	1:B:170:ASP:N	2.17	0.60
1:B:102:TRP:CH2	1:B:129:SER:HB3	2.34	0.60
1:B:118:MET:O	1:B:122:MET:HG3	2.02	0.59
1:C:103:TYR:CE1	1:C:158:LEU:HD22	2.38	0.59
2:A:132:GLN:HE22	2:D:153:ASN:ND2	1.92	0.59
1:B:25:LYS:HD2	1:B:75:PHE:O	2.03	0.59
1:B:109:ILE:O	1:B:109:ILE:HG22	2.03	0.58
1:B:166:GLY:O	1:B:168:ILE:HG12	2.03	0.58
1:C:4:VAL:HG22	1:C:29:THR:HG22	1.85	0.58
1:B:4:VAL:HG22	1:B:29:THR:HG22	1.85	0.58
1:C:14:TYR:CD1	1:C:64:ARG:HD3	2.38	0.58
2:A:120:ASP:O	2:A:122:LYS:N	2.37	0.57
2:D:83:PRO:HB2	2:D:84:PRO:HD2	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:MET:CE	1:B:74:ARG:HH22	2.16	0.57
2:D:186:ASN:O	2:D:190:GLU:HG3	2.03	0.57
2:D:197:GLU:O	2:D:201:ILE:HD12	2.05	0.57
1:C:126:LEU:HB3	1:C:167:PHE:CZ	2.39	0.57
2:D:47:LEU:HG	2:D:48:ASN:N	2.20	0.56
2:D:94:ALA:O	2:D:98:ILE:HG13	2.05	0.56
1:C:24:GLU:HG2	1:C:80:LEU:HD21	1.88	0.56
2:D:124:ALA:O	2:D:128:LYS:HG3	2.06	0.56
1:C:5:THR:HA	1:C:30:ASP:O	2.06	0.56
2:A:105:LYS:HE2	2:A:109:GLN:HE22	1.70	0.56
2:A:122:LYS:HE3	2:A:126:GLU:OE1	2.06	0.55
2:A:35:VAL:HG12	2:A:44:LEU:HD12	1.86	0.55
2:D:135:ILE:HD11	2:D:170:PHE:HE1	1.71	0.55
2:A:35:VAL:CG1	2:A:44:LEU:HD12	2.35	0.55
1:B:169:ILE:O	1:B:170:ASP:CB	2.54	0.55
1:B:18:ALA:HA	1:B:68:LEU:HD11	1.89	0.55
1:C:1:ASN:O	1:C:3:MET:N	2.39	0.55
2:A:184:ASP:CB	2:A:187:ILE:HD12	2.37	0.55
2:D:44:LEU:HG	2:D:47:LEU:HD23	1.87	0.55
2:D:187:ILE:O	2:D:191:ILE:HG13	2.07	0.55
1:C:59:TYR:CE2	1:C:61:ILE:HB	2.42	0.54
2:D:34:ASP:OD1	2:D:36:SER:HB3	2.06	0.54
2:D:7:LEU:HD12	2:D:58:THR:O	2.06	0.54
2:A:20:ARG:NH2	2:A:203:THR:O	2.37	0.54
1:B:106:LEU:HD11	1:B:110:ARG:HD3	1.90	0.54
1:B:51:PHE:HB3	1:B:52:PRO:HA	1.89	0.54
1:C:103:TYR:O	1:C:106:LEU:HB3	2.08	0.54
2:A:129:MET:O	2:A:133:ARG:HG3	2.08	0.54
1:C:50:VAL:HG21	1:C:62:ASN:OD1	2.07	0.54
2:A:34:ASP:OD1	2:A:36:SER:HB3	2.07	0.54
1:B:97:LYS:O	1:B:101:GLU:HB2	2.08	0.54
1:B:187:ASP:HA	1:B:190:LYS:CB	2.29	0.53
1:C:178:GLU:O	1:C:182:ARG:HG3	2.08	0.53
1:C:53:VAL:HG22	1:C:62:ASN:HB3	1.90	0.53
1:B:169:ILE:HG22	1:B:177:TYR:CD2	2.43	0.53
2:A:200:PHE:O	2:A:204:ILE:HG12	2.08	0.53
1:C:98:ILE:HG22	1:C:99:ASP:N	2.22	0.53
2:A:197:GLU:O	2:A:201:ILE:HD12	2.09	0.53
2:D:119:LEU:HD23	2:D:124:ALA:CB	2.35	0.52
1:C:170:ASP:HB2	1:C:172:GLU:HG2	1.90	0.52
2:A:139:TYR:O	2:A:143:VAL:HG23	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:THR:HG23	1:C:46:THR:O	2.08	0.52
1:B:108:GLN:HE21	1:B:112:HIS:CE1	2.28	0.52
1:C:86:ASN:O	1:C:90:LYS:HG2	2.10	0.52
1:B:91:ILE:O	1:B:95:LEU:HG	2.10	0.52
2:A:105:LYS:HE2	2:A:109:GLN:NE2	2.25	0.52
2:A:161:VAL:O	2:A:165:ILE:HG13	2.10	0.52
2:A:44:LEU:O	2:A:44:LEU:HD23	2.09	0.52
2:A:132:GLN:NE2	2:D:153:ASN:HD21	1.96	0.51
1:B:102:TRP:O	1:B:105:VAL:HG12	2.10	0.51
2:D:72:ILE:O	2:D:76:ILE:HG13	2.10	0.51
1:B:24:GLU:HG2	1:B:80:LEU:HD21	1.92	0.51
1:B:93:LEU:O	1:B:97:LYS:HG3	2.11	0.51
2:D:106:THR:HG22	2:D:107:PHE:CD2	2.46	0.51
2:A:107:PHE:O	2:A:111:ILE:HG13	2.10	0.51
2:A:164:LEU:HD23	2:A:164:LEU:C	2.31	0.50
2:D:106:THR:CG2	2:D:107:PHE:CE2	2.94	0.50
1:B:4:VAL:CG2	1:B:29:THR:HG22	2.41	0.49
1:C:65:LYS:HZ1	1:C:100:ASN:ND2	2.03	0.49
1:C:103:TYR:CD1	1:C:158:LEU:HD22	2.48	0.49
2:D:156:ALA:C	2:D:157:GLN:HG2	2.33	0.49
2:A:75:ALA:O	2:A:79:LEU:HG	2.12	0.49
2:A:11:LYS:HE2	1:B:83:ASN:HD21	1.78	0.49
1:B:126:LEU:HD11	1:B:167:PHE:CE2	2.48	0.48
1:C:45:ILE:HD12	1:C:55:MET:HE1	1.93	0.48
2:A:21:MET:HG2	2:A:164:LEU:HD11	1.96	0.48
1:C:25:LYS:HD2	1:C:75:PHE:O	2.13	0.48
1:C:182:ARG:O	1:C:185:ALA:HB3	2.13	0.48
2:A:123:GLN:CD	2:A:124:ALA:N	2.66	0.48
1:B:180:LYS:HG3	1:B:184:PHE:CE2	2.49	0.48
1:C:93:LEU:HD22	2:A:66:VAL:HG12	1.96	0.48
2:A:67:TYR:O	2:A:68:ARG:HB2	2.14	0.48
1:B:115:ASP:O	1:B:119:LEU:HG	2.13	0.48
1:C:50:VAL:HG21	1:C:62:ASN:CB	2.43	0.48
2:D:164:LEU:O	2:D:167:THR:HB	2.14	0.48
2:D:167:THR:OG1	2:D:194:LEU:CD1	2.62	0.48
2:A:104:ASN:HA	2:A:108:LEU:HB2	1.96	0.48
1:B:3:MET:HE2	1:B:74:ARG:HH22	1.77	0.48
2:D:156:ALA:O	2:D:157:GLN:HG2	2.13	0.48
1:B:54:LEU:HD23	1:B:54:LEU:C	2.34	0.47
1:C:126:LEU:HD12	1:C:126:LEU:HA	1.78	0.47
2:D:175:LYS:HE3	2:D:175:LYS:HB2	1.74	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:123:GLN:CG	2:A:124:ALA:N	2.77	0.47
1:B:131:LEU:HD21	1:B:173:TYR:CE1	2.49	0.47
2:D:20:ARG:HG2	2:D:32:ILE:HD13	1.95	0.47
1:B:85:VAL:HG13	2:D:78:ASP:HB3	1.95	0.47
1:B:103:TYR:N	1:B:104:PRO:CD	2.77	0.47
1:C:184:PHE:O	1:C:190:LYS:HE2	2.13	0.47
1:B:93:LEU:HD22	2:D:66:VAL:HA	1.95	0.47
1:C:26:LYS:HE3	1:C:77:ALA:O	2.14	0.47
2:D:26:LYS:NZ	2:D:83:PRO:O	2.46	0.47
2:A:119:LEU:C	2:A:123:GLN:HE21	2.18	0.47
1:B:149:THR:HG22	1:B:150:LEU:N	2.29	0.47
2:A:186:ASN:HA	2:A:189:LYS:CG	2.28	0.47
2:A:125:ASN:O	2:A:129:MET:HG3	2.15	0.47
1:C:85:VAL:HG13	2:A:78:ASP:HB3	1.96	0.47
2:D:112:ILE:HD13	2:D:112:ILE:HA	1.77	0.47
2:A:61:THR:HG23	2:A:63:ASP:N	2.22	0.47
2:A:130:LEU:C	2:A:130:LEU:HD23	2.35	0.47
1:B:25:LYS:NZ	1:B:72:ASP:HA	2.30	0.47
1:C:67:LEU:O	1:C:70:TYR:HB3	2.15	0.47
1:C:194:ILE:O	1:C:194:ILE:HD12	2.16	0.46
1:B:95:LEU:O	1:B:98:ILE:HB	2.15	0.46
2:D:138:THR:O	2:D:141:LYS:N	2.48	0.46
1:B:177:TYR:O	1:B:180:LYS:HB3	2.16	0.46
2:D:189:LYS:O	2:D:189:LYS:HD2	2.15	0.46
2:D:197:GLU:HG3	2:D:198:PRO:HD2	1.98	0.46
1:B:34:ALA:C	1:B:36:ASP:H	2.18	0.46
2:D:86:PHE:HB3	2:D:87:PRO:CD	2.46	0.46
2:D:104:ASN:HA	2:D:108:LEU:HD12	1.98	0.46
2:D:153:ASN:CG	2:D:156:ALA:HB2	2.36	0.46
1:B:36:ASP:O	1:B:36:ASP:OD1	2.34	0.46
1:C:27:MET:HA	1:C:27:MET:CE	2.45	0.46
1:B:72:ASP:OD1	1:B:80:LEU:HB2	2.17	0.46
2:D:180:ILE:HA	2:D:181:PRO:HD2	1.83	0.46
2:D:128:LYS:O	2:D:132:GLN:HG3	2.16	0.45
2:D:77:GLU:OE1	2:D:77:GLU:HA	2.16	0.45
2:A:100:LEU:C	2:A:100:LEU:HD23	2.35	0.45
1:C:8:THR:HB	1:C:15:SER:HB3	1.97	0.45
2:D:93:ARG:O	2:D:96:ALA:HB3	2.17	0.45
1:B:68:LEU:HD12	1:B:71:ILE:HD12	1.98	0.45
1:C:129:SER:O	1:C:133:MET:HG2	2.17	0.45
2:A:89:PHE:N	2:A:89:PHE:CD1	2.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:151:ALA:O	1:C:155:ILE:HG13	2.17	0.45
2:D:87:PRO:HD2	2:D:93:ARG:HG2	1.98	0.45
1:B:169:ILE:HG22	1:B:177:TYR:HD2	1.81	0.44
2:D:141:LYS:O	2:D:144:SER:HB3	2.17	0.44
1:B:89:ILE:CD1	2:D:79:LEU:HD21	2.47	0.44
2:A:119:LEU:C	2:A:123:GLN:NE2	2.70	0.44
2:A:21:MET:O	2:A:21:MET:CG	2.65	0.44
1:C:18:ALA:HA	1:C:68:LEU:HD11	1.99	0.44
2:D:44:LEU:HG	2:D:47:LEU:CD2	2.47	0.44
1:B:105:VAL:O	1:B:109:ILE:HG13	2.17	0.44
1:B:106:LEU:O	1:B:106:LEU:HD12	2.17	0.44
2:D:75:ALA:O	2:D:79:LEU:HG	2.16	0.44
1:C:9:THR:HG23	1:C:12:CYS:CB	2.48	0.44
2:A:84:PRO:HG2	2:A:158:ASN:HA	1.99	0.44
1:B:149:THR:HG22	1:B:150:LEU:H	1.82	0.44
2:D:171:TYR:CD1	2:D:200:PHE:HZ	2.36	0.44
1:B:0:SER:OG	1:B:1:ASN:N	2.49	0.44
1:C:106:LEU:O	1:C:110:ARG:HG3	2.18	0.44
2:A:6:LEU:HD12	2:A:31:LYS:O	2.18	0.44
2:A:44:LEU:HD23	2:A:44:LEU:C	2.39	0.44
1:B:108:GLN:NE2	1:B:112:HIS:CE1	2.86	0.44
1:B:86:ASN:O	1:B:90:LYS:HG3	2.18	0.44
1:C:103:TYR:N	1:C:104:PRO:CD	2.81	0.44
2:D:44:LEU:O	2:D:47:LEU:HD23	2.18	0.44
2:D:46:GLU:HG3	2:D:46:GLU:H	1.58	0.44
2:A:104:ASN:O	2:A:109:GLN:HG2	2.18	0.44
2:A:10:LYS:HG3	2:A:56:ILE:HD12	2.00	0.44
2:A:123:GLN:N	2:A:123:GLN:OE1	2.51	0.44
1:B:4:VAL:HG12	1:B:56:GLU:HG2	1.99	0.44
2:A:61:THR:HG21	2:A:64:PHE:CE2	2.52	0.43
1:B:167:PHE:CZ	1:B:173:TYR:OH	2.70	0.43
1:B:33:GLU:O	1:B:36:ASP:HB3	2.18	0.43
1:C:113:ARG:C	1:C:115:ASP:H	2.20	0.43
2:A:127:ILE:HG22	2:A:131:MET:HE3	2.00	0.43
2:D:11:LYS:HZ3	2:D:12:ASP:CG	2.22	0.43
2:A:170:PHE:O	2:A:174:ILE:HG13	2.18	0.43
2:D:168:PHE:O	2:D:172:TYR:HD1	2.01	0.43
2:A:43:HIS:O	2:A:46:GLU:HB3	2.18	0.43
1:C:166:GLY:O	1:C:168:ILE:HD12	2.19	0.43
1:C:25:LYS:O	1:C:26:LYS:C	2.57	0.43
2:A:124:ALA:O	2:A:125:ASN:C	2.56	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:82:PHE:HA	2:A:83:PRO:HA	1.82	0.43
1:B:109:ILE:O	1:B:113:ARG:HB3	2.19	0.43
1:C:26:LYS:HE2	1:C:78:PRO:HD3	2.00	0.43
2:D:106:THR:HG22	2:D:107:PHE:CE2	2.54	0.43
2:D:197:GLU:HG3	2:D:198:PRO:CD	2.49	0.43
2:D:82:PHE:HA	2:D:83:PRO:HA	1.88	0.43
2:D:111:ILE:O	2:D:114:LEU:HB2	2.19	0.43
2:D:168:PHE:HA	2:D:171:TYR:HB3	2.01	0.42
1:B:27:MET:O	1:B:29:THR:HG23	2.19	0.42
1:C:105:VAL:O	1:C:109:ILE:HG13	2.18	0.42
2:A:40:ASN:C	2:A:42:LYS:N	2.71	0.42
1:B:169:ILE:HG13	1:B:169:ILE:H	1.52	0.42
2:D:61:THR:C	2:D:63:ASP:H	2.22	0.42
2:A:183:LYS:O	2:A:184:ASP:C	2.57	0.42
1:C:38:GLU:C	1:C:40:ALA:H	2.22	0.42
2:D:89:PHE:HA	2:D:90:PRO:HD3	1.79	0.42
1:C:163:GLU:HA	1:C:167:PHE:O	2.20	0.42
1:C:3:MET:CE	1:C:27:MET:HE2	2.49	0.42
1:C:48:ASN:HD22	1:C:50:VAL:CG1	2.32	0.42
1:B:106:LEU:HD13	1:B:162:LEU:HD13	2.02	0.42
1:C:110:ARG:O	1:C:113:ARG:HG2	2.20	0.42
1:C:180:LYS:O	1:C:184:PHE:HD2	2.02	0.42
1:C:14:TYR:CE1	1:C:64:ARG:HD3	2.54	0.42
2:A:85:MET:HE2	2:A:159:ILE:HB	2.02	0.42
1:B:180:LYS:HB2	1:B:180:LYS:HE3	1.77	0.42
2:D:91:LYS:HG3	2:D:92:GLN:N	2.35	0.42
2:A:164:LEU:HD23	2:A:165:ILE:N	2.35	0.42
1:B:116:GLN:O	1:B:120:GLU:HG2	2.20	0.42
1:C:110:ARG:HB3	1:C:113:ARG:NH1	2.35	0.42
1:C:108:GLN:O	1:C:122:MET:CE	2.68	0.42
2:D:44:LEU:HD23	2:D:44:LEU:C	2.40	0.42
2:D:144:SER:O	2:D:148:VAL:HG23	2.19	0.41
1:C:80:LEU:HA	1:C:80:LEU:HD23	1.83	0.41
2:D:39:GLU:H	2:D:39:GLU:HG2	1.65	0.41
1:B:67:LEU:O	1:B:70:TYR:HB3	2.20	0.41
1:C:9:THR:HG22	1:C:51:PHE:HB3	2.03	0.41
2:D:107:PHE:O	2:D:111:ILE:HG13	2.21	0.41
2:D:104:ASN:O	2:D:109:GLN:HG2	2.20	0.41
1:B:194:ILE:HD12	1:B:194:ILE:C	2.41	0.41
1:C:109:ILE:HG22	1:C:109:ILE:O	2.21	0.41
1:B:19:ARG:HH12	1:B:192:ALA:HA	1.80	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:143:VAL:O	2:A:144:SER:C	2.59	0.41
1:B:7:TYR:CD1	1:B:32:VAL:HB	2.55	0.41
1:B:97:LYS:O	1:B:101:GLU:HG3	2.21	0.41
1:B:106:LEU:C	1:B:106:LEU:HD12	2.40	0.41
1:C:25:LYS:C	1:C:26:LYS:HD2	2.39	0.41
1:C:84:VAL:O	1:C:85:VAL:C	2.59	0.41
2:D:135:ILE:HG23	2:D:136:ILE:H	1.86	0.41
2:D:171:TYR:CG	2:D:200:PHE:HZ	2.39	0.41
2:A:89:PHE:HD1	2:A:89:PHE:N	2.19	0.41
2:D:161:VAL:CG1	2:D:162:LEU:N	2.84	0.41
2:A:132:GLN:O	2:A:135:ILE:HG22	2.21	0.40
2:A:89:PHE:HA	2:A:90:PRO:HD3	1.95	0.40
1:C:34:ALA:C	1:C:36:ASP:H	2.24	0.40
2:D:10:LYS:HD2	2:D:56:ILE:HG13	2.03	0.40
1:C:56:GLU:O	1:C:57:LYS:C	2.59	0.40
2:D:176:LEU:O	2:D:177:LYS:HB2	2.20	0.40
2:A:159:ILE:HG22	2:A:160:ASN:N	2.37	0.40
1:B:13:PRO:HB2	1:B:103:TYR:CE2	2.56	0.40
1:C:109:ILE:O	1:C:113:ARG:HB3	2.21	0.40
2:D:161:VAL:HG13	2:D:162:LEU:N	2.35	0.40
2:D:167:THR:OG1	2:D:194:LEU:HD13	2.21	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	B	193/211 (92%)	174 (90%)	15 (8%)	4 (2%)	8	38
1	C	193/211 (92%)	174 (90%)	14 (7%)	5 (3%)	6	33
2	A	201/204 (98%)	184 (92%)	13 (6%)	4 (2%)	9	39
2	D	201/204 (98%)	178 (89%)	21 (10%)	2 (1%)	18	53

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	788/830 (95%)	710 (90%)	63 (8%)	15 (2%)	9	41

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	2	ALA
1	C	52	PRO
2	A	121	GLU
1	B	170	ASP
1	B	171	ASP
1	B	169	ILE
1	C	57	LYS
2	A	185	LYS
1	B	2	ALA
2	D	184	ASP
2	A	87	PRO
2	D	181	PRO
1	C	114	SER
1	C	85	VAL
2	A	148	VAL

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	B	160/184 (87%)	146 (91%)	14 (9%)	12	40
1	C	161/184 (88%)	145 (90%)	16 (10%)	9	34
2	A	176/192 (92%)	170 (97%)	6 (3%)	42	73
2	D	174/192 (91%)	170 (98%)	4 (2%)	56	79
All	All	671/752 (89%)	631 (94%)	40 (6%)	22	58

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	9	THR
1	C	26	LYS
1	C	41	MET
1	C	48	ASN
1	C	50	VAL
1	C	51	PHE
1	C	56	GLU
1	C	91	ILE
1	C	114	SER
1	C	128	GLU
1	C	130	LEU
1	C	149	THR
1	C	163	GLU
1	C	171	ASP
1	C	187	ASP
1	C	188	SER
2	A	56	ILE
2	A	68	ARG
2	A	91	LYS
2	A	123	GLN
2	A	162	LEU
2	A	166	ILE
1	B	9	THR
1	B	19	ARG
1	B	68	LEU
1	B	96	ASP
1	B	106	LEU
1	B	107	ASP
1	B	108	GLN
1	B	114	SER
1	B	126	LEU
1	B	131	LEU
1	B	137	PHE
1	B	162	LEU
1	B	171	ASP
1	B	180	LYS
2	D	47	LEU
2	D	68	ARG
2	D	146	ARG
2	D	197	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	100	ASN
1	C	112	HIS
2	A	109	GLN
2	A	132	GLN
1	B	108	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	B	195/211 (92%)	0.32	11 (5%) 25 23	59, 110, 191, 240	0
1	C	195/211 (92%)	-0.04	3 (1%) 74 70	49, 90, 162, 194	0
2	A	203/204 (99%)	-0.16	2 (0%) 82 81	47, 79, 132, 189	0
2	D	203/204 (99%)	-0.10	3 (1%) 74 70	50, 80, 137, 181	0
All	All	796/830 (95%)	0.00	19 (2%) 59 56	47, 89, 170, 240	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	2	SER	8.0
1	B	167	PHE	5.6
1	B	169	ILE	4.3
1	B	131	LEU	3.5
2	D	3	ASN	3.3
1	B	130	LEU	3.1
1	C	121	SER	2.7
1	B	166	GLY	2.6
1	C	112	HIS	2.6
1	B	117	LYS	2.5
1	B	102	TRP	2.4
1	C	0	SER	2.3
1	B	155	ILE	2.3
1	B	32	VAL	2.3
1	B	112	HIS	2.2
2	D	56	ILE	2.2
2	D	8	TYR	2.2
2	A	3	ASN	2.1
1	B	34	ALA	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.