



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:25 PM UTC

PDB ID : 4MXI / pdb_00004mxi
Title : ClpP Ser98dhA
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Deposited on : 2013-09-26
Resolution : 2.30 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

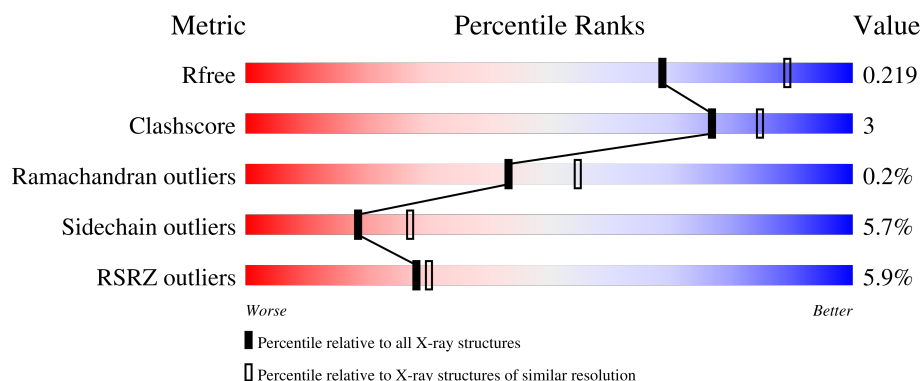
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 $\leq 5\%$. 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	195	4% 78% 9% • 12%
1	B	195	7% 75% 9% • 15%
1	C	195	6% 76% 10% • 13%
1	D	195	4% 74% 10% • 14%
1	E	195	6% 77% 9% • 13%

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Mol	Chain	Length	Quality of chain
1	F	195	
1	G	195	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	DHA	G	98	-	X	-	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9634 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 ATP-dependent Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	0	0	0
			1330	837	226	261	6			
1	B	166	Total	C	N	O	S	0	0	0
			1290	814	219	251	6			
1	C	169	Total	C	N	O	S	0	0	0
			1314	829	222	257	6			
1	D	167	Total	C	N	O	S	0	0	0
			1302	823	220	253	6			
1	E	170	Total	C	N	O	S	0	0	0
			1324	837	224	257	6			
1	F	170	Total	C	N	O	S	0	0	0
			1324	837	224	257	6			
1	G	171	Total	C	N	O	S	0	0	0
			1333	842	225	260	6			

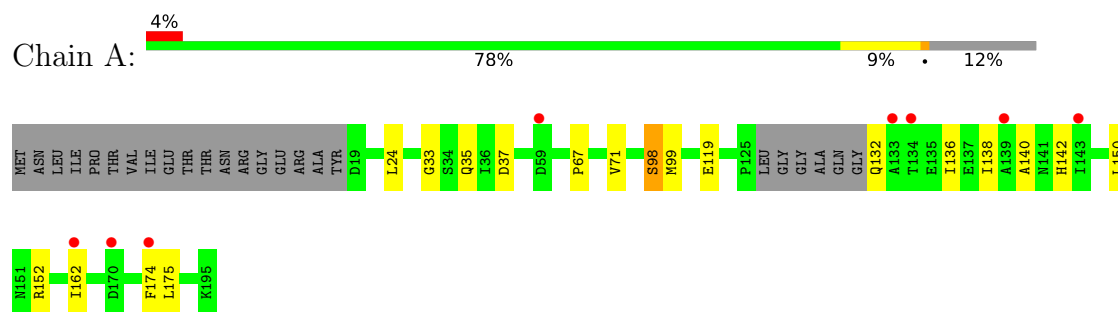
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	64	Total	O	0	0
			64	64		
2	B	56	Total	O	0	0
			56	56		
2	C	59	Total	O	0	0
			59	59		
2	D	62	Total	O	0	0
			62	62		
2	E	69	Total	O	0	0
			69	69		
2	F	66	Total	O	0	0
			66	66		
2	G	41	Total	O	0	0
			41	41		

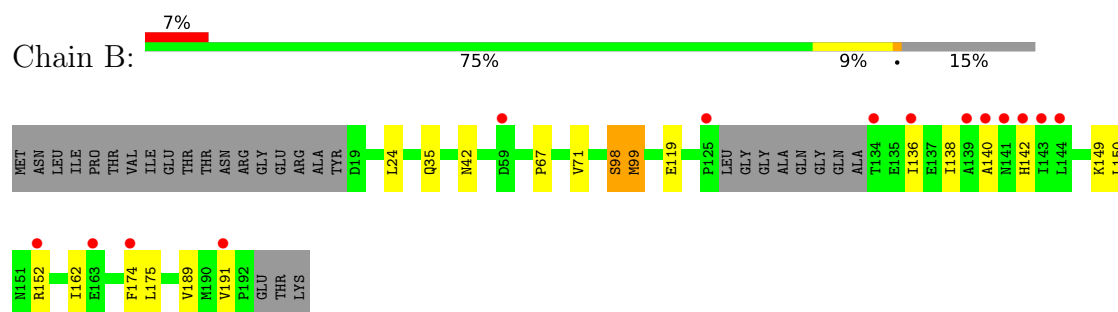
3 Residue-property plots [i](#)

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.

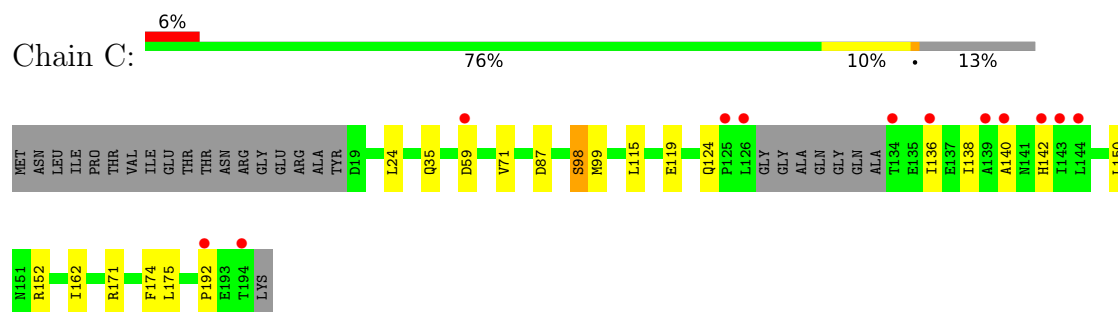
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



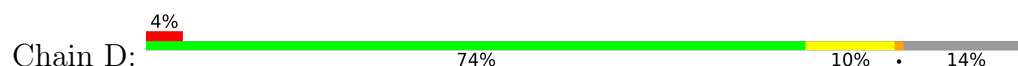
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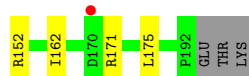
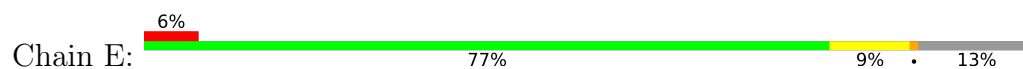


- Molecule 1: ATP-dependent Clp protease proteolytic subunit

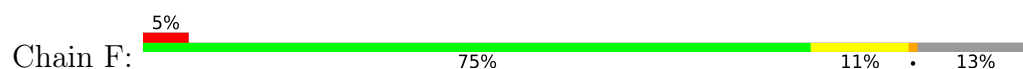




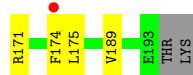
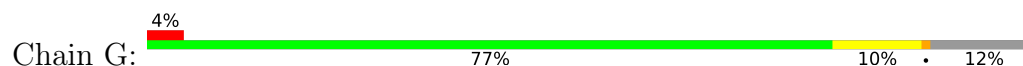
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.59Å 121.59Å 402.95Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.30 15.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (15.00-2.30) 99.5 (15.00-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.84 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.196 , 0.221 0.198 , 0.219	Depositor DCC
R_{free} test set	3942 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	50.3	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 67.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9634	wwPDB-VP
Average B, all atoms (Å ²)	61.0	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: DHA

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	0.38	0/1340	0.71	0/1803
1	B	0.38	0/1300	0.71	0/1751
1	C	0.38	0/1324	0.72	0/1784
1	D	0.37	0/1313	0.71	0/1769
1	E	0.39	0/1335	0.71	0/1799
1	F	0.38	0/1335	0.72	0/1799
1	G	0.38	0/1344	0.71	0/1811
All	All	0.38	0/9291	0.71	0/12516

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
1	C	0	1
1	E	0	1
1	F	0	1
1	G	0	1
All	All	0	7

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	98	DHA	Mainchain
1	B	98	DHA	Mainchain,Peptide
1	C	98	DHA	Mainchain
1	E	98	DHA	Mainchain
1	F	98	DHA	Mainchain
1	G	98	DHA	Mainchain

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	1330	0	1343	10	0
1	B	1290	0	1304	11	0
1	C	1314	0	1328	9	0
1	D	1302	0	1313	12	0
1	E	1324	0	1337	8	0
1	F	1324	0	1337	10	0
1	G	1333	0	1343	9	0
2	A	64	0	0	0	0
2	B	56	0	0	0	1
2	C	59	0	0	0	0
2	D	62	0	0	1	0
2	E	69	0	0	0	1
2	F	66	0	0	0	0
2	G	41	0	0	0	0
All	All	9634	0	9305	54	1

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 3.

All (54) 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:F:18:TYR:HB3	1:F:20:ILE:HG22	1.61	0.82
1:A:174:PHE:CE1	1:B:140:ALA:HB2	2.30	0.67
1:B:152:ARG:HG3	1:B:162:ILE:HD11	1.79	0.64
1:B:174:PHE:CE1	1:C:140:ALA:HB2	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:ARG:HG3	1:A:162:ILE:HD11	1.86	0.58
1:D:174:PHE:CE1	1:E:140:ALA:HB2	2.40	0.56
1:G:152:ARG:HG3	1:G:162:ILE:HD11	1.86	0.56
1:C:152:ARG:HG3	1:C:162:ILE:HD11	1.88	0.56
1:E:152:ARG:HG3	1:E:162:ILE:HD11	1.88	0.55
1:D:18:TYR:HB3	1:D:20:ILE:HG22	1.87	0.55
1:D:152:ARG:HG3	1:D:162:ILE:HD11	1.89	0.55
1:F:138:ILE:O	1:F:142:HIS:HD2	1.90	0.55
1:F:149:LYS:HE3	1:F:152:ARG:HH12	1.73	0.54
1:A:37:ASP:HB3	1:A:132:GLN:HE21	1.75	0.52
1:A:67:PRO:HB2	1:B:136:ILE:HD12	1.92	0.51
1:A:138:ILE:O	1:A:142:HIS:HD2	1.93	0.51
1:E:138:ILE:O	1:E:142:HIS:HD2	1.93	0.51
1:C:174:PHE:CE1	1:D:140:ALA:HB2	2.45	0.51
1:D:138:ILE:O	1:D:142:HIS:HD2	1.94	0.51
1:A:140:ALA:HB2	1:G:174:PHE:CE1	2.46	0.50
1:F:174:PHE:CE1	1:G:140:ALA:HB2	2.47	0.49
1:C:138:ILE:O	1:C:142:HIS:HD2	1.96	0.48
1:D:67:PRO:HB2	1:E:136:ILE:HD12	1.96	0.47
1:G:138:ILE:O	1:G:142:HIS:HD2	1.98	0.46
1:B:149:LYS:HE3	1:B:152:ARG:HH12	1.81	0.46
1:D:99:MET:HE1	1:D:150:LEU:CD2	2.46	0.46
1:B:99:MET:HE1	1:B:150:LEU:CD2	2.46	0.46
1:D:134:THR:OG1	2:D:237:HOH:O	2.20	0.46
1:B:138:ILE:O	1:B:142:HIS:HD2	1.99	0.45
1:C:99:MET:HE1	1:C:150:LEU:CD2	2.47	0.45
1:B:67:PRO:HB2	1:C:136:ILE:HD12	1.98	0.44
1:A:99:MET:HE1	1:A:150:LEU:HD22	1.99	0.44
1:A:99:MET:HE1	1:A:150:LEU:CD2	2.48	0.44
1:C:99:MET:HE1	1:C:150:LEU:HD22	1.99	0.44
1:F:138:ILE:O	1:F:142:HIS:CD2	2.71	0.43
1:C:115:LEU:HD13	1:D:79:ASP:HB3	2.00	0.43
1:E:99:MET:HE1	1:E:150:LEU:HD22	2.02	0.42
1:F:67:PRO:HB2	1:G:136:ILE:HD12	2.01	0.42
1:F:99:MET:HE1	1:F:150:LEU:HD22	2.00	0.42
1:E:99:MET:HE1	1:E:150:LEU:CD2	2.49	0.42
1:D:99:MET:HE1	1:D:150:LEU:HD22	2.01	0.42
1:F:99:MET:HE1	1:F:150:LEU:CD2	2.50	0.42
1:D:149:LYS:HE3	1:D:152:ARG:HH12	1.85	0.41
1:B:99:MET:HE1	1:B:150:LEU:HD22	2.02	0.41
1:G:99:MET:HE1	1:G:150:LEU:CD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:59:ASP:HB2	1:C:87:ASP:O	2.20	0.41
1:A:33:GLY:O	1:B:42:ASN:ND2	2.54	0.41
1:B:149:LYS:HE3	1:B:152:ARG:HH22	1.84	0.41
1:D:115:LEU:HD13	1:E:79:ASP:HB3	2.03	0.41
1:E:149:LYS:HE3	1:E:152:ARG:HH12	1.86	0.41
1:A:136:ILE:HD12	1:G:67:PRO:HB2	2.04	0.40
1:G:99:MET:HE1	1:G:150:LEU:HD22	2.03	0.40
1:F:93:ILE:HD12	1:G:45:VAL:HG11	2.04	0.40
1:F:149:LYS:HE3	1:F:152:ARG:NH1	2.36	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:250:HOH:O	2:E:201:HOH:O[5_554]	2.14	0.06

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	166/195 (85%)	164 (99%)	2 (1%)	0	100	100
1	B	161/195 (83%)	156 (97%)	4 (2%)	1 (1%)	21	27
1	C	164/195 (84%)	159 (97%)	4 (2%)	1 (1%)	21	27
1	D	162/195 (83%)	159 (98%)	3 (2%)	0	100	100
1	E	165/195 (85%)	162 (98%)	3 (2%)	0	100	100
1	F	165/195 (85%)	162 (98%)	3 (2%)	0	100	100
1	G	166/195 (85%)	161 (97%)	5 (3%)	0	100	100
All	All	1149/1365 (84%)	1123 (98%)	24 (2%)	2 (0%)	43	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	99	MET
1	C	192	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	144/162 (89%)	139 (96%)	5 (4%)	32	48
1	B	140/162 (86%)	133 (95%)	7 (5%)	22	33
1	C	143/162 (88%)	136 (95%)	7 (5%)	22	34
1	D	141/162 (87%)	133 (94%)	8 (6%)	18	27
1	E	143/162 (88%)	135 (94%)	8 (6%)	19	28
1	F	143/162 (88%)	130 (91%)	13 (9%)	9	11
1	G	144/162 (89%)	135 (94%)	9 (6%)	16	23
All	All	998/1134 (88%)	941 (94%)	57 (6%)	18	27

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

Mol	Chain	Res	Type
1	A	24	LEU
1	A	35	GLN
1	A	71	VAL
1	A	119	GLU
1	A	175	LEU
1	B	24	LEU
1	B	35	GLN
1	B	71	VAL
1	B	119	GLU
1	B	175	LEU
1	B	189	VAL
1	B	191	VAL
1	C	24	LEU
1	C	35	GLN
1	C	71	VAL
1	C	119	GLU

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Mol	Chain	Res	Type
1	C	124	GLN
1	C	171	ARG
1	C	175	LEU
1	D	18	TYR
1	D	24	LEU
1	D	35	GLN
1	D	71	VAL
1	D	119	GLU
1	D	134	THR
1	D	171	ARG
1	D	175	LEU
1	E	24	LEU
1	E	35	GLN
1	E	71	VAL
1	E	119	GLU
1	E	124	GLN
1	E	126	LEU
1	E	171	ARG
1	E	175	LEU
1	F	18	TYR
1	F	24	LEU
1	F	35	GLN
1	F	71	VAL
1	F	119	GLU
1	F	124	GLN
1	F	126	LEU
1	F	134	THR
1	F	162	ILE
1	F	171	ARG
1	F	175	LEU
1	F	189	VAL
1	F	191	VAL
1	G	24	LEU
1	G	35	GLN
1	G	71	VAL
1	G	119	GLU
1	G	124	GLN
1	G	126	LEU
1	G	171	ARG
1	G	175	LEU
1	G	189	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	82	GLN
1	A	132	GLN
1	A	141	ASN
1	A	142	HIS
1	B	35	GLN
1	B	52	GLN
1	B	65	ASN
1	B	82	GLN
1	B	142	HIS
1	C	35	GLN
1	C	42	ASN
1	C	82	GLN
1	C	142	HIS
1	D	35	GLN
1	D	52	GLN
1	D	82	GLN
1	D	83	HIS
1	D	142	HIS
1	E	35	GLN
1	E	82	GLN
1	E	117	ASN
1	E	142	HIS
1	F	35	GLN
1	F	54	GLN
1	F	82	GLN
1	F	142	HIS
1	G	35	GLN
1	G	54	GLN
1	G	82	GLN
1	G	83	HIS
1	G	142	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

7 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
1	DHA	E	98	1	3,4,5	1.61	1 (33%)	2,4,6	2.67	2 (100%)
1	DHA	A	98	1	3,4,5	2.22	1 (33%)	2,4,6	2.66	2 (100%)
1	DHA	G	98	1	3,4,5	3.91	2 (66%)	2,4,6	2.64	2 (100%)
1	DHA	F	98	1	3,4,5	0.66	0	2,4,6	2.68	2 (100%)
1	DHA	C	98	1	3,4,5	2.24	1 (33%)	2,4,6	2.65	2 (100%)
1	DHA	D	98	1	3,4,5	2.03	1 (33%)	2,4,6	2.66	2 (100%)
1	DHA	B	98	1	3,4,5	2.10	1 (33%)	2,4,6	2.66	2 (100%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	DHA	E	98	1	-	0/0/2/4	-
1	DHA	A	98	1	-	0/0/2/4	-
1	DHA	G	98	1	-	0/0/2/4	-
1	DHA	F	98	1	-	0/0/2/4	-
1	DHA	C	98	1	-	0/0/2/4	-
1	DHA	D	98	1	-	0/0/2/4	-
1	DHA	B	98	1	-	0/0/2/4	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	98	DHA	C-CA	-6.12	1.34	1.45
1	C	98	DHA	C-CA	3.86	1.51	1.45
1	A	98	DHA	O-C	-3.46	1.14	1.22
1	B	98	DHA	C-CA	-3.36	1.39	1.45
1	D	98	DHA	C-CA	-3.24	1.39	1.45
1	G	98	DHA	CB-CA	-2.82	1.22	1.34
1	E	98	DHA	C-CA	-2.64	1.40	1.45

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	98	DHA	O-C-CA	-2.92	120.05	125.53
1	F	98	DHA	O-C-CA	-2.92	120.05	125.53
1	B	98	DHA	O-C-CA	-2.91	120.07	125.53
1	E	98	DHA	O-C-CA	-2.91	120.08	125.53
1	A	98	DHA	O-C-CA	-2.90	120.08	125.53
1	C	98	DHA	O-C-CA	-2.90	120.09	125.53
1	G	98	DHA	O-C-CA	-2.88	120.12	125.53
1	F	98	DHA	CB-CA-N	-2.41	119.96	125.76
1	E	98	DHA	CB-CA-N	-2.40	119.97	125.76
1	B	98	DHA	CB-CA-N	-2.39	120.01	125.76
1	A	98	DHA	CB-CA-N	-2.39	120.01	125.76
1	G	98	DHA	CB-CA-N	-2.38	120.03	125.76
1	C	98	DHA	CB-CA-N	-2.37	120.05	125.76
1	D	98	DHA	CB-CA-N	-2.37	120.05	125.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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	A	170/195 (87%)	0.12	8 (4%) 36 38	37, 55, 93, 110	0
1	B	165/195 (84%)	0.12	14 (8%) 16 18	37, 52, 93, 108	0
1	C	168/195 (86%)	0.08	12 (7%) 22 24	35, 52, 91, 120	0
1	D	166/195 (85%)	0.05	7 (4%) 40 42	37, 55, 94, 121	0
1	E	169/195 (86%)	0.13	12 (7%) 22 24	37, 56, 97, 115	0
1	F	169/195 (86%)	0.06	10 (5%) 28 30	39, 53, 82, 112	0
1	G	170/195 (87%)	0.16	7 (4%) 41 43	45, 63, 102, 118	0
All	All	1177/1365 (86%)	0.10	70 (5%) 28 30	35, 55, 94, 121	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	194	THR	5.7
1	C	143	ILE	4.6
1	G	126	LEU	4.4
1	A	133	ALA	4.3
1	E	126	LEU	4.2
1	B	139	ALA	3.9
1	C	126	LEU	3.8
1	A	143	ILE	3.6
1	B	134	THR	3.5
1	F	18	TYR	3.4
1	C	139	ALA	3.3
1	F	191	VAL	3.3
1	B	140	ALA	3.2
1	D	144	LEU	3.2
1	F	126	LEU	3.1
1	B	191	VAL	3.1
1	B	143	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	134	THR	3.0
1	B	152	ARG	3.0
1	B	144	LEU	2.9
1	C	142	HIS	2.9
1	E	125	PRO	2.9
1	F	148	GLU	2.8
1	D	191	VAL	2.7
1	F	59	ASP	2.7
1	E	144	LEU	2.7
1	C	140	ALA	2.6
1	C	136	ILE	2.6
1	A	162	ILE	2.6
1	B	136	ILE	2.6
1	F	136	ILE	2.6
1	A	134	THR	2.6
1	C	192	PRO	2.6
1	C	59	ASP	2.6
1	B	125	PRO	2.6
1	C	125	PRO	2.6
1	G	174	PHE	2.6
1	E	93	ILE	2.5
1	A	170	ASP	2.5
1	E	133	ALA	2.5
1	D	140	ALA	2.5
1	G	18	TYR	2.4
1	E	18	TYR	2.4
1	E	124	GLN	2.4
1	G	59	ASP	2.4
1	E	140	ALA	2.4
1	B	174	PHE	2.3
1	B	59	ASP	2.3
1	E	134	THR	2.3
1	B	163	GLU	2.3
1	A	59	ASP	2.3
1	F	162	ILE	2.3
1	F	141	ASN	2.2
1	B	142	HIS	2.2
1	G	132	GLN	2.2
1	C	144	LEU	2.2
1	E	19	ASP	2.2
1	D	192	PRO	2.2
1	F	142	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	124	GLN	2.2
1	A	139	ALA	2.2
1	D	134	THR	2.2
1	D	139	ALA	2.2
1	B	141	ASN	2.1
1	A	174	PHE	2.1
1	D	18	TYR	2.1
1	E	139	ALA	2.0
1	F	31	MET	2.0
1	E	170	ASP	2.0
1	G	125	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	DHA	E	98	5/6	0.88	0.11	57,57,61,67	0
1	DHA	B	98	5/6	0.89	0.12	52,52,55,59	0
1	DHA	G	98	5/6	0.89	0.09	60,60,62,65	0
1	DHA	F	98	5/6	0.92	0.09	50,53,56,61	0
1	DHA	D	98	5/6	0.94	0.08	55,60,62,63	0
1	DHA	C	98	5/6	0.94	0.07	50,50,52,59	0
1	DHA	A	98	5/6	0.95	0.07	53,57,62,66	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.