



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 03:37 PM GMT

PDB ID : 4AAL
Title : MacA wild-type oxidized
Authors : Seidel, J.
Deposited on : 2011-12-05
Resolution : 1.84 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

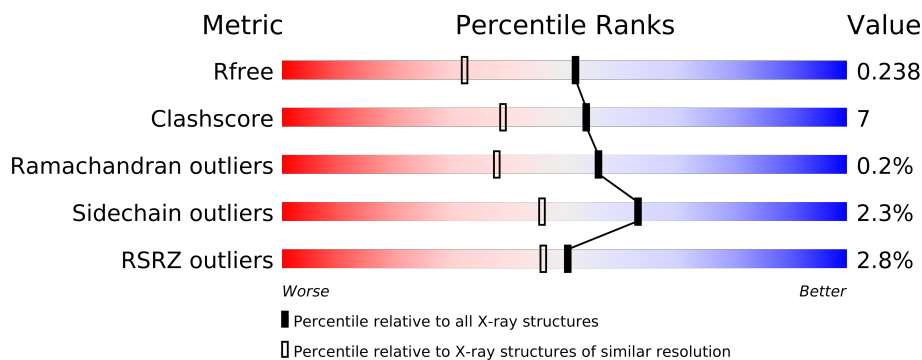
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.84 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}	66092	1857 (1.86-1.82)
Clashscore	79885	2149 (1.86-1.82)
Ramachandran outliers	78287	2124 (1.86-1.82)
Sidechain outliers	78261	2125 (1.86-1.82)
RSRZ outliers	66119	1857 (1.86-1.82)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	341	
1	B	341	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	PO4	A	1348	-	X
4	PO4	A	1349	-	X
4	PO4	A	1351	-	X
4	PO4	A	1352	-	X
4	PO4	A	1355	-	X
4	PO4	A	1362	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
4	PO4	B	1350	-	X
4	PO4	B	1357	-	X
4	PO4	B	1358	-	X
4	PO4	B	1359	-	X
5	ACT	A	1360	-	X
5	ACT	B	1360	-	X
5	ACT	B	1361	-	X
5	ACT	B	1362	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5841 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CYTOCHROME C551 PEROXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	3	0
			2447	1553	426	455	13			
1	B	319	Total	C	N	O	S	0	4	0
			2449	1553	426	457	13			

There are 34 discrepancies between the modelled and reference sequences:

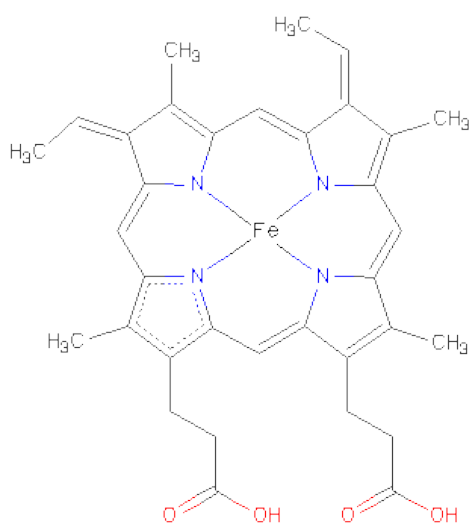
Chain	Residue	Modelled	Actual	Comment	Reference
A	6	TRP	-	EXPRESSION TAG	UNP Q74FY6
A	7	SER	-	EXPRESSION TAG	UNP Q74FY6
A	8	HIS	-	EXPRESSION TAG	UNP Q74FY6
A	9	PRO	-	EXPRESSION TAG	UNP Q74FY6
A	10	GLN	-	EXPRESSION TAG	UNP Q74FY6
A	11	PHE	-	EXPRESSION TAG	UNP Q74FY6
A	12	GLU	-	EXPRESSION TAG	UNP Q74FY6
A	13	LYS	-	EXPRESSION TAG	UNP Q74FY6
A	14	GLY	-	EXPRESSION TAG	UNP Q74FY6
A	15	ALA	-	EXPRESSION TAG	UNP Q74FY6
A	16	GLU	-	EXPRESSION TAG	UNP Q74FY6
A	17	THR	-	EXPRESSION TAG	UNP Q74FY6
A	18	ALA	-	EXPRESSION TAG	UNP Q74FY6
A	19	VAL	-	EXPRESSION TAG	UNP Q74FY6
A	20	PRO	-	EXPRESSION TAG	UNP Q74FY6
A	21	ASN	-	EXPRESSION TAG	UNP Q74FY6
A	22	SER	-	EXPRESSION TAG	UNP Q74FY6
B	6	TRP	-	EXPRESSION TAG	UNP Q74FY6
B	7	SER	-	EXPRESSION TAG	UNP Q74FY6
B	8	HIS	-	EXPRESSION TAG	UNP Q74FY6
B	9	PRO	-	EXPRESSION TAG	UNP Q74FY6
B	10	GLN	-	EXPRESSION TAG	UNP Q74FY6
B	11	PHE	-	EXPRESSION TAG	UNP Q74FY6
B	12	GLU	-	EXPRESSION TAG	UNP Q74FY6
B	13	LYS	-	EXPRESSION TAG	UNP Q74FY6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	14	GLY	-	EXPRESSION TAG	UNP Q74FY6
B	15	ALA	-	EXPRESSION TAG	UNP Q74FY6
B	16	GLU	-	EXPRESSION TAG	UNP Q74FY6
B	17	THR	-	EXPRESSION TAG	UNP Q74FY6
B	18	ALA	-	EXPRESSION TAG	UNP Q74FY6
B	19	VAL	-	EXPRESSION TAG	UNP Q74FY6
B	20	PRO	-	EXPRESSION TAG	UNP Q74FY6
B	21	ASN	-	EXPRESSION TAG	UNP Q74FY6
B	22	SER	-	EXPRESSION TAG	UNP Q74FY6

- Molecule 2 is HEME C (three-letter code: HEC) (formula: $C_{34}H_{34}FeN_4O_4$).

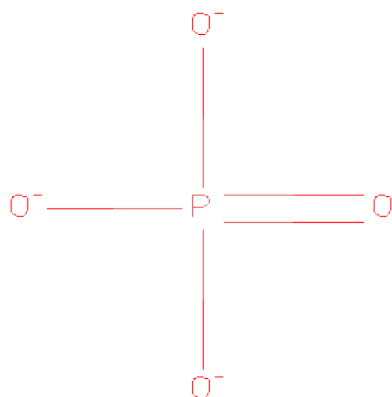


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		
3	A	1	Total	Ca	0	0
			1	1		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



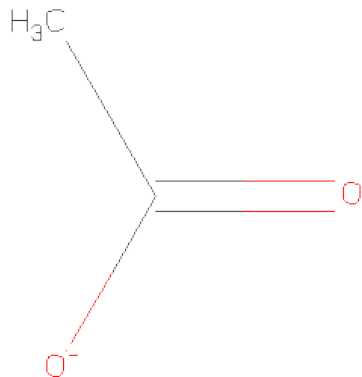
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		

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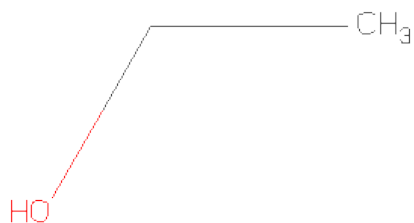
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is ETHANOL (three-letter code: EOH) (formula: C₂H₆O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			3	2	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	344	Total	O	0	0
			344	344		
7	B	285	Total	O	0	0
			285	285		

3 Residue-property plots

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

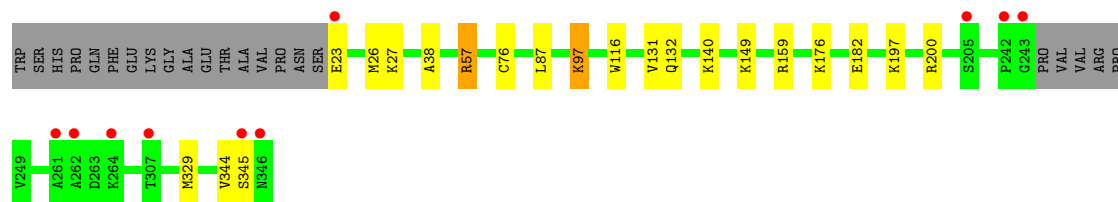
• Molecule 1: CYTOCHROME C551 PEROXIDASE

Chain A: 



• Molecule 1: CYTOCHROME C551 PEROXIDASE

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	118.03Å 118.03Å 242.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	102.06 – 1.84 47.08 – 1.84	Depositor EDS
% Data completeness (in resolution range)	98.0 (102.06-1.84) 98.2 (47.08-1.84)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 1.84Å)	Xtriage
Refinement program	REFMAC 5.6.0113	Depositor
R, R_{free}	0.189 , 0.237 0.193 , 0.238	Depositor DCC
R_{free} test set	4284 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.003	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 85453 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5841	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEC, CA, EOH, PO4, ACT

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	0.96	2/2514 (0.1%)	1.00	13/3413 (0.4%)
1	B	0.94	1/2515 (0.0%)	0.96	2/3414 (0.1%)
All	All	0.95	3/5029 (0.1%)	0.98	15/6827 (0.2%)

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	2
1	B	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	224	SER	CA-CB	6.21	1.62	1.52
1	A	288	TRP	CD2-CE2	5.81	1.48	1.41
1	B	116	TRP	CD2-CE2	5.08	1.47	1.41

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	57	ARG	NE-CZ-NH1	-11.38	114.61	120.30
1	A	268	ARG	NE-CZ-NH2	-7.46	116.57	120.30
1	A	268	ARG	NE-CZ-NH1	5.95	123.28	120.30
1	A	342	ARG	NE-CZ-NH1	-5.52	117.54	120.30
1	A	28	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	B	57	ARG	NE-CZ-NH2	5.29	122.95	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	200[A]	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	A	200[B]	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	A	200[A]	ARG	NE-CZ-NH1	5.15	122.87	120.30
1	A	200[B]	ARG	NE-CZ-NH1	5.15	122.87	120.30
1	A	65	LEU	CB-CG-CD1	5.14	119.74	111.00
1	A	249	VAL	N-CA-C	5.14	124.88	111.00
1	A	329[A]	MET	CG-SD-CE	-5.07	92.09	100.20
1	A	329[B]	MET	CG-SD-CE	-5.07	92.09	100.20
1	A	167	ASP	CB-CG-OD1	5.00	122.80	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	128	LYS	Peptide
1	A	249	VAL	Peptide
1	B	344	VAL	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2447	0	0	15	1
1	B	2449	0	0	17	1
2	A	86	0	0	6	0
2	B	86	0	0	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	55	0	0	3	0
4	B	60	0	0	3	0
5	A	12	0	9	0	0
5	B	12	0	9	1	0
6	A	3	0	6	0	0
7	A	344	0	0	7	1
7	B	285	0	0	8	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5841	0	24	34	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (34) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:76:CYS:SG	2:B:423:HEC:CAC	2.07	1.43
5:B:1362:ACT:H3	7:B:2282:HOH:O	1.50	1.10
1:A:329[B]:MET:SD	7:A:2306:HOH:O	2.24	0.95
1:B:23:GLU:CB	7:B:2001:HOH:O	2.24	0.86
1:B:76:CYS:SG	2:B:423:HEC:C3C	2.65	0.84
1:A:222:CYS:SG	2:A:424:HEC:C3C	2.67	0.81
1:B:23:GLU:CG	1:B:200:ARG:NH2	2.45	0.80
1:B:57:ARG:NH1	7:B:2040:HOH:O	2.15	0.80
1:B:57:ARG:NH1	7:B:2041:HOH:O	2.17	0.78
1:A:76:CYS:SG	2:A:423:HEC:C3C	2.71	0.78
1:B:76:CYS:SG	2:B:423:HEC:CBC	2.74	0.76
1:A:76:CYS:SG	2:A:423:HEC:CBC	2.73	0.75
1:A:167:ASP:OD1	7:A:2195:HOH:O	2.05	0.73
1:A:222:CYS:SG	2:A:424:HEC:CBC	2.75	0.73
1:A:165:GLN:OE1	7:A:2191:HOH:O	2.11	0.68
1:B:26:MET:CE	7:B:2007:HOH:O	2.44	0.64
1:B:149:LYS:NZ	7:B:2148:HOH:O	2.33	0.62
1:A:339:ASP:OD1	7:A:2323:HOH:O	2.17	0.56
1:B:23:GLU:CG	1:B:200:ARG:CZ	2.83	0.56
1:B:159:ARG:NH2	4:B:1350:PO4:O4	2.39	0.55
1:A:97:LYS:N	7:A:2104:HOH:O	2.41	0.53
1:A:346:ASN:N	7:A:2331:HOH:O	2.41	0.53
4:A:1351:PO4:O4	7:A:2278:HOH:O	2.19	0.52
1:B:97:LYS:N	7:B:2088:HOH:O	2.42	0.52
1:A:28:ARG:NH2	1:A:216:ASP:OD1	2.43	0.52
1:B:38:ALA:N	4:B:1357:PO4:O4	2.44	0.50
1:B:76:CYS:CB	2:B:423:HEC:C3C	2.93	0.47
1:A:159:ARG:NE	4:A:1352:PO4:O2	2.48	0.46
1:B:176:LYS:NZ	4:B:1352:PO4:O2	2.49	0.46
1:A:283:HIS:ND1	2:A:424:HEC:O2D	2.50	0.45
1:A:159:ARG:NH2	4:A:1352:PO4:O1	2.50	0.44
1:B:197:LYS:NZ	7:B:2011:HOH:O	2.51	0.43
1:A:76:CYS:CB	2:A:423:HEC:C3C	2.96	0.43
1:B:132:GLN:NE2	1:B:182:GLU:OE1	2.52	0.43

All (2) 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	Distance(Å)	Clash(Å)
1:A:329[A]:MET:SD	1:B:329[A]:MET:SD[8_445]	1.47	0.73
7:A:2081:HOH:O	7:B:2264:HOH:O[8_445]	1.93	0.27

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	319/341 (94%)	307 (96%)	11 (3%)	1 (0%)	50	32
1	B	319/341 (94%)	310 (97%)	9 (3%)	0	100	100
All	All	638/682 (94%)	617 (97%)	20 (3%)	1 (0%)	56	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	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	266/281 (95%)	259 (97%)	7 (3%)	59	40
1	B	266/281 (95%)	260 (98%)	6 (2%)	63	46
All	All	532/562 (95%)	519 (98%)	13 (2%)	63	44

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	LYS
1	A	160[A]	LYS
1	A	160[B]	LYS
1	A	205	SER
1	A	249	VAL
1	A	345	SER
1	A	346	ASN
1	B	27	LYS
1	B	87	LEU
1	B	97	LYS
1	B	131	VAL
1	B	140	LYS
1	B	345	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 36 ligands modelled in this entry, 2 are monoatomic - leaving 34 for Mogul analysis.

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
4	PO4	A	1348	-	4,4,4	0.27	0	6,6,6	0.32	0
4	PO4	A	1349	-	4,4,4	0.40	0	6,6,6	0.30	0
4	PO4	A	1350	-	4,4,4	0.33	0	6,6,6	0.38	0
4	PO4	A	1351	-	4,4,4	0.27	0	6,6,6	0.33	0
4	PO4	A	1352	-	4,4,4	0.21	0	6,6,6	0.32	0
4	PO4	A	1353	-	4,4,4	0.31	0	6,6,6	0.32	0
4	PO4	A	1354	-	4,4,4	0.18	0	6,6,6	0.33	0
4	PO4	A	1355	-	4,4,4	0.34	0	6,6,6	0.32	0
4	PO4	A	1356	-	4,4,4	0.29	0	6,6,6	0.37	0
4	PO4	A	1357	-	4,4,4	0.49	0	6,6,6	0.38	0
5	ACT	A	1358	-	1,3,3	1.58	0	0,3,3	0.00	-
5	ACT	A	1359	-	1,3,3	1.34	0	0,3,3	0.00	-
5	ACT	A	1360	-	1,3,3	2.61	1 (100%)	0,3,3	0.00	-
6	EOH	A	1361	-	2,2,2	0.36	0	1,1,1	0.66	0
4	PO4	A	1362	-	4,4,4	0.30	0	6,6,6	0.34	0
2	HEC	A	423	1	50,50,50	2.37	10 (20%)	56,82,82	3.04	31 (55%)
2	HEC	A	424	1	50,50,50	2.55	13 (26%)	56,82,82	2.89	26 (46%)
4	PO4	B	1348	-	4,4,4	0.24	0	6,6,6	0.32	0
4	PO4	B	1349	-	4,4,4	0.28	0	6,6,6	0.30	0
4	PO4	B	1350	-	4,4,4	0.20	0	6,6,6	0.32	0
4	PO4	B	1351	-	4,4,4	0.48	0	6,6,6	0.32	0
4	PO4	B	1352	-	4,4,4	0.35	0	6,6,6	0.38	0
4	PO4	B	1353	-	4,4,4	0.40	0	6,6,6	0.30	0
4	PO4	B	1354	-	4,4,4	0.39	0	6,6,6	0.35	0
4	PO4	B	1355	-	4,4,4	0.19	0	6,6,6	0.34	0
4	PO4	B	1356	-	4,4,4	0.32	0	6,6,6	0.33	0
4	PO4	B	1357	-	4,4,4	0.46	0	6,6,6	0.28	0
4	PO4	B	1358	-	4,4,4	0.31	0	6,6,6	0.35	0
4	PO4	B	1359	-	4,4,4	0.37	0	6,6,6	0.34	0
5	ACT	B	1360	-	1,3,3	2.85	1 (100%)	0,3,3	0.00	-
5	ACT	B	1361	-	1,3,3	2.12	1 (100%)	0,3,3	0.00	-
5	ACT	B	1362	-	1,3,3	0.86	0	0,3,3	0.00	-
2	HEC	B	423	1	50,50,50	2.32	10 (20%)	56,82,82	3.25	36 (64%)
2	HEC	B	424	1	50,50,50	2.45	12 (24%)	56,82,82	2.74	25 (44%)

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
4	PO4	A	1348	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1349	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1350	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1351	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1352	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1353	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1354	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1355	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1356	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1357	-	-	0/0/0/0	0/0/0/0
5	ACT	A	1358	-	-	0/0/0/0	0/0/0/0
5	ACT	A	1359	-	-	0/0/0/0	0/0/0/0
5	ACT	A	1360	-	-	0/0/0/0	0/0/0/0
6	EOH	A	1361	-	-	0/0/0/0	0/0/0/0
4	PO4	A	1362	-	-	0/0/0/0	0/0/0/0
2	HEC	A	423	1	-	0/10/54/54	0/0/8/8
2	HEC	A	424	1	-	0/10/54/54	0/0/8/8
4	PO4	B	1348	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1349	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1350	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1351	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1352	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1353	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1354	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1355	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1356	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1357	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1358	-	-	0/0/0/0	0/0/0/0
4	PO4	B	1359	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1360	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1361	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1362	-	-	0/0/0/0	0/0/0/0
2	HEC	B	423	1	-	0/10/54/54	0/0/8/8
2	HEC	B	424	1	-	0/10/54/54	0/0/8/8

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	424	HEC	C3C-CAC	8.50	1.52	1.35
2	B	424	HEC	FE-NB	8.05	2.27	1.92
2	A	423	HEC	C3C-CAC	7.87	1.51	1.35
2	B	423	HEC	C1D-C2D	7.26	1.49	1.40
2	B	424	HEC	C3C-CAC	7.22	1.50	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	424	HEC	FE-NB	6.80	2.21	1.92
2	B	423	HEC	FE-NB	6.72	2.21	1.92
2	B	423	HEC	C3B-CAB	6.60	1.49	1.35
2	A	423	HEC	C3B-CAB	6.52	1.48	1.35
2	B	424	HEC	C3B-CAB	6.47	1.48	1.35
2	A	424	HEC	C1D-C2D	6.44	1.48	1.40
2	A	424	HEC	C3B-CAB	6.44	1.48	1.35
2	A	423	HEC	C1D-C2D	6.26	1.48	1.40
2	B	423	HEC	C3C-CAC	6.26	1.48	1.35
2	A	423	HEC	FE-NB	6.14	2.18	1.92
2	B	424	HEC	C1D-C2D	5.46	1.47	1.40
2	B	424	HEC	C3B-C4B	5.26	1.48	1.41
2	A	424	HEC	C1B-C2B	5.03	1.46	1.40
2	A	423	HEC	C4B-NB	-4.68	1.30	1.37
2	A	423	HEC	C3B-C4B	4.54	1.47	1.41
2	A	424	HEC	C3B-C4B	4.18	1.46	1.41
2	B	423	HEC	C3B-C4B	3.91	1.46	1.41
2	A	423	HEC	FE-NC	3.57	2.07	1.92
2	B	424	HEC	FE-NC	3.53	2.07	1.92
2	B	423	HEC	C4B-NB	-3.45	1.32	1.37
2	B	424	HEC	C1B-C2B	3.36	1.44	1.40
2	A	424	HEC	FE-NC	3.35	2.06	1.92
2	A	424	HEC	CBC-CAC	-3.28	1.35	1.49
2	B	424	HEC	CBC-CAC	-3.10	1.36	1.49
2	B	423	HEC	FE-NC	3.09	2.05	1.92
2	B	423	HEC	CBC-CAC	-2.95	1.36	1.49
5	B	1360	ACT	CH3-C	2.85	1.52	1.48
2	A	424	HEC	C4D-C3D	2.83	1.48	1.43
2	A	424	HEC	FE-NA	2.62	2.03	1.92
5	A	1360	ACT	CH3-C	2.61	1.52	1.48
2	A	424	HEC	C1C-C2C	2.59	1.43	1.40
2	B	424	HEC	C4D-C3D	2.51	1.47	1.43
2	B	423	HEC	CBB-CAB	-2.50	1.38	1.49
2	B	423	HEC	C4D-C3D	2.42	1.47	1.43
2	A	424	HEC	C4B-NB	-2.41	1.33	1.37
2	B	424	HEC	C1B-NB	-2.26	1.32	1.36
2	A	423	HEC	CBB-CAB	-2.23	1.39	1.49
2	A	423	HEC	CBD-CGD	2.14	1.56	1.50
5	B	1361	ACT	CH3-C	2.12	1.51	1.48
2	A	424	HEC	CBB-CAB	-2.11	1.40	1.49
2	B	424	HEC	CBB-CAB	-2.09	1.40	1.49
2	A	423	HEC	C4C-NC	-2.05	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	424	HEC	C3C-C4C	2.01	1.44	1.41

All (118) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	423	HEC	C4B-NB-C1B	8.57	118.05	106.76
2	B	423	HEC	CBD-CAD-C3D	-7.90	98.77	112.69
2	A	424	HEC	C4B-NB-C1B	7.60	116.77	106.76
2	A	423	HEC	C4B-NB-C1B	7.46	116.60	106.76
2	A	424	HEC	CBB-CAB-C3B	-6.73	109.49	128.44
2	B	424	HEC	C1A-CHA-C4D	-6.66	118.71	127.47
2	A	424	HEC	C4C-CHD-C1D	-6.65	118.73	127.47
2	B	423	HEC	C2B-C1B-NB	-6.61	104.42	109.41
2	B	424	HEC	C4B-NB-C1B	6.52	115.35	106.76
2	B	423	HEC	C3B-C4B-NB	-6.40	102.28	111.52
2	A	423	HEC	CBB-CAB-C3B	-6.39	110.44	128.44
2	A	423	HEC	C1A-CHA-C4D	-6.11	119.43	127.47
2	B	423	HEC	CBB-CAB-C3B	-6.05	111.41	128.44
2	A	424	HEC	C2B-C1B-NB	-6.00	104.88	109.41
2	B	424	HEC	CBB-CAB-C3B	-5.83	112.03	128.44
2	B	423	HEC	C1A-CHA-C4D	-5.80	119.84	127.47
2	B	424	HEC	C4C-CHD-C1D	-5.79	119.85	127.47
2	A	423	HEC	C4C-CHD-C1D	-5.77	119.88	127.47
2	A	423	HEC	C3B-C4B-NB	-5.72	103.26	111.52
2	B	423	HEC	C4C-CHD-C1D	-5.69	119.98	127.47
2	B	424	HEC	C3B-C4B-NB	-5.55	103.51	111.52
2	A	424	HEC	C1A-CHA-C4D	-5.46	120.29	127.47
2	A	423	HEC	C2B-C1B-NB	-5.12	105.54	109.41
2	A	423	HEC	CBD-CAD-C3D	-5.06	103.77	112.69
2	A	424	HEC	C3B-C4B-NB	-5.00	104.29	111.52
2	B	423	HEC	C2A-C1A-NA	-4.54	106.23	109.64
2	A	424	HEC	CMB-C2B-C3B	4.50	130.68	125.72
2	A	424	HEC	CBD-CAD-C3D	-4.46	104.84	112.69
2	B	423	HEC	C4A-CHB-C1B	-4.27	121.85	127.47
2	B	424	HEC	C3D-C4D-CHA	-4.14	118.15	126.00
2	A	423	HEC	CMB-C2B-C3B	4.11	130.26	125.72
2	B	424	HEC	CBD-CAD-C3D	-3.95	105.73	112.69
2	A	424	HEC	CMB-C2B-C1B	-3.94	122.56	128.62
2	B	423	HEC	C3D-C4D-CHA	-3.85	118.69	126.00
2	A	423	HEC	CMB-C2B-C1B	-3.83	122.73	128.62
2	B	424	HEC	C2B-C1B-NB	-3.81	106.53	109.41
2	A	423	HEC	C2D-C1D-ND	3.77	112.26	109.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	424	HEC	C4C-NC-C1C	3.72	111.67	106.76
2	B	424	HEC	C3C-C4C-NC	-3.72	106.14	111.52
2	B	424	HEC	C2D-C1D-ND	3.59	112.12	109.41
2	A	423	HEC	C4A-CHB-C1B	-3.58	122.76	127.47
2	A	423	HEC	C4B-CHC-C1C	-3.58	122.76	127.47
2	B	423	HEC	CMB-C2B-C1B	-3.55	123.15	128.62
2	A	423	HEC	O2A-CGA-CBA	3.46	126.45	114.22
2	B	424	HEC	C2A-C1A-NA	-3.45	107.05	109.64
2	A	424	HEC	C2D-C1D-CHD	-3.39	119.57	126.00
2	A	423	HEC	C2A-C1A-NA	-3.39	107.09	109.64
2	B	424	HEC	CHA-C4D-ND	3.39	130.23	124.58
2	A	424	HEC	C3D-C4D-CHA	-3.38	119.58	126.00
2	A	423	HEC	C4C-NC-C1C	3.38	111.21	106.76
2	B	423	HEC	C4D-C3D-C2D	-3.37	103.43	106.92
2	A	424	HEC	C4C-NC-C1C	3.37	111.20	106.76
2	B	423	HEC	C3C-C4C-NC	-3.36	106.67	111.52
2	A	423	HEC	C1D-C2D-C3D	-3.35	104.67	107.00
2	B	424	HEC	O2D-CGD-CBD	3.27	125.76	114.22
2	A	423	HEC	CMC-C2C-C1C	-3.24	123.63	128.62
2	A	424	HEC	O2D-CGD-CBD	3.20	125.54	114.22
2	A	423	HEC	C4A-C3A-C2A	3.19	108.92	106.89
2	B	423	HEC	C3D-C4D-ND	3.19	114.16	109.73
2	A	424	HEC	CBA-CAA-C2A	-3.16	106.45	112.35
2	A	423	HEC	CBA-CAA-C2A	-3.13	106.49	112.35
2	B	424	HEC	CBC-CAC-C3C	-3.10	119.71	128.44
2	A	424	HEC	CMD-C2D-C1D	-3.07	123.90	128.62
2	A	423	HEC	O2D-CGD-O1D	-3.05	115.54	123.30
2	B	423	HEC	CMA-C3A-C2A	3.03	130.66	124.94
2	B	423	HEC	CBC-CAC-C3C	-3.03	119.90	128.44
2	B	423	HEC	C2D-C1D-ND	3.02	111.69	109.41
2	B	423	HEC	CMC-C2C-C3C	2.99	129.02	125.72
2	B	423	HEC	CMB-C2B-C3B	2.97	129.00	125.72
2	B	423	HEC	CMC-C2C-C1C	-2.96	124.07	128.62
2	A	423	HEC	C2C-C1C-NC	-2.91	107.22	109.41
2	A	423	HEC	O1A-CGA-CBA	-2.86	113.19	123.03
2	B	424	HEC	CBA-CAA-C2A	-2.84	107.05	112.35
2	B	424	HEC	C2C-C1C-NC	-2.83	107.27	109.41
2	B	423	HEC	C4A-NA-C1A	2.79	109.76	105.58
2	A	424	HEC	CHD-C1D-ND	2.76	129.19	124.58
2	A	423	HEC	CAA-C2A-C1A	2.76	129.66	124.67
2	B	423	HEC	C4C-NC-C1C	2.76	110.40	106.76
2	A	424	HEC	O2D-CGD-O1D	-2.74	116.32	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	424	HEC	CMC-C2C-C1C	-2.64	124.57	128.62
2	A	423	HEC	C3D-C4D-CHA	-2.63	121.01	126.00
2	A	423	HEC	C2D-C1D-CHD	-2.63	121.02	126.00
2	A	423	HEC	C4C-C3C-C2C	-2.62	105.01	106.68
2	A	424	HEC	C4A-CHB-C1B	-2.61	124.04	127.47
2	B	423	HEC	CAA-CBA-CGA	-2.60	105.10	113.47
2	B	423	HEC	C4B-CHC-C1C	-2.60	124.05	127.47
2	B	424	HEC	C4A-CHB-C1B	-2.53	124.15	127.47
2	B	423	HEC	C3A-C4A-NA	-2.52	107.51	109.41
2	B	424	HEC	O2A-CGA-CBA	2.49	123.02	114.22
2	B	423	HEC	O2D-CGD-CBD	2.47	122.96	114.22
2	A	423	HEC	CAA-CBA-CGA	-2.47	105.53	113.47
2	A	423	HEC	C3C-C4C-NC	-2.47	107.95	111.52
2	B	423	HEC	CMA-C3A-C4A	-2.43	121.36	126.16
2	A	424	HEC	C2D-C1D-ND	2.42	111.24	109.41
2	B	424	HEC	C2D-C1D-CHD	-2.40	121.45	126.00
2	B	424	HEC	CMC-C2C-C1C	-2.38	124.96	128.62
2	A	423	HEC	C4D-ND-C1D	-2.35	103.66	106.76
2	A	424	HEC	C2C-C1C-NC	-2.35	107.64	109.41
2	B	424	HEC	C1A-C2A-C3A	2.31	108.51	106.69
2	A	424	HEC	CHA-C4D-ND	2.30	128.42	124.58
2	B	423	HEC	CHB-C1B-NB	2.28	128.39	124.58
2	A	424	HEC	C3C-C4C-NC	-2.28	108.23	111.52
2	A	424	HEC	CMC-C2C-C3C	2.25	128.21	125.72
2	B	423	HEC	C1A-C2A-C3A	2.24	108.45	106.69
2	B	424	HEC	C4A-NA-C1A	2.23	108.92	105.58
2	B	423	HEC	C4C-C3C-C2C	2.22	108.10	106.68
2	B	423	HEC	O1A-CGA-CBA	-2.22	115.39	123.03
2	A	424	HEC	CMA-C3A-C2A	2.15	129.00	124.94
2	B	423	HEC	C2D-C1D-CHD	-2.13	121.96	126.00
2	A	423	HEC	CAA-C2A-C3A	-2.13	122.93	129.00
2	B	424	HEC	O1D-CGD-CBD	-2.12	115.75	123.03
2	B	423	HEC	O2A-CGA-CBA	2.10	121.66	114.22
2	B	423	HEC	CAA-C2A-C3A	-2.09	123.03	129.00
2	B	423	HEC	C2C-C1C-NC	-2.09	107.83	109.41
2	A	424	HEC	C4D-ND-C1D	-2.09	104.02	106.76
2	B	423	HEC	CAA-C2A-C1A	2.08	128.42	124.67
2	B	424	HEC	CMC-C2C-C3C	2.08	128.01	125.72
2	A	423	HEC	CMC-C2C-C3C	2.03	127.96	125.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	320/341 (93%)	-0.11	9 (2%) 50 46	21, 31, 56, 98	0
1	B	319/341 (93%)	-0.08	10 (3%) 47 42	22, 32, 54, 124	0
All	All	639/682 (93%)	-0.10	19 (2%) 50 44	21, 32, 55, 124	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	346	ASN	9.4
1	A	249	VAL	5.9
1	B	345	SER	5.5
1	B	23	GLU	4.8
1	A	248	PRO	3.7
1	A	242	PRO	3.4
1	A	346	ASN	3.4
1	B	242	PRO	2.8
1	A	253	GLY	2.5
1	A	252	THR	2.5
1	B	243	GLY	2.4
1	B	262	ALA	2.4
1	A	250	ASP	2.2
1	B	205	SER	2.1
1	A	251	ASP	2.1
1	B	264	LYS	2.1
1	B	261	ALA	2.1
1	A	243	GLY	2.1
1	B	307	THR	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	PO4	A	1348	5/5	0.23	10.80	58,60,77,90	0
4	PO4	A	1355	5/5	0.20	10.73	60,67,81,84	0
4	PO4	A	1352	5/5	0.25	9.29	66,73,81,82	0
5	ACT	B	1362	4/4	0.20	6.47	43,45,51,68	0
4	PO4	B	1359	5/5	0.18	6.13	53,60,68,80	0
4	PO4	B	1358	5/5	0.22	5.25	55,61,74,77	0
4	PO4	B	1350	5/5	0.23	4.45	74,79,86,89	0
5	ACT	A	1360	4/4	0.20	3.77	54,61,67,72	0
4	PO4	A	1362	5/5	0.21	3.32	68,78,83,97	0
4	PO4	A	1351	5/5	0.25	3.30	67,74,96,99	0
5	ACT	B	1360	4/4	0.23	2.57	50,59,61,66	0
4	PO4	B	1357	5/5	0.21	2.49	60,81,98,103	0
5	ACT	B	1361	4/4	0.25	2.37	78,82,83,89	0
4	PO4	A	1349	5/5	0.21	2.32	65,71,89,102	0
4	PO4	B	1351	5/5	0.13	1.36	53,64,66,69	0
4	PO4	B	1356	5/5	0.14	1.25	68,91,97,97	0
5	ACT	A	1358	4/4	0.10	1.00	58,58,66,69	0
4	PO4	A	1353	5/5	0.24	0.93	42,56,63,68	0
4	PO4	B	1352	5/5	0.12	0.82	62,63,67,71	0
4	PO4	A	1357	5/5	0.13	0.76	43,53,60,65	0
4	PO4	B	1348	5/5	0.27	0.68	81,89,105,107	0
2	HEC	A	423	43/43	0.13	0.59	19,22,29,31	0
4	PO4	B	1355	5/5	0.12	0.51	66,68,75,81	0
4	PO4	B	1353	5/5	0.16	0.49	69,74,84,86	0
4	PO4	A	1350	5/5	0.13	0.37	61,71,86,87	0
2	HEC	B	423	43/43	0.11	0.32	21,23,31,40	0
2	HEC	A	424	43/43	0.10	0.20	23,27,31,32	0
4	PO4	B	1354	5/5	0.10	0.08	51,59,60,62	0
4	PO4	A	1356	5/5	0.10	0.00	63,63,66,73	0
2	HEC	B	424	43/43	0.09	-0.07	23,29,33,35	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	PO4	B	1349	5/5	0.12	-0.34	67,70,79,80	0
3	CA	A	425	1/1	0.08	-0.55	25,25,25,25	0
4	PO4	A	1354	5/5	0.12	-0.67	70,70,81,93	0
3	CA	B	425	1/1	0.10	-0.67	25,25,25,25	0
5	ACT	A	1359	4/4	0.09	-1.26	56,58,59,61	0
6	EOH	A	1361	3/3	0.11	-1.85	50,50,54,55	0

6.5 Other polymers ⓘ

There are no such residues in this entry.