Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 06:53 PM EDT

PDB ID	:	3M1Y
Title	:	Crystal Structure of a Phosphoserine phosphatase (SerB) from Helicobacter pylori
Authors	:	Syed Ibrahim, B.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2010-03-05
Resolution	:	2.40 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

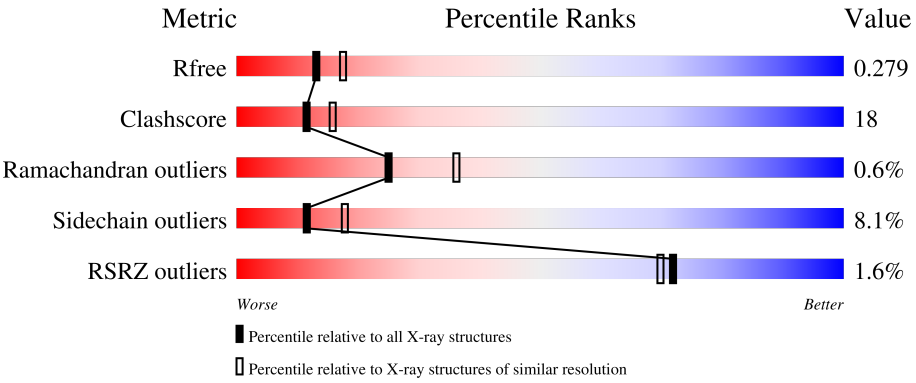
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=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	217	65% 26% • 6%
1	B	217	% 61% 31% • 5%
1	C	217	% 55% 36% 5% •
1	D	217	3% 51% 40% • 5%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6500 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 Phosphoserine phosphatase (SerB).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	Se	0	0	0
			1580	1015	267	290	3	5			
1	B	206	Total	C	N	O	S	Se	0	0	0
			1599	1026	270	294	3	6			
1	C	208	Total	C	N	O	S	Se	0	0	0
			1611	1035	273	294	3	6			
1	D	206	Total	C	N	O	S	Se	0	0	0
			1598	1024	271	294	3	6			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	expression tag	UNP O25367
A	0	SER	-	expression tag	UNP O25367
A	1	LEU	-	expression tag	UNP O25367
A	208	GLU	-	expression tag	UNP O25367
A	209	GLY	-	expression tag	UNP O25367
A	210	HIS	-	expression tag	UNP O25367
A	211	HIS	-	expression tag	UNP O25367
A	212	HIS	-	expression tag	UNP O25367
A	213	HIS	-	expression tag	UNP O25367
A	214	HIS	-	expression tag	UNP O25367
A	215	HIS	-	expression tag	UNP O25367
B	-1	MSE	-	expression tag	UNP O25367
B	0	SER	-	expression tag	UNP O25367
B	1	LEU	-	expression tag	UNP O25367
B	208	GLU	-	expression tag	UNP O25367
B	209	GLY	-	expression tag	UNP O25367
B	210	HIS	-	expression tag	UNP O25367
B	211	HIS	-	expression tag	UNP O25367
B	212	HIS	-	expression tag	UNP O25367
B	213	HIS	-	expression tag	UNP O25367
B	214	HIS	-	expression tag	UNP O25367

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Chain	Residue	Modelled	Actual	Comment	Reference
B	215	HIS	-	expression tag	UNP O25367
C	-1	MSE	-	expression tag	UNP O25367
C	0	SER	-	expression tag	UNP O25367
C	1	LEU	-	expression tag	UNP O25367
C	208	GLU	-	expression tag	UNP O25367
C	209	GLY	-	expression tag	UNP O25367
C	210	HIS	-	expression tag	UNP O25367
C	211	HIS	-	expression tag	UNP O25367
C	212	HIS	-	expression tag	UNP O25367
C	213	HIS	-	expression tag	UNP O25367
C	214	HIS	-	expression tag	UNP O25367
C	215	HIS	-	expression tag	UNP O25367
D	-1	MSE	-	expression tag	UNP O25367
D	0	SER	-	expression tag	UNP O25367
D	1	LEU	-	expression tag	UNP O25367
D	208	GLU	-	expression tag	UNP O25367
D	209	GLY	-	expression tag	UNP O25367
D	210	HIS	-	expression tag	UNP O25367
D	211	HIS	-	expression tag	UNP O25367
D	212	HIS	-	expression tag	UNP O25367
D	213	HIS	-	expression tag	UNP O25367
D	214	HIS	-	expression tag	UNP O25367
D	215	HIS	-	expression tag	UNP O25367

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Cl 2 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	C	1	Total Cl 1 1	0	0
3	D	1	Total Cl 1 1	0	0

- Molecule 4 is water.

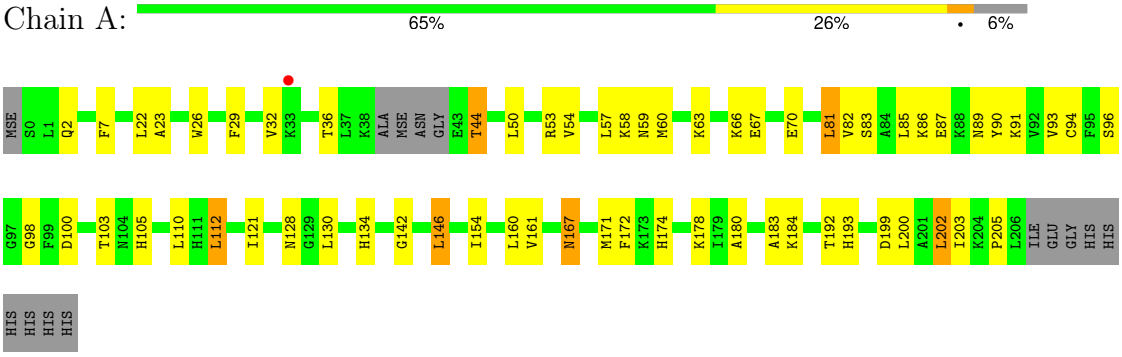
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	30	Total O 30 30	0	0
4	B	24	Total O 24 24	0	0
4	C	35	Total O 35 35	0	0
4	D	13	Total O 13 13	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

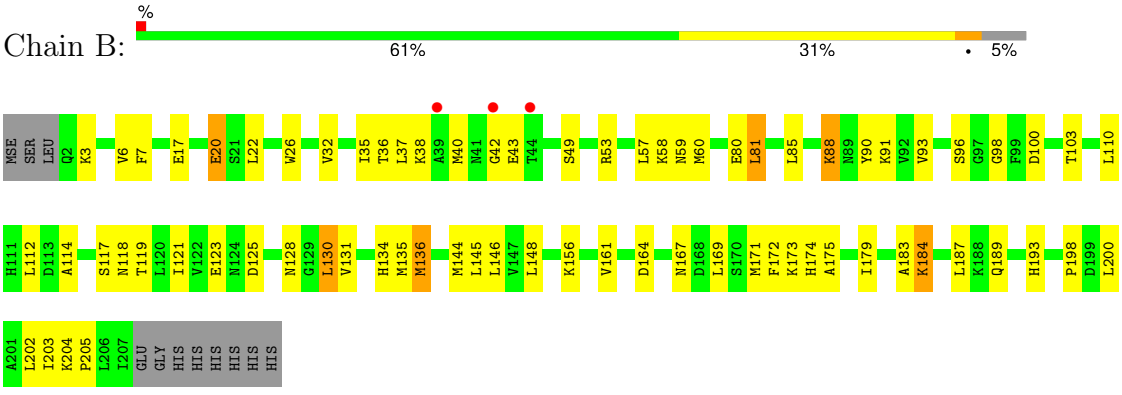
• Molecule 1: Phosphoserine phosphatase (SerB)

Chain A:



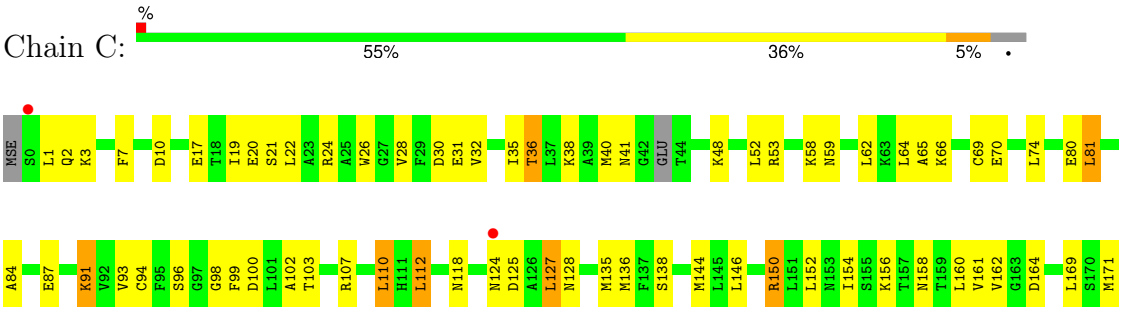
• Molecule 1: Phosphoserine phosphatase (SerB)

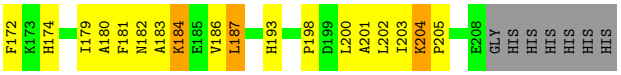
Chain B:



• Molecule 1: Phosphoserine phosphatase (SerB)

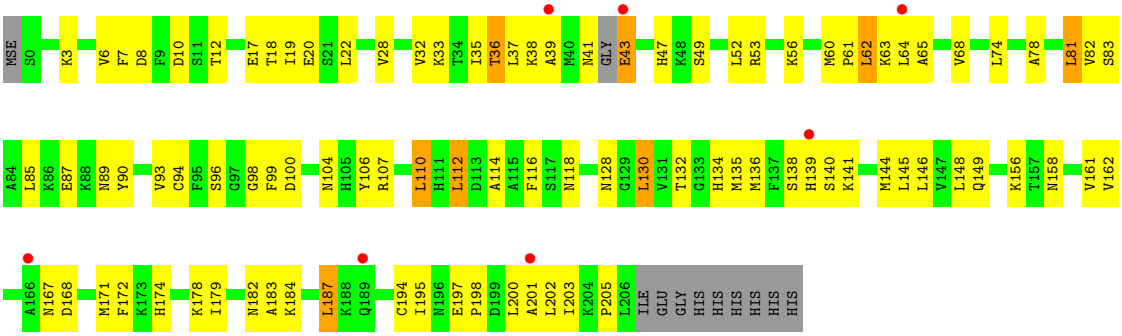
Chain C:





• Molecule 1: Phosphoserine phosphatase (SerB)

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	51.61Å 84.72Å 204.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.07 – 2.40 44.07 – 2.39	Depositor EDS
% Data completeness (in resolution range)	94.0 (44.07-2.40) 93.4 (44.07-2.39)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.60 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.229 , 0.271 0.246 , 0.279	Depositor DCC
R_{free} test set	1673 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å^2)	34.7	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 39.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	6500	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: MG, CL

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	0.52	0/1600	0.71	2/2150 (0.1%)
1	B	0.48	1/1620 (0.1%)	0.64	0/2177
1	C	0.51	0/1631	0.72	1/2190 (0.0%)
1	D	0.49	0/1617	0.68	1/2170 (0.0%)
All	All	0.50	1/6468 (0.0%)	0.69	4/8687 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	136	MSE	CG-SE	-5.69	1.76	1.95

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	205	PRO	N-CA-CB	5.97	110.47	103.30
1	C	91	LYS	N-CA-C	-5.58	95.94	111.00
1	A	91	LYS	N-CA-C	-5.49	96.18	111.00
1	A	205	PRO	N-CA-CB	5.14	109.46	103.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1580	0	1624	45	0
1	B	1599	0	1637	49	0
1	C	1611	0	1655	69	0
1	D	1598	0	1630	72	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	30	0	0	0	0
4	B	24	0	0	1	0
4	C	35	0	0	1	0
4	D	13	0	0	1	0
All	All	6500	0	6546	228	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 18.

All (228) 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:A:167:ASN:ND2	1:A:167:ASN:H	1.66	0.89
1:A:161:VAL:HG13	1:A:171:MSE:HE3	1.53	0.87
1:C:17:GLU:HB3	1:C:20:GLU:HG2	1.59	0.85
1:D:161:VAL:HG21	1:D:171:MSE:HB3	1.58	0.82
1:C:38:LYS:HA	1:C:41:ASN:ND2	1.97	0.80
1:D:110:LEU:HB3	1:D:112:LEU:HD13	1.64	0.78
1:C:107:ARG:HA	1:C:112:LEU:HD22	1.67	0.76
1:A:167:ASN:H	1:A:167:ASN:HD22	1.36	0.73
1:D:81:LEU:HD12	1:D:200:LEU:HB3	1.70	0.73
1:D:83:SER:O	1:D:87:GLU:HG2	1.89	0.73
1:B:145:LEU:HD23	1:B:174:HIS:HB2	1.71	0.72
1:C:36:THR:HG22	1:C:53:ARG:HD3	1.71	0.72
1:A:105:HIS:NE2	1:C:193:HIS:HD2	1.86	0.72
1:A:128:ASN:HD21	1:A:130:LEU:HD22	1.55	0.72
1:C:10:ASP:HB2	4:C:317:HOH:O	1.90	0.71
1:B:184:LYS:N	1:B:184:LYS:HE2	2.06	0.71
1:D:38:LYS:O	1:D:43:GLU:HB3	1.92	0.69
1:B:200:LEU:O	1:B:203:ILE:HD13	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:VAL:HG13	1:B:53:ARG:HG2	1.76	0.68
1:B:37:LEU:O	1:B:38:LYS:HB2	1.93	0.68
1:D:32:VAL:HG13	1:D:53:ARG:HG2	1.76	0.68
1:D:61:PRO:HG2	1:D:64:LEU:HD13	1.76	0.68
1:C:150:ARG:CG	1:C:150:ARG:HH11	2.07	0.67
1:D:182:ASN:HD21	1:D:198:PRO:HA	1.60	0.67
1:A:199:ASP:HB3	1:A:202:LEU:HD22	1.76	0.67
1:A:121:ILE:HD12	1:A:130:LEU:HB3	1.77	0.67
1:C:64:LEU:O	1:C:65:ALA:HB3	1.95	0.66
1:D:49:SER:O	1:D:53:ARG:HD3	1.96	0.66
1:C:31:GLU:O	1:C:35:ILE:HG13	1.95	0.65
1:D:35:ILE:HD13	1:D:52:LEU:HB3	1.79	0.65
1:C:66:LYS:O	1:C:70:GLU:HG3	1.98	0.64
1:B:164:ASP:HA	1:B:183:ALA:HB2	1.80	0.63
1:B:100:ASP:HA	1:B:103:THR:OG1	1.98	0.61
1:D:85:LEU:O	1:D:90:TYR:HB2	1.99	0.61
1:A:66:LYS:O	1:A:70:GLU:HG3	2.00	0.60
1:D:10:ASP:HB2	4:D:328:HOH:O	2.01	0.60
1:C:135:MSE:HE1	1:C:144:MSE:HB2	1.84	0.60
1:D:172:PHE:O	1:D:178:LYS:HD3	2.02	0.60
1:D:162:VAL:HG22	1:D:179:ILE:HB	1.84	0.60
1:A:83:SER:O	1:A:87:GLU:HG3	2.03	0.59
1:A:100:ASP:HA	1:A:103:THR:OG1	2.01	0.59
1:A:134:HIS:HE2	1:B:144:MSE:HE1	1.67	0.59
1:C:172:PHE:CE2	1:C:187:LEU:HD22	2.37	0.59
1:C:164:ASP:O	1:C:184:LYS:HD2	2.03	0.58
1:A:167:ASN:ND2	1:A:167:ASN:N	2.45	0.58
1:C:48:LYS:O	1:C:52:LEU:HG	2.03	0.58
1:D:145:LEU:O	1:D:149:GLN:HG3	2.03	0.58
1:A:110:LEU:HB2	1:A:112:LEU:HD13	1.85	0.58
1:C:36:THR:O	1:C:40:MSE:HG3	2.03	0.58
1:A:63:LYS:NZ	1:A:63:LYS:HB3	2.19	0.57
1:B:6:VAL:HG13	1:B:161:VAL:HG22	1.85	0.57
1:D:32:VAL:O	1:D:36:THR:HG22	2.03	0.57
1:D:138:SER:O	1:D:139:HIS:HB2	2.03	0.57
1:B:128:ASN:OD1	1:B:130:LEU:HD22	2.05	0.57
1:B:6:VAL:CG1	1:B:161:VAL:HG22	2.35	0.57
1:A:128:ASN:ND2	1:A:130:LEU:HD22	2.19	0.56
1:B:203:ILE:N	1:B:203:ILE:HD12	2.20	0.56
1:A:32:VAL:HG13	1:A:53:ARG:HG2	1.86	0.56
1:B:32:VAL:O	1:B:36:THR:HG23	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:PHE:HZ	1:C:81:LEU:HD13	1.72	0.55
1:C:150:ARG:HH11	1:C:150:ARG:HG3	1.70	0.55
1:B:161:VAL:HG11	1:B:171:MSE:HG3	1.88	0.55
1:D:28:VAL:HG12	1:D:28:VAL:O	2.07	0.54
1:B:17:GLU:HB3	1:B:20:GLU:HG3	1.89	0.54
1:B:85:LEU:O	1:B:90:TYR:HB2	2.07	0.54
1:C:74:LEU:N	1:C:74:LEU:HD12	2.23	0.54
1:D:28:VAL:HG11	1:D:56:LYS:HB3	1.90	0.54
1:A:85:LEU:O	1:A:90:TYR:HB2	2.07	0.53
1:B:136:MSE:HB2	4:B:333:HOH:O	2.08	0.53
1:C:38:LYS:NZ	1:C:38:LYS:HB3	2.23	0.53
1:D:184:LYS:HD3	1:D:184:LYS:N	2.24	0.53
1:D:144:MSE:O	1:D:148:LEU:HD13	2.08	0.53
1:A:167:ASN:HD22	1:A:167:ASN:N	2.06	0.53
1:D:93:VAL:HG12	1:D:94:CYS:N	2.23	0.53
1:D:197:GLU:OE1	1:D:197:GLU:HA	2.08	0.53
1:B:81:LEU:HD22	1:B:85:LEU:HG	1.91	0.53
1:C:204:LYS:N	1:C:205:PRO:HD2	2.24	0.53
1:B:135:MSE:HE1	1:B:144:MSE:HB2	1.91	0.53
1:A:184:LYS:HD2	1:A:184:LYS:N	2.23	0.53
1:B:7:PHE:HZ	1:B:81:LEU:HD13	1.73	0.53
1:D:106:TYR:O	1:D:110:LEU:HB2	2.09	0.53
1:D:19:ILE:HG23	1:D:20:GLU:N	2.25	0.52
1:A:161:VAL:HG21	1:A:171:MSE:HG2	1.91	0.52
1:A:192:THR:HG22	1:A:193:HIS:CE1	2.44	0.52
1:C:36:THR:HG22	1:C:53:ARG:CD	2.38	0.52
1:D:6:VAL:HG13	1:D:161:VAL:HG23	1.91	0.51
1:C:26:TRP:HA	1:C:26:TRP:CE3	2.45	0.51
1:C:93:VAL:HG12	1:C:94:CYS:N	2.25	0.51
1:B:26:TRP:CH2	1:B:60:MSE:HA	2.45	0.51
1:A:50:LEU:O	1:A:54:VAL:HG23	2.11	0.51
1:D:172:PHE:CE2	1:D:187:LEU:HD22	2.46	0.51
1:C:96:SER:C	1:C:98:GLY:H	2.14	0.51
1:D:141:LYS:NZ	1:D:167:ASN:ND2	2.59	0.51
1:B:179:ILE:HD12	1:B:193:HIS:HB2	1.94	0.50
1:C:7:PHE:CZ	1:C:81:LEU:HD13	2.46	0.50
1:C:2:GLN:HG2	1:C:154:ILE:HD11	1.94	0.50
1:C:180:ALA:HB1	1:C:183:ALA:HB3	1.92	0.50
1:C:17:GLU:HB3	1:C:20:GLU:CG	2.37	0.50
1:B:98:GLY:O	1:B:118:ASN:HB2	2.11	0.50
1:C:100:ASP:HA	1:C:103:THR:OG1	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:SER:HB3	1:D:53:ARG:HH11	1.77	0.50
1:D:128:ASN:OD1	1:D:130:LEU:HB2	2.12	0.50
1:C:69:CYS:HB3	1:C:102:ALA:HB2	1.94	0.49
1:C:150:ARG:CG	1:C:150:ARG:NH1	2.69	0.49
1:A:142:GLY:O	1:A:146:LEU:HD22	2.13	0.49
1:A:82:VAL:O	1:A:86:LYS:HG3	2.13	0.49
1:A:58:LYS:O	1:A:59:ASN:HB2	2.13	0.49
1:B:161:VAL:HG13	1:B:171:MSE:HE3	1.94	0.49
1:C:201:ALA:O	1:C:204:LYS:HG2	2.12	0.49
1:B:134:HIS:N	1:B:134:HIS:CD2	2.81	0.48
1:C:110:LEU:HB3	1:C:112:LEU:HD13	1.94	0.48
1:D:18:THR:O	1:D:22:LEU:HB2	2.13	0.48
1:D:184:LYS:N	1:D:184:LYS:CD	2.75	0.48
1:A:2:GLN:HG3	1:A:154:ILE:HD11	1.94	0.48
1:A:81:LEU:HD12	1:A:200:LEU:HB3	1.95	0.48
1:D:135:MSE:HE2	1:D:140:SER:HB3	1.95	0.48
1:A:7:PHE:CZ	1:A:81:LEU:HD13	2.49	0.48
1:A:161:VAL:CG2	1:A:171:MSE:HG2	2.44	0.48
1:C:162:VAL:HG22	1:C:179:ILE:HB	1.95	0.48
1:C:32:VAL:O	1:C:36:THR:HG23	2.13	0.48
1:B:156:LYS:HG3	1:B:174:HIS:O	2.13	0.48
1:D:141:LYS:NZ	1:D:167:ASN:HD22	2.12	0.48
1:A:172:PHE:O	1:A:178:LYS:HE2	2.14	0.47
1:B:35:ILE:O	1:B:37:LEU:O	2.32	0.47
1:B:119:THR:O	1:B:131:VAL:HG23	2.14	0.47
1:C:160:LEU:C	1:C:160:LEU:HD23	2.35	0.47
1:A:134:HIS:NE2	1:B:144:MSE:HE1	2.30	0.47
1:B:36:THR:O	1:B:40:MSE:HG3	2.14	0.47
1:B:169:LEU:O	1:B:173:LYS:HG3	2.13	0.47
1:C:144:MSE:HE1	1:D:134:HIS:HE1	1.80	0.47
1:D:8:ASP:O	1:D:12:THR:HB	2.15	0.47
1:C:144:MSE:HE1	1:D:134:HIS:CE1	2.50	0.47
1:B:17:GLU:HB3	1:B:20:GLU:CG	2.45	0.46
1:D:96:SER:C	1:D:98:GLY:H	2.17	0.46
1:C:3:LYS:HG2	1:C:158:ASN:HB3	1.97	0.46
1:D:3:LYS:HG2	1:D:158:ASN:HB3	1.96	0.46
1:D:81:LEU:HD11	1:D:203:ILE:HD12	1.97	0.46
1:D:62:LEU:HD12	1:D:63:LYS:N	2.31	0.46
1:D:107:ARG:HA	1:D:112:LEU:HD22	1.97	0.46
1:C:62:LEU:O	1:C:64:LEU:O	2.34	0.46
1:C:184:LYS:HD2	1:C:184:LYS:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:TRP:HA	1:C:26:TRP:HE3	1.80	0.46
1:D:78:ALA:O	1:D:82:VAL:HG23	2.16	0.45
1:D:161:VAL:HG21	1:D:171:MSE:CB	2.38	0.45
1:D:168:ASP:HB2	1:D:187:LEU:HD11	1.98	0.45
1:C:91:LYS:HD3	1:C:91:LYS:HA	1.80	0.45
1:B:110:LEU:HB2	1:B:112:LEU:HD13	1.97	0.45
1:C:99:PHE:HA	1:C:118:ASN:O	2.17	0.45
1:A:26:TRP:CH2	1:A:60:MSE:HA	2.52	0.45
1:D:200:LEU:O	1:D:201:ALA:HB3	2.16	0.45
1:A:105:HIS:NE2	1:C:193:HIS:CD2	2.77	0.44
1:C:172:PHE:CZ	1:C:187:LEU:HD22	2.52	0.44
1:D:33:LYS:O	1:D:36:THR:HG23	2.18	0.44
1:B:7:PHE:CZ	1:B:81:LEU:HD13	2.52	0.44
1:B:49:SER:O	1:B:53:ARG:HD3	2.17	0.44
1:B:96:SER:C	1:B:98:GLY:H	2.20	0.44
1:D:60:MSE:HE2	1:D:65:ALA:N	2.32	0.44
1:D:182:ASN:ND2	1:D:198:PRO:HA	2.31	0.44
1:A:93:VAL:HG12	1:A:94:CYS:N	2.32	0.44
1:A:161:VAL:HG13	1:A:171:MSE:CE	2.37	0.44
1:A:180:ALA:HB1	1:A:183:ALA:HB3	1.99	0.44
1:B:134:HIS:CD2	1:B:134:HIS:H	2.35	0.44
1:D:203:ILE:HG22	1:D:203:ILE:O	2.18	0.44
1:C:21:SER:HA	1:C:24:ARG:NH1	2.32	0.44
1:C:181:PHE:O	1:C:182:ASN:C	2.56	0.44
1:D:64:LEU:O	1:D:68:VAL:HG23	2.18	0.44
1:C:38:LYS:HA	1:C:41:ASN:HD21	1.81	0.43
1:B:93:VAL:HG22	1:B:114:ALA:HB3	2.00	0.43
1:B:58:LYS:O	1:B:59:ASN:HB2	2.17	0.43
1:D:138:SER:O	1:D:139:HIS:CB	2.67	0.43
1:D:156:LYS:HG3	1:D:174:HIS:C	2.39	0.43
1:D:183:ALA:C	1:D:184:LYS:HD3	2.39	0.43
1:C:58:LYS:O	1:C:59:ASN:HB2	2.18	0.43
1:C:96:SER:C	1:C:98:GLY:N	2.70	0.43
1:A:54:VAL:O	1:A:57:LEU:HB2	2.19	0.43
1:C:81:LEU:HD12	1:C:200:LEU:HB3	2.00	0.43
1:D:17:GLU:HB3	1:D:20:GLU:CG	2.48	0.43
1:D:134:HIS:N	1:D:134:HIS:CD2	2.85	0.42
1:D:99:PHE:HA	1:D:118:ASN:O	2.19	0.42
1:C:58:LYS:HG3	1:C:128:ASN:O	2.18	0.42
1:C:127:LEU:HD12	1:C:127:LEU:HA	1.77	0.42
1:C:180:ALA:HB1	1:C:183:ALA:CB	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:GLU:HG2	1:C:204:LYS:HE3	2.02	0.42
1:D:33:LYS:HA	1:D:36:THR:CG2	2.49	0.42
1:B:156:LYS:HE3	1:B:175:ALA:N	2.35	0.42
1:D:74:LEU:HD12	1:D:74:LEU:N	2.35	0.42
1:D:194:CYS:O	1:D:195:ILE:HD13	2.19	0.42
1:C:19:ILE:HG23	1:C:20:GLU:N	2.34	0.42
1:C:28:VAL:HG12	1:C:28:VAL:O	2.19	0.42
1:A:146:LEU:HD21	1:A:174:HIS:CE1	2.55	0.42
1:C:136:MSE:HG3	1:D:132:THR:OG1	2.20	0.42
1:D:110:LEU:HB3	1:D:112:LEU:CD1	2.42	0.41
1:D:47:HIS:HD2	1:D:136:MSE:O	2.03	0.41
1:D:200:LEU:C	1:D:202:LEU:N	2.73	0.41
1:B:88:LYS:HD2	1:B:88:LYS:N	2.36	0.41
1:C:161:VAL:CG2	1:C:171:MSE:HE3	2.51	0.41
1:C:169:LEU:HD11	1:C:186:VAL:HG12	2.01	0.41
1:D:93:VAL:HG13	1:D:114:ALA:O	2.21	0.41
1:A:23:ALA:HB1	1:A:29:PHE:HA	2.02	0.41
1:C:64:LEU:O	1:C:65:ALA:CB	2.60	0.41
1:B:37:LEU:O	1:B:38:LYS:CB	2.64	0.41
1:C:203:ILE:C	1:C:205:PRO:HD2	2.41	0.41
1:B:37:LEU:HD23	1:B:40:MSE:HE2	2.02	0.41
1:A:93:VAL:CG1	1:A:94:CYS:N	2.84	0.41
1:B:3:LYS:HB3	1:B:90:TYR:CE1	2.55	0.41
1:B:38:LYS:HB3	1:B:43:GLU:HB2	2.02	0.41
1:C:64:LEU:O	1:C:66:LYS:N	2.47	0.41
1:A:160:LEU:C	1:A:160:LEU:HD23	2.41	0.41
1:B:121:ILE:HD12	1:B:130:LEU:HB3	2.02	0.41
1:B:204:LYS:HB2	1:B:205:PRO:HD3	2.03	0.41
1:C:36:THR:HG22	1:C:53:ARG:HH11	1.85	0.41
1:C:41:ASN:OD1	1:C:41:ASN:C	2.60	0.41
1:A:192:THR:HG22	1:A:193:HIS:ND1	2.36	0.41
1:C:156:LYS:HG3	1:C:174:HIS:O	2.20	0.41
1:A:96:SER:C	1:A:98:GLY:H	2.23	0.40
1:C:84:ALA:O	1:C:87:GLU:HG2	2.20	0.40
1:D:37:LEU:C	1:D:39:ALA:H	2.24	0.40
1:A:203:ILE:O	1:A:203:ILE:CG2	2.69	0.40
1:B:35:ILE:HG22	1:B:53:ARG:HD2	2.02	0.40
1:C:156:LYS:HG3	1:C:174:HIS:C	2.42	0.40
1:D:74:LEU:HD12	1:D:74:LEU:H	1.86	0.40
1:B:26:TRP:CZ2	1:B:60:MSE:HA	2.57	0.40
1:D:100:ASP:O	1:D:104:ASN:HB2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:116:PHE:HD1	1:D:134:HIS:ND1	2.19	0.40
1:A:63:LYS:HB3	1:A:63:LYS:HZ3	1.87	0.40
1:D:141:LYS:HZ2	1:D:167:ASN:ND2	2.19	0.40
1:D:200:LEU:C	1:D:202:LEU:H	2.23	0.40
1:D:7:PHE:CD1	1:D:162:VAL:HB	2.56	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	199/217 (92%)	190 (96%)	8 (4%)	1 (0%)	29	41
1	B	204/217 (94%)	195 (96%)	7 (3%)	2 (1%)	15	23
1	C	204/217 (94%)	192 (94%)	10 (5%)	2 (1%)	15	23
1	D	202/217 (93%)	183 (91%)	19 (9%)	0	100	100
All	All	809/868 (93%)	760 (94%)	44 (5%)	5 (1%)	25	36

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	THR
1	B	42	GLY
1	C	124	ASN
1	B	198	PRO
1	C	198	PRO

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	175/181 (97%)	165 (94%)	10 (6%)	20 33
1	B	176/181 (97%)	157 (89%)	19 (11%)	6 9
1	C	177/181 (98%)	160 (90%)	17 (10%)	8 12
1	D	175/181 (97%)	164 (94%)	11 (6%)	18 28
All	All	703/724 (97%)	646 (92%)	57 (8%)	11 18

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	36	THR
1	A	44	THR
1	A	67	GLU
1	A	81	LEU
1	A	89	ASN
1	A	112	LEU
1	A	146	LEU
1	A	167	ASN
1	A	202	LEU
1	B	20	GLU
1	B	22	LEU
1	B	57	LEU
1	B	80	GLU
1	B	81	LEU
1	B	88	LYS
1	B	91	LYS
1	B	117	SER
1	B	123	GLU
1	B	125	ASP
1	B	130	LEU
1	B	146	LEU
1	B	148	LEU
1	B	167	ASN
1	B	172	PHE
1	B	184	LYS
1	B	187	LEU
1	B	189	GLN

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Mol	Chain	Res	Type
1	B	202	LEU
1	C	1	LEU
1	C	22	LEU
1	C	30	ASP
1	C	36	THR
1	C	81	LEU
1	C	110	LEU
1	C	112	LEU
1	C	125	ASP
1	C	127	LEU
1	C	138	SER
1	C	146	LEU
1	C	150	ARG
1	C	152	LEU
1	C	184	LYS
1	C	187	LEU
1	C	202	LEU
1	C	204	LYS
1	D	36	THR
1	D	41	ASN
1	D	43	GLU
1	D	62	LEU
1	D	81	LEU
1	D	89	ASN
1	D	110	LEU
1	D	112	LEU
1	D	130	LEU
1	D	146	LEU
1	D	187	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	HIS
1	A	89	ASN
1	A	167	ASN
1	B	15	ASN
1	B	134	HIS
1	B	189	GLN
1	B	196	ASN
1	C	2	GLN
1	C	89	ASN

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Mol	Chain	Res	Type
1	C	189	GLN
1	C	193	HIS
1	C	196	ASN
1	D	41	ASN
1	D	47	HIS
1	D	89	ASN
1	D	167	ASN
1	D	182	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 10 ligands modelled in this entry, 10 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 (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	198/217 (91%)	-0.17	1 (0%)	91	89	18, 29, 48, 60	0
1	B	200/217 (92%)	0.06	3 (1%)	73	72	22, 40, 58, 64	0
1	C	202/217 (93%)	0.00	2 (0%)	82	80	17, 31, 61, 69	0
1	D	200/217 (92%)	0.27	7 (3%)	44	43	29, 48, 65, 70	0
All	All	800/868 (92%)	0.04	13 (1%)	72	70	17, 37, 61, 70	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	201	ALA	3.3
1	D	64	LEU	3.0
1	D	166	ALA	3.0
1	A	33	LYS	2.8
1	B	42	GLY	2.7
1	D	43	GLU	2.5
1	B	39	ALA	2.4
1	D	39	ALA	2.1
1	D	189	GLN	2.1
1	C	124	ASN	2.1
1	B	44	THR	2.1
1	D	139	HIS	2.1
1	C	0	SER	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 monosaccharides 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. The B-factors column lists the minimum, median, 95^{th} 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	B-factors(Å^2)	Q<0.9
2	MG	C	300	1/1	0.69	0.10	36,36,36,36	0
2	MG	A	300	1/1	0.85	0.14	31,31,31,31	0
2	MG	C	301	1/1	0.89	0.24	45,45,45,45	0
3	CL	C	302	1/1	0.90	0.12	55,55,55,55	0
2	MG	B	300	1/1	0.91	0.09	28,28,28,28	0
3	CL	D	302	1/1	0.93	0.08	55,55,55,55	0
3	CL	A	303	1/1	0.94	0.15	46,46,46,46	0
3	CL	B	302	1/1	0.94	0.08	55,55,55,55	0
2	MG	D	300	1/1	0.95	0.13	44,44,44,44	0
3	CL	A	302	1/1	0.97	0.06	38,38,38,38	0

6.5 Other polymers (i)

There are no such residues in this entry.