



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 06:30 am BST

PDB ID : 5AF0
Title : MAEL domain from Bombyx mori Maelstrom
Authors : Chen, K.; Campbell, E.; Pandey, R.R.; Yang, Z.; McCarthy, A.A.; Pillai, R.S.
Deposited on : 2015-01-13
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 ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

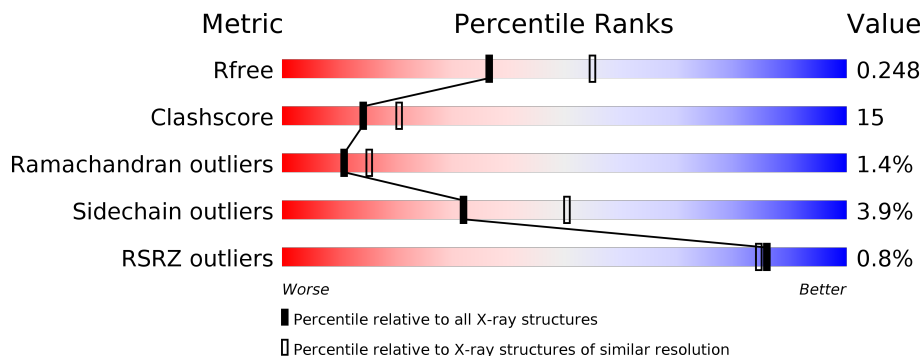
1 Overall quality at a glance i

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 on 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 $\leq 5\%$. 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	261	 % 60% 29% 9%
1	B	261	 % 62% 23% 5% 10%
1	C	261	 % 59% 31% 8%
1	D	261	 60% 29% 9%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7297 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 MAELSTROM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	237	1829	1170	297	350	12	0	0	0
1	B	236	1797	1147	292	345	13	0	0	0
1	C	239	1809	1163	291	342	13	0	0	0
1	D	238	1823	1170	295	345	13	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	298	LYS	ARG	conflict	UNP H9JFX7
B	298	LYS	ARG	conflict	UNP H9JFX7
C	298	LYS	ARG	conflict	UNP H9JFX7
D	298	LYS	ARG	conflict	UNP H9JFX7

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

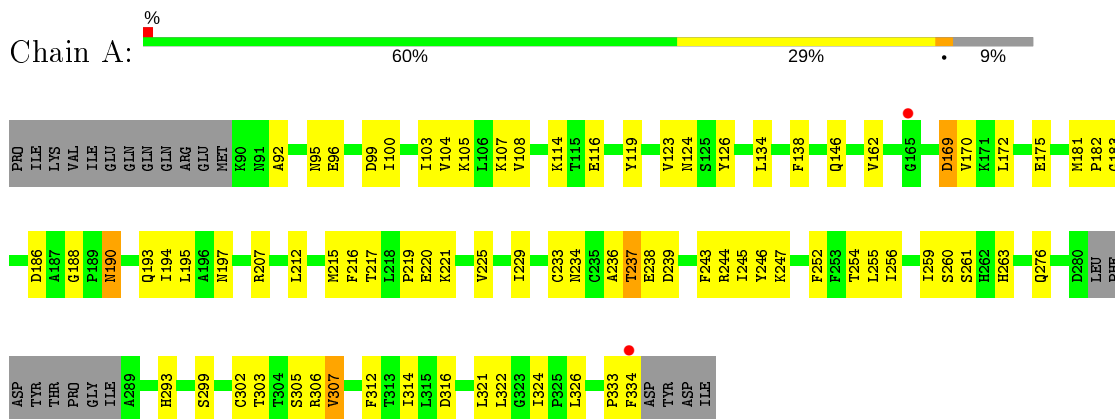
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	14	Total 14	O 14	0	0
3	B	9	Total 9	O 9	0	0
3	C	7	Total 7	O 7	0	0
3	D	5	Total 5	O 5	0	0

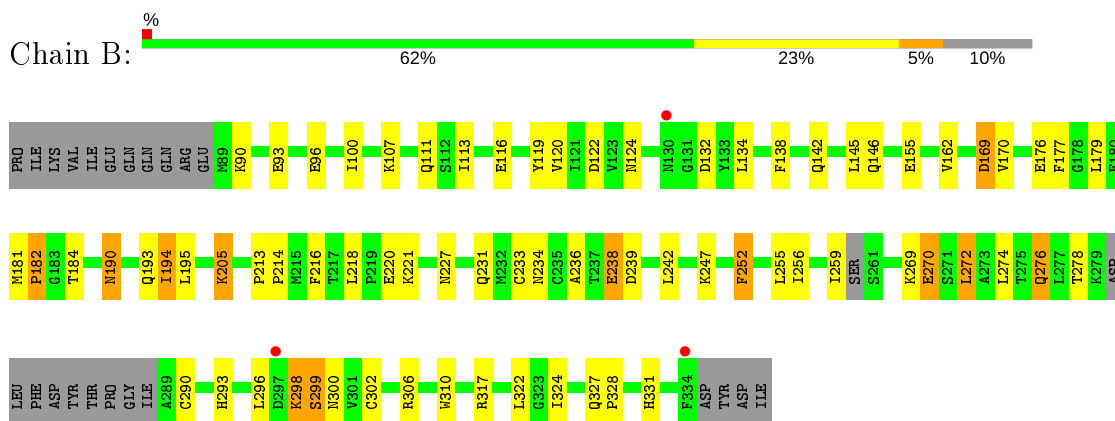
3 Residue-property plots [i](#)

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.

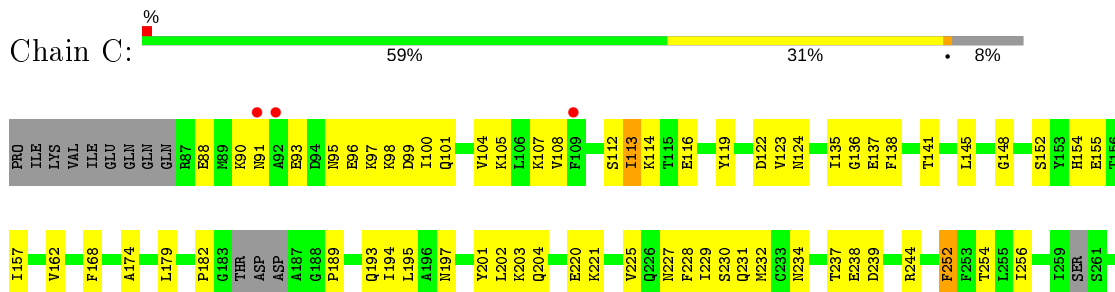
• Molecule 1: MAELSTROM



• Molecule 1: MAELSTROM



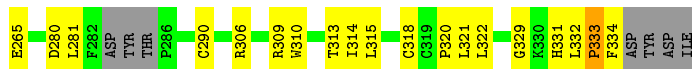
• Molecule 1: MAELSTROM





- Molecule 1: MAELSTROM

Chain D: 60% 29% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.37Å 101.21Å 101.64Å 90.00° 90.04° 90.00°	Depositor
Resolution (Å)	45.42 – 2.40 45.42 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.2 (45.42-2.40) 81.1 (45.42-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.96 (at 2.39Å)	Xtrriage
Refinement program	PHENIX dev_1839	Depositor
R, R_{free}	0.227 , 0.263 0.219 , 0.248	Depositor DCC
R_{free} test set	2870 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	35.4	Xtrriage
Anisotropy	0.951	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 59.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	0.000 for -h,l,k 0.000 for -h,-l,-k 0.426 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7297	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.39 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.5621e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 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:
ZN

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	0.51	0/1869	0.71	0/2544
1	B	0.49	0/1833	0.71	0/2495
1	C	0.53	0/1846	0.69	0/2509
1	D	0.50	0/1862	0.68	0/2528
All	All	0.51	0/7410	0.70	0/10076

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	B	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	B	298	LYS	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	1829	0	1731	60	0
1	B	1797	0	1693	49	0
1	C	1809	0	1698	57	0
1	D	1823	0	1698	54	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	14	0	0	1	0
3	B	9	0	0	1	0
3	C	7	0	0	0	0
3	D	5	0	0	0	0
All	All	7297	0	6820	218	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 15.

All (218) 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:262:HIS:HB2	1:D:265:GLU:HG3	1.58	0.85
1:A:260:SER:HA	1:A:321:LEU:HD13	1.59	0.83
1:D:91:ASN:O	1:D:95:ASN:ND2	2.15	0.80
1:C:99:ASP:OD2	1:C:244:ARG:NH1	2.15	0.80
1:A:124:ASN:OD1	1:A:306:ARG:NH2	2.14	0.80
1:B:124:ASN:OD1	1:B:306:ARG:NH2	2.15	0.79
1:B:259:ILE:HB	1:B:322:LEU:HD11	1.63	0.79
1:D:227:ASN:OD1	1:D:231:GLN:NE2	2.16	0.78
1:C:95:ASN:O	1:C:99:ASP:N	2.11	0.78
1:A:259:ILE:HB	1:A:322:LEU:HD11	1.65	0.76
1:D:122:ASP:OD2	1:D:306:ARG:NH1	2.19	0.76
1:B:132:ASP:HB3	1:B:231:GLN:HE21	1.53	0.72
1:B:276:GLN:HE22	1:B:317:ARG:HH11	1.38	0.72
1:B:146:GLN:HA	1:B:324:ILE:HD13	1.70	0.72
1:B:181:MET:HE3	1:B:182:PRO:HD2	1.71	0.70
1:C:152:SER:HA	1:C:201:TYR:OH	1.91	0.70
1:A:306:ARG:NH1	3:A:2006:HOH:O	2.21	0.70
1:A:212:LEU:O	1:A:244:ARG:NH2	2.25	0.69
1:B:190:ASN:HB2	1:B:193:GLN:HB3	1.73	0.69
1:D:94:ASP:OD1	1:D:95:ASN:ND2	2.26	0.69
1:B:145:LEU:HD11	1:B:255:LEU:HD11	1.75	0.69
1:A:252:PHE:CZ	1:A:314:ILE:HA	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:96:GLU:O	1:C:100:ILE:HG13	1.93	0.68
1:C:104:VAL:HG21	1:C:254:THR:HG22	1.76	0.68
1:B:146:GLN:O	1:B:331:HIS:NE2	2.25	0.67
1:A:194:ILE:HA	1:A:197:ASN:HD22	1.58	0.67
1:B:293:HIS:HB3	1:B:299:SER:HA	1.76	0.67
1:D:104:VAL:HG11	1:D:254:THR:HG22	1.78	0.66
1:B:111:GLN:OE1	3:B:2001:HOH:O	2.13	0.66
1:A:302:CYS:HB3	1:A:305:SER:HB2	1.79	0.65
1:C:193:GLN:HG3	1:C:197:ASN:HD21	1.62	0.64
1:C:268:PRO:HD2	1:C:272:LEU:HD22	1.80	0.63
1:D:138:PHE:O	1:D:154:HIS:HD2	1.82	0.63
1:A:221:LYS:O	1:A:225:VAL:HG22	1.99	0.63
1:C:227:ASN:OD1	1:C:231:GLN:NE2	2.31	0.63
1:C:194:ILE:HA	1:C:197:ASN:HD22	1.63	0.62
1:D:239:ASP:HB3	1:D:242:LEU:HD13	1.80	0.62
1:A:312:PHE:O	1:A:316:ASP:HB2	2.00	0.62
1:B:96:GLU:O	1:B:100:ILE:HG13	2.00	0.61
1:A:126:TYR:HB2	1:A:134:LEU:HD23	1.83	0.59
1:A:162:VAL:HG22	1:A:182:PRO:HG2	1.85	0.59
1:B:274:LEU:O	1:B:278:THR:N	2.34	0.59
1:B:119:TYR:OH	1:B:142:GLN:NE2	2.33	0.59
1:D:126:TYR:OH	1:D:180:GLU:O	2.07	0.59
1:A:169:ASP:OD1	1:A:169:ASP:N	2.36	0.58
1:D:105:LYS:HG2	1:D:106:LEU:HD12	1.86	0.58
1:A:225:VAL:HG23	1:A:245:ILE:HG21	1.86	0.58
1:D:329:GLY:N	1:D:332:LEU:O	2.30	0.57
1:C:135:ILE:HG13	1:C:228:PHE:CD1	2.39	0.57
1:B:270:GLU:CD	1:B:270:GLU:H	2.04	0.57
1:A:162:VAL:HG13	1:A:170:VAL:HG21	1.85	0.56
1:B:274:LEU:H	1:B:274:LEU:HD12	1.68	0.56
1:D:281:LEU:HG	1:D:310:TRP:HE1	1.69	0.56
1:B:190:ASN:O	1:B:194:ILE:HG23	2.06	0.56
1:C:95:ASN:HA	1:C:98:LYS:HB2	1.88	0.55
1:A:220:GLU:HG3	1:A:221:LYS:N	2.22	0.55
1:B:169:ASP:N	1:B:169:ASP:OD1	2.39	0.55
1:D:107:LYS:HE3	1:D:116:GLU:HG3	1.87	0.55
1:D:99:ASP:OD2	1:D:244:ARG:NH1	2.40	0.55
1:A:303:THR:O	1:A:307:VAL:HB	2.07	0.54
1:B:220:GLU:HG3	1:B:221:LYS:HG2	1.90	0.54
1:C:138:PHE:O	1:C:154:HIS:HD2	1.90	0.54
1:C:112:SER:O	1:C:116:GLU:HG2	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:100:ILE:O	1:C:104:VAL:HG22	2.08	0.54
1:C:124:ASN:HB2	1:C:137:GLU:HB3	1.90	0.54
1:C:193:GLN:O	1:C:197:ASN:ND2	2.41	0.54
1:D:104:VAL:HG21	1:D:255:LEU:HA	1.89	0.53
1:D:163:PRO:HD2	1:D:170:VAL:HG21	1.89	0.53
1:C:91:ASN:O	1:C:95:ASN:N	2.23	0.53
1:A:162:VAL:CG2	1:A:182:PRO:HG2	2.39	0.53
1:D:136:GLY:HA2	1:D:157:ILE:O	2.08	0.52
1:A:96:GLU:O	1:A:100:ILE:HG13	2.10	0.52
1:C:202:LEU:O	1:C:204:GLN:NE2	2.39	0.52
1:C:333:PRO:O	1:C:334:PHE:HB2	2.09	0.52
1:D:194:ILE:HD11	1:D:232:MET:SD	2.49	0.52
1:D:108:VAL:HG12	1:D:113:ILE:HG21	1.92	0.52
1:D:152:SER:HA	1:D:201:TYR:OH	2.09	0.52
1:C:195:LEU:HD22	1:C:232:MET:O	2.10	0.52
1:A:186:ASP:O	1:A:188:GLY:N	2.44	0.51
1:C:264:ASP:N	1:C:264:ASP:OD1	2.33	0.51
1:D:100:ILE:O	1:D:104:VAL:HG12	2.10	0.51
1:C:289:ALA:HB1	1:C:302:CYS:SG	2.51	0.51
1:B:162:VAL:HG13	1:B:170:VAL:HG21	1.93	0.51
1:A:193:GLN:O	1:A:197:ASN:ND2	2.44	0.50
1:C:230:SER:O	1:C:234:ASN:N	2.43	0.50
1:B:122:ASP:OD2	1:B:306:ARG:NH1	2.43	0.50
1:A:172:LEU:O	1:A:175:GLU:HB3	2.12	0.50
1:B:90:LYS:HA	1:B:93:GLU:HB3	1.92	0.50
1:B:145:LEU:HD21	1:B:255:LEU:HD21	1.94	0.50
1:A:181:MET:HE3	1:A:182:PRO:HD2	1.92	0.50
1:A:188:GLY:O	1:A:190:ASN:ND2	2.45	0.50
1:D:112:SER:O	1:D:116:GLU:HG2	2.10	0.50
1:C:148:GLY:HA2	1:C:315:LEU:HD11	1.93	0.50
1:D:122:ASP:OD1	1:D:221:LYS:HE3	2.12	0.50
1:C:108:VAL:HB	1:C:113:ILE:HD13	1.93	0.49
1:D:190:ASN:HA	1:D:193:GLN:HB3	1.95	0.49
1:D:333:PRO:O	1:D:334:PHE:HB2	2.13	0.49
1:A:236:ALA:O	1:A:238:GLU:N	2.45	0.49
1:D:111:GLN:O	1:D:114:LYS:NZ	2.45	0.49
1:A:219:PRO:HG3	1:A:247:LYS:HE3	1.93	0.49
1:D:315:LEU:HD13	1:D:331:HIS:CD2	2.48	0.48
1:A:236:ALA:C	1:A:238:GLU:H	2.16	0.48
1:A:233:CYS:O	1:A:238:GLU:N	2.46	0.48
1:C:225:VAL:O	1:C:229:ILE:HG12	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:296:LEU:O	1:B:298:LYS:N	2.42	0.48
1:B:290:CYS:HB2	1:B:302:CYS:SG	2.54	0.48
1:B:255:LEU:O	1:B:259:ILE:HG12	2.13	0.48
1:C:107:LYS:HD3	1:C:116:GLU:HG3	1.94	0.48
1:A:123:VAL:HG12	1:A:138:PHE:HB3	1.95	0.48
1:A:193:GLN:HG2	1:A:197:ASN:HD21	1.79	0.47
1:B:162:VAL:HG23	1:B:182:PRO:HG2	1.96	0.47
1:A:99:ASP:OD2	1:A:246:TYR:OH	2.28	0.47
1:D:280:ASP:HB3	1:D:313:THR:HG21	1.97	0.47
1:A:146:GLN:HA	1:A:324:ILE:HD13	1.97	0.47
1:B:239:ASP:O	1:B:242:LEU:HD13	2.15	0.47
1:C:124:ASN:ND2	1:C:301:VAL:O	2.39	0.47
1:B:227:ASN:O	1:B:231:GLN:HB2	2.15	0.47
1:A:92:ALA:HA	1:A:95:ASN:ND2	2.30	0.47
1:B:107:LYS:HZ3	1:B:116:GLU:CD	2.17	0.47
1:C:136:GLY:HA2	1:C:157:ILE:O	2.14	0.47
1:C:220:GLU:HG2	1:C:221:LYS:N	2.29	0.47
1:A:195:LEU:HD11	1:A:243:PHE:CZ	2.50	0.46
1:A:104:VAL:HG11	1:A:254:THR:HG22	1.98	0.46
1:B:177:PHE:HB2	1:B:179:LEU:HG	1.96	0.46
1:B:276:GLN:NE2	1:B:317:ARG:HH11	2.10	0.46
1:C:138:PHE:CE2	1:C:155:GLU:HB2	2.50	0.46
1:A:107:LYS:HD3	1:A:116:GLU:HG2	1.97	0.46
1:C:315:LEU:HD13	1:C:331:HIS:CD2	2.51	0.46
1:D:318:CYS:O	1:D:321:LEU:HB2	2.16	0.46
1:C:93:GLU:O	1:C:97:LYS:HG2	2.16	0.46
1:B:269:LYS:O	1:B:272:LEU:HB2	2.16	0.45
1:C:154:HIS:CE1	1:C:290:CYS:HB2	2.51	0.45
1:D:118:PHE:HB2	1:D:143:PHE:CE1	2.51	0.45
1:D:252:PHE:CZ	1:D:314:ILE:HA	2.51	0.45
1:D:136:GLY:O	1:D:157:ILE:N	2.46	0.45
1:C:124:ASN:OD1	1:C:306:ARG:NH1	2.49	0.45
1:D:162:VAL:HG22	1:D:182:PRO:HG2	1.99	0.45
1:C:114:LYS:HA	1:C:145:LEU:HB2	1.98	0.45
1:C:138:PHE:O	1:C:154:HIS:CD2	2.69	0.45
1:A:107:LYS:HE3	1:C:168:PHE:CD2	2.52	0.45
1:C:195:LEU:HD13	1:C:232:MET:HB3	1.99	0.44
1:A:252:PHE:O	1:A:256:ILE:HG22	2.18	0.44
1:D:281:LEU:O	1:D:309:ARG:NE	2.51	0.43
1:B:233:CYS:HB3	1:B:238:GLU:O	2.17	0.43
1:B:252:PHE:O	1:B:256:ILE:HG22	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:ASN:O	1:C:96:GLU:C	2.56	0.43
1:B:90:LYS:NZ	1:B:93:GLU:OE2	2.44	0.43
1:D:227:ASN:O	1:D:231:GLN:HG3	2.18	0.43
1:B:138:PHE:CE2	1:B:155:GLU:HB2	2.54	0.43
1:B:190:ASN:H	1:B:190:ASN:ND2	2.15	0.43
1:B:218:LEU:HD23	1:B:218:LEU:HA	1.74	0.43
1:C:154:HIS:CE1	1:C:290:CYS:CB	3.01	0.43
1:C:193:GLN:HG3	1:C:197:ASN:ND2	2.31	0.43
1:A:263:HIS:CE1	1:D:191:TYR:CE2	3.06	0.43
1:C:145:LEU:HD13	1:C:322:LEU:CD1	2.48	0.43
1:C:329:GLY:N	1:C:332:LEU:O	2.52	0.43
1:A:225:VAL:O	1:A:229:ILE:HG12	2.19	0.43
1:A:293:HIS:HB3	1:A:299:SER:HA	2.00	0.43
1:A:123:VAL:HG12	1:A:138:PHE:CB	2.48	0.43
1:A:293:HIS:HB3	1:A:299:SER:N	2.34	0.43
1:A:255:LEU:O	1:A:259:ILE:HG12	2.19	0.43
1:C:98:LYS:HA	1:C:101:GLN:OE1	2.19	0.43
1:C:123:VAL:HG12	1:C:138:PHE:HB3	2.01	0.43
1:D:102:ASN:O	1:D:106:LEU:HD13	2.18	0.43
1:A:236:ALA:C	1:A:238:GLU:N	2.72	0.42
1:A:293:HIS:HB3	1:A:299:SER:H	1.83	0.42
1:C:122:ASP:OD1	1:C:221:LYS:HE3	2.19	0.42
1:D:94:ASP:OD1	1:D:95:ASN:N	2.52	0.42
1:B:122:ASP:OD2	1:B:306:ARG:HD2	2.19	0.42
1:C:157:ILE:HG23	1:C:189:PRO:CB	2.48	0.42
1:A:234:ASN:C	1:A:236:ALA:H	2.23	0.42
1:A:238:GLU:HG3	1:A:239:ASP:N	2.34	0.42
1:A:216:PHE:HA	1:A:246:TYR:O	2.19	0.42
1:A:293:HIS:HB3	1:A:299:SER:CA	2.49	0.42
1:D:96:GLU:O	1:D:100:ILE:HG13	2.19	0.42
1:D:109:PHE:HD1	1:D:109:PHE:HA	1.71	0.42
1:D:214:PRO:HG3	1:D:244:ARG:CZ	2.49	0.42
1:A:333:PRO:O	1:A:334:PHE:HB2	2.20	0.42
1:A:114:LYS:HG2	1:A:322:LEU:HD23	2.01	0.42
1:B:234:ASN:C	1:B:236:ALA:H	2.23	0.42
1:C:119:TYR:HA	1:C:141:THR:O	2.20	0.42
1:C:267:PHE:HA	1:C:268:PRO:HD3	1.87	0.42
1:A:103:ILE:O	1:A:107:LYS:HG3	2.20	0.42
1:A:316:ASP:HA	1:A:326:LEU:HD21	2.01	0.42
1:B:194:ILE:HG13	1:B:195:LEU:N	2.34	0.42
1:D:154:HIS:CE1	1:D:290:CYS:CB	3.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:190:ASN:O	1:D:194:ILE:N	2.38	0.42
1:B:120:VAL:HG23	1:B:216:PHE:O	2.20	0.41
1:C:298:LYS:HA	1:C:299:SER:CB	2.49	0.41
1:D:138:PHE:CE2	1:D:155:GLU:HB2	2.55	0.41
1:A:119:TYR:O	1:A:215:MET:HA	2.20	0.41
1:C:162:VAL:HG22	1:C:182:PRO:HG2	2.02	0.41
1:D:89:MET:HG3	1:D:89:MET:O	2.12	0.41
1:D:96:GLU:HG3	1:D:246:TYR:CE1	2.55	0.41
1:C:321:LEU:HA	1:C:321:LEU:HD23	1.85	0.41
1:D:204:GLN:HB2	1:D:207:ARG:HB2	2.01	0.41
1:C:174:ALA:HA	1:C:179:LEU:HB2	2.03	0.41
1:D:229:ILE:HD13	1:D:229:ILE:HA	1.88	0.41
1:A:105:LYS:O	1:A:108:VAL:HG12	2.20	0.41
1:C:252:PHE:O	1:C:256:ILE:HG22	2.21	0.41
1:D:92:ALA:HA	1:D:95:ASN:HB2	2.02	0.41
1:B:213:PRO:HA	1:B:214:PRO:HD3	1.95	0.41
1:C:227:ASN:O	1:C:231:GLN:HG3	2.21	0.41
1:D:148:GLY:HA2	1:D:315:LEU:HD11	2.03	0.41
1:A:260:SER:HA	1:A:321:LEU:CD1	2.40	0.41
1:B:162:VAL:CG2	1:B:182:PRO:HG2	2.51	0.41
1:B:218:LEU:HG	1:B:310:TRP:CZ2	2.56	0.41
1:A:186:ASP:C	1:A:188:GLY:N	2.75	0.41
1:B:107:LYS:NZ	1:B:116:GLU:OE2	2.54	0.40
1:C:104:VAL:HG23	1:C:105:LYS:N	2.36	0.40
1:A:217:THR:HG21	1:A:225:VAL:HG21	2.04	0.40
1:D:132:ASP:HA	1:D:227:ASN:OD1	2.22	0.40
1:D:135:ILE:HG13	1:D:228:PHE:CD1	2.57	0.40
1:B:247:LYS:HE2	1:B:247:LYS:HB3	1.85	0.40
1:D:320:PRO:C	1:D:322:LEU:H	2.25	0.40
1:A:220:GLU:HG3	1:A:221:LYS:CG	2.52	0.40
1:B:327:GLN:HA	1:B:328:PRO:HD3	1.79	0.40
1:D:107:LYS:HB2	1:D:113:ILE:HD12	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	233/261 (89%)	220 (94%)	10 (4%)	3 (1%)	12	17
1	B	230/261 (88%)	216 (94%)	10 (4%)	4 (2%)	9	11
1	C	231/261 (88%)	218 (94%)	9 (4%)	4 (2%)	9	11
1	D	230/261 (88%)	215 (94%)	13 (6%)	2 (1%)	17	25
All	All	924/1044 (88%)	869 (94%)	42 (4%)	13 (1%)	11	15

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	205	LYS
1	B	299	SER
1	B	184	THR
1	C	88	GLU
1	C	300	ASN
1	A	237	THR
1	C	90	LYS
1	D	87	ARG
1	A	261	SER
1	D	333	PRO
1	C	299	SER
1	A	183	GLY
1	B	182	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	198/239 (83%)	192 (97%)	6 (3%)	41	61
1	B	192/239 (80%)	179 (93%)	13 (7%)	16	25
1	C	189/239 (79%)	183 (97%)	6 (3%)	39	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	191/239 (80%)	186 (97%)	5 (3%)	46	66
All	All	770/956 (80%)	740 (96%)	30 (4%)	32	50

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	ASP
1	A	190	ASN
1	A	207	ARG
1	A	237	THR
1	A	276	GLN
1	A	307	VAL
1	B	113	ILE
1	B	134	LEU
1	B	169	ASP
1	B	176	GLU
1	B	190	ASN
1	B	194	ILE
1	B	205	LYS
1	B	238	GLU
1	B	252	PHE
1	B	270	GLU
1	B	272	LEU
1	B	276	GLN
1	B	300	ASN
1	C	113	ILE
1	C	203	LYS
1	C	237	THR
1	C	238	GLU
1	C	239	ASP
1	C	252	PHE
1	D	89	MET
1	D	109	PHE
1	D	114	LYS
1	D	238	GLU
1	D	252	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	ASN

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Mol	Chain	Res	Type
1	A	197	ASN
1	B	142	GLN
1	B	190	ASN
1	B	204	GLN
1	B	231	GLN
1	C	95	ASN
1	C	154	HIS
1	C	197	ASN
1	D	102	ASN
1	D	154	HIS
1	D	193	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](#)

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 [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	237/261 (90%)	-0.15	2 (0%) 86 84	32, 53, 98, 120	0
1	B	236/261 (90%)	-0.01	3 (1%) 77 75	33, 60, 109, 134	0
1	C	239/261 (91%)	-0.09	3 (1%) 77 75	32, 57, 91, 135	0
1	D	238/261 (91%)	-0.08	0 100 100	32, 58, 92, 136	0
All	All	950/1044 (90%)	-0.08	8 (0%) 86 84	32, 57, 98, 136	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	92	ALA	4.1
1	B	297	ASP	3.4
1	B	334	PHE	3.3
1	B	130	ASN	3.0
1	C	91	ASN	2.6
1	C	109	PHE	2.4
1	A	334	PHE	2.3
1	A	165	GLY	2.1

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. 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	B-factors(Å ²)	Q<0.9
2	ZN	B	501	1/1	0.98	0.19	86,86,86,86	0
2	ZN	D	501	1/1	0.98	0.13	73,73,73,73	0
2	ZN	A	501	1/1	0.99	0.14	85,85,85,85	0
2	ZN	C	501	1/1	0.99	0.16	73,73,73,73	0

6.5 Other polymers [i](#)

There are no such residues in this entry.