



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:06 PM GMT

PDB ID : 4BIZ  
Title : Crystal structure of CpxAHDC (orthorhombic form 2)  
Authors : Mechaly, A.E.; Sassoon, N.; Betton, J.M.; Alzari, P.M.  
Deposited on : 2013-04-13  
Resolution : 2.65 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

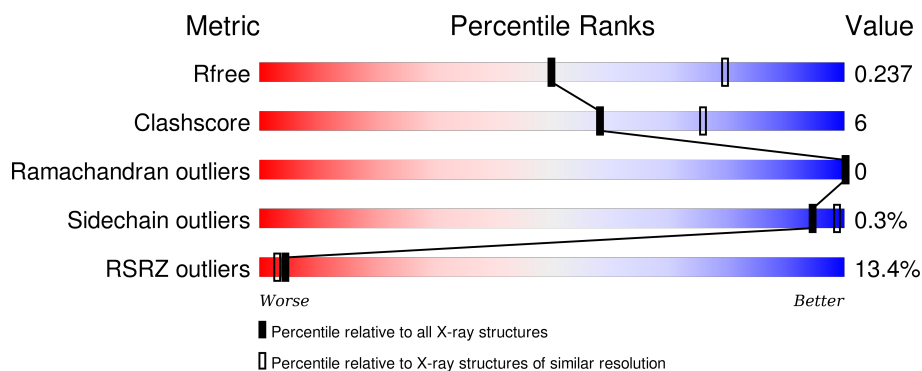
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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))
$R_{free}$	91344	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	298	<div> <div>7%</div> <div>67%</div> <div>11%</div> <div>22%</div> </div>
1	B	298	<div> <div>13%</div> <div>65%</div> <div>11%</div> <div>24%</div> </div>
1	C	298	<div> <div>6%</div> <div>73%</div> <div>7%</div> <div>20%</div> </div>
1	D	298	<div> <div>17%</div> <div>62%</div> <div>9%</div> <div>28%</div> </div>
1	E	298	<div> <div>8%</div> <div>67%</div> <div>9%</div> <div>23%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	298	<p>10% 69% 6% 24%</p>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11128 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 SENSOR PROTEIN CPXA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1838	1146	329	357	6			
1	B	227	Total	C	N	O	S	0	0	0
			1787	1120	317	344	6			
1	C	239	Total	C	N	O	S	0	0	0
			1881	1174	335	366	6			
1	D	214	Total	C	N	O	S	0	0	0
			1674	1049	297	322	6			
1	E	229	Total	C	N	O	S	0	0	0
			1809	1130	322	351	6			
1	F	225	Total	C	N	O	S	0	0	0
			1783	1116	318	343	6			

There are 174 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	160	MET	-	EXPRESSION TAG	UNP P0AE82
A	161	GLY	-	EXPRESSION TAG	UNP P0AE82
A	162	SER	-	EXPRESSION TAG	UNP P0AE82
A	163	SER	-	EXPRESSION TAG	UNP P0AE82
A	164	HIS	-	EXPRESSION TAG	UNP P0AE82
A	165	HIS	-	EXPRESSION TAG	UNP P0AE82
A	166	HIS	-	EXPRESSION TAG	UNP P0AE82
A	167	HIS	-	EXPRESSION TAG	UNP P0AE82
A	168	HIS	-	EXPRESSION TAG	UNP P0AE82
A	169	HIS	-	EXPRESSION TAG	UNP P0AE82
A	170	SER	-	EXPRESSION TAG	UNP P0AE82
A	171	SER	-	EXPRESSION TAG	UNP P0AE82
A	172	GLY	-	EXPRESSION TAG	UNP P0AE82
A	173	LEU	-	EXPRESSION TAG	UNP P0AE82
A	174	VAL	-	EXPRESSION TAG	UNP P0AE82
A	175	PRO	-	EXPRESSION TAG	UNP P0AE82
A	176	ARG	-	EXPRESSION TAG	UNP P0AE82

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Chain	Residue	Modelled	Actual	Comment	Reference
A	177	GLY	-	EXPRESSION TAG	UNP P0AE82
A	178	SER	-	EXPRESSION TAG	UNP P0AE82
A	179	HIS	-	EXPRESSION TAG	UNP P0AE82
A	180	MET	-	EXPRESSION TAG	UNP P0AE82
A	181	GLU	-	EXPRESSION TAG	UNP P0AE82
A	182	ASN	-	EXPRESSION TAG	UNP P0AE82
A	183	LEU	-	EXPRESSION TAG	UNP P0AE82
A	184	TYR	-	EXPRESSION TAG	UNP P0AE82
A	185	PHE	-	EXPRESSION TAG	UNP P0AE82
A	186	GLN	-	EXPRESSION TAG	UNP P0AE82
A	187	GLY	-	EXPRESSION TAG	UNP P0AE82
A	248	GLU	HIS	ENGINEERED MUTATION	UNP P0AE82
B	160	MET	-	EXPRESSION TAG	UNP P0AE82
B	161	GLY	-	EXPRESSION TAG	UNP P0AE82
B	162	SER	-	EXPRESSION TAG	UNP P0AE82
B	163	SER	-	EXPRESSION TAG	UNP P0AE82
B	164	HIS	-	EXPRESSION TAG	UNP P0AE82
B	165	HIS	-	EXPRESSION TAG	UNP P0AE82
B	166	HIS	-	EXPRESSION TAG	UNP P0AE82
B	167	HIS	-	EXPRESSION TAG	UNP P0AE82
B	168	HIS	-	EXPRESSION TAG	UNP P0AE82
B	169	HIS	-	EXPRESSION TAG	UNP P0AE82
B	170	SER	-	EXPRESSION TAG	UNP P0AE82
B	171	SER	-	EXPRESSION TAG	UNP P0AE82
B	172	GLY	-	EXPRESSION TAG	UNP P0AE82
B	173	LEU	-	EXPRESSION TAG	UNP P0AE82
B	174	VAL	-	EXPRESSION TAG	UNP P0AE82
B	175	PRO	-	EXPRESSION TAG	UNP P0AE82
B	176	ARG	-	EXPRESSION TAG	UNP P0AE82
B	177	GLY	-	EXPRESSION TAG	UNP P0AE82
B	178	SER	-	EXPRESSION TAG	UNP P0AE82
B	179	HIS	-	EXPRESSION TAG	UNP P0AE82
B	180	MET	-	EXPRESSION TAG	UNP P0AE82
B	181	GLU	-	EXPRESSION TAG	UNP P0AE82
B	182	ASN	-	EXPRESSION TAG	UNP P0AE82
B	183	LEU	-	EXPRESSION TAG	UNP P0AE82
B	184	TYR	-	EXPRESSION TAG	UNP P0AE82
B	185	PHE	-	EXPRESSION TAG	UNP P0AE82
B	186	GLN	-	EXPRESSION TAG	UNP P0AE82
B	187	GLY	-	EXPRESSION TAG	UNP P0AE82
B	248	GLU	HIS	ENGINEERED MUTATION	UNP P0AE82
C	160	MET	-	EXPRESSION TAG	UNP P0AE82

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Chain	Residue	Modelled	Actual	Comment	Reference
C	161	GLY	-	EXPRESSION TAG	UNP P0AE82
C	162	SER	-	EXPRESSION TAG	UNP P0AE82
C	163	SER	-	EXPRESSION TAG	UNP P0AE82
C	164	HIS	-	EXPRESSION TAG	UNP P0AE82
C	165	HIS	-	EXPRESSION TAG	UNP P0AE82
C	166	HIS	-	EXPRESSION TAG	UNP P0AE82
C	167	HIS	-	EXPRESSION TAG	UNP P0AE82
C	168	HIS	-	EXPRESSION TAG	UNP P0AE82
C	169	HIS	-	EXPRESSION TAG	UNP P0AE82
C	170	SER	-	EXPRESSION TAG	UNP P0AE82
C	171	SER	-	EXPRESSION TAG	UNP P0AE82
C	172	GLY	-	EXPRESSION TAG	UNP P0AE82
C	173	LEU	-	EXPRESSION TAG	UNP P0AE82
C	174	VAL	-	EXPRESSION TAG	UNP P0AE82
C	175	PRO	-	EXPRESSION TAG	UNP P0AE82
C	176	ARG	-	EXPRESSION TAG	UNP P0AE82
C	177	GLY	-	EXPRESSION TAG	UNP P0AE82
C	178	SER	-	EXPRESSION TAG	UNP P0AE82
C	179	HIS	-	EXPRESSION TAG	UNP P0AE82
C	180	MET	-	EXPRESSION TAG	UNP P0AE82
C	181	GLU	-	EXPRESSION TAG	UNP P0AE82
C	182	ASN	-	EXPRESSION TAG	UNP P0AE82
C	183	LEU	-	EXPRESSION TAG	UNP P0AE82
C	184	TYR	-	EXPRESSION TAG	UNP P0AE82
C	185	PHE	-	EXPRESSION TAG	UNP P0AE82
C	186	GLN	-	EXPRESSION TAG	UNP P0AE82
C	187	GLY	-	EXPRESSION TAG	UNP P0AE82
C	248	GLU	HIS	ENGINEERED MUTATION	UNP P0AE82
D	160	MET	-	EXPRESSION TAG	UNP P0AE82
D	161	GLY	-	EXPRESSION TAG	UNP P0AE82
D	162	SER	-	EXPRESSION TAG	UNP P0AE82
D	163	SER	-	EXPRESSION TAG	UNP P0AE82
D	164	HIS	-	EXPRESSION TAG	UNP P0AE82
D	165	HIS	-	EXPRESSION TAG	UNP P0AE82
D	166	HIS	-	EXPRESSION TAG	UNP P0AE82
D	167	HIS	-	EXPRESSION TAG	UNP P0AE82
D	168	HIS	-	EXPRESSION TAG	UNP P0AE82
D	169	HIS	-	EXPRESSION TAG	UNP P0AE82
D	170	SER	-	EXPRESSION TAG	UNP P0AE82
D	171	SER	-	EXPRESSION TAG	UNP P0AE82
D	172	GLY	-	EXPRESSION TAG	UNP P0AE82
D	173	LEU	-	EXPRESSION TAG	UNP P0AE82

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Chain	Residue	Modelled	Actual	Comment	Reference
D	174	VAL	-	EXPRESSION TAG	UNP P0AE82
D	175	PRO	-	EXPRESSION TAG	UNP P0AE82
D	176	ARG	-	EXPRESSION TAG	UNP P0AE82
D	177	GLY	-	EXPRESSION TAG	UNP P0AE82
D	178	SER	-	EXPRESSION TAG	UNP P0AE82
D	179	HIS	-	EXPRESSION TAG	UNP P0AE82
D	180	MET	-	EXPRESSION TAG	UNP P0AE82
D	181	GLU	-	EXPRESSION TAG	UNP P0AE82
D	182	ASN	-	EXPRESSION TAG	UNP P0AE82
D	183	LEU	-	EXPRESSION TAG	UNP P0AE82
D	184	TYR	-	EXPRESSION TAG	UNP P0AE82
D	185	PHE	-	EXPRESSION TAG	UNP P0AE82
D	186	GLN	-	EXPRESSION TAG	UNP P0AE82
D	187	GLY	-	EXPRESSION TAG	UNP P0AE82
D	248	GLU	HIS	ENGINEERED MUTATION	UNP P0AE82
E	160	MET	-	EXPRESSION TAG	UNP P0AE82
E	161	GLY	-	EXPRESSION TAG	UNP P0AE82
E	162	SER	-	EXPRESSION TAG	UNP P0AE82
E	163	SER	-	EXPRESSION TAG	UNP P0AE82
E	164	HIS	-	EXPRESSION TAG	UNP P0AE82
E	165	HIS	-	EXPRESSION TAG	UNP P0AE82
E	166	HIS	-	EXPRESSION TAG	UNP P0AE82
E	167	HIS	-	EXPRESSION TAG	UNP P0AE82
E	168	HIS	-	EXPRESSION TAG	UNP P0AE82
E	169	HIS	-	EXPRESSION TAG	UNP P0AE82
E	170	SER	-	EXPRESSION TAG	UNP P0AE82
E	171	SER	-	EXPRESSION TAG	UNP P0AE82
E	172	GLY	-	EXPRESSION TAG	UNP P0AE82
E	173	LEU	-	EXPRESSION TAG	UNP P0AE82
E	174	VAL	-	EXPRESSION TAG	UNP P0AE82
E	175	PRO	-	EXPRESSION TAG	UNP P0AE82
E	176	ARG	-	EXPRESSION TAG	UNP P0AE82
E	177	GLY	-	EXPRESSION TAG	UNP P0AE82
E	178	SER	-	EXPRESSION TAG	UNP P0AE82
E	179	HIS	-	EXPRESSION TAG	UNP P0AE82
E	180	MET	-	EXPRESSION TAG	UNP P0AE82
E	181	GLU	-	EXPRESSION TAG	UNP P0AE82
E	182	ASN	-	EXPRESSION TAG	UNP P0AE82
E	183	LEU	-	EXPRESSION TAG	UNP P0AE82
E	184	TYR	-	EXPRESSION TAG	UNP P0AE82
E	185	PHE	-	EXPRESSION TAG	UNP P0AE82
E	186	GLN	-	EXPRESSION TAG	UNP P0AE82

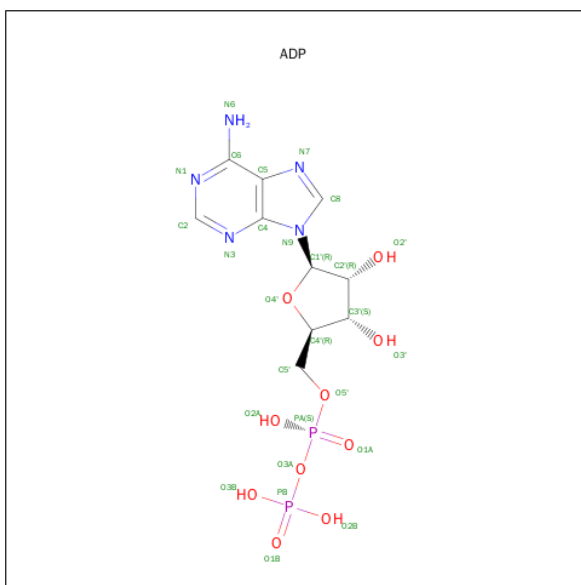
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Chain	Residue	Modelled	Actual	Comment	Reference
E	187	GLY	-	EXPRESSION TAG	UNP P0AE82
E	248	GLU	HIS	ENGINEERED MUTATION	UNP P0AE82
F	160	MET	-	EXPRESSION TAG	UNP P0AE82
F	161	GLY	-	EXPRESSION TAG	UNP P0AE82
F	162	SER	-	EXPRESSION TAG	UNP P0AE82
F	163	SER	-	EXPRESSION TAG	UNP P0AE82
F	164	HIS	-	EXPRESSION TAG	UNP P0AE82
F	165	HIS	-	EXPRESSION TAG	UNP P0AE82
F	166	HIS	-	EXPRESSION TAG	UNP P0AE82
F	167	HIS	-	EXPRESSION TAG	UNP P0AE82
F	168	HIS	-	EXPRESSION TAG	UNP P0AE82
F	169	HIS	-	EXPRESSION TAG	UNP P0AE82
F	170	SER	-	EXPRESSION TAG	UNP P0AE82
F	171	SER	-	EXPRESSION TAG	UNP P0AE82
F	172	GLY	-	EXPRESSION TAG	UNP P0AE82
F	173	LEU	-	EXPRESSION TAG	UNP P0AE82
F	174	VAL	-	EXPRESSION TAG	UNP P0AE82
F	175	PRO	-	EXPRESSION TAG	UNP P0AE82
F	176	ARG	-	EXPRESSION TAG	UNP P0AE82
F	177	GLY	-	EXPRESSION TAG	UNP P0AE82
F	178	SER	-	EXPRESSION TAG	UNP P0AE82
F	179	HIS	-	EXPRESSION TAG	UNP P0AE82
F	180	MET	-	EXPRESSION TAG	UNP P0AE82
F	181	GLU	-	EXPRESSION TAG	UNP P0AE82
F	182	ASN	-	EXPRESSION TAG	UNP P0AE82
F	183	LEU	-	EXPRESSION TAG	UNP P0AE82
F	184	TYR	-	EXPRESSION TAG	UNP P0AE82
F	185	PHE	-	EXPRESSION TAG	UNP P0AE82
F	186	GLN	-	EXPRESSION TAG	UNP P0AE82
F	187	GLY	-	EXPRESSION TAG	UNP P0AE82
F	248	GLU	HIS	ENGINEERED MUTATION	UNP P0AE82

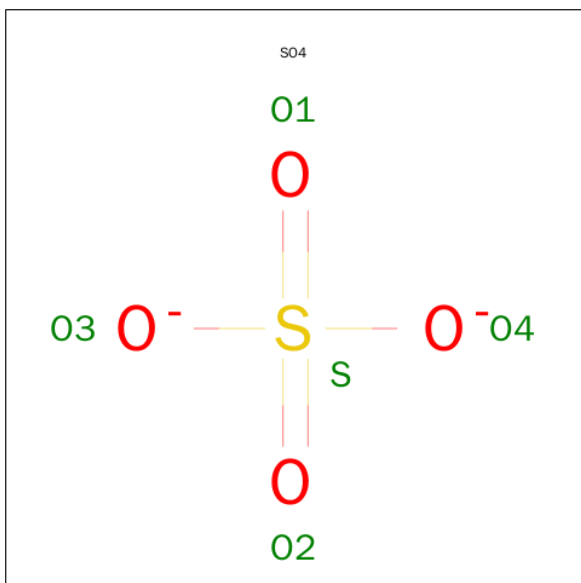
- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0
3	F	1	Total O S 5 4 1	0	0

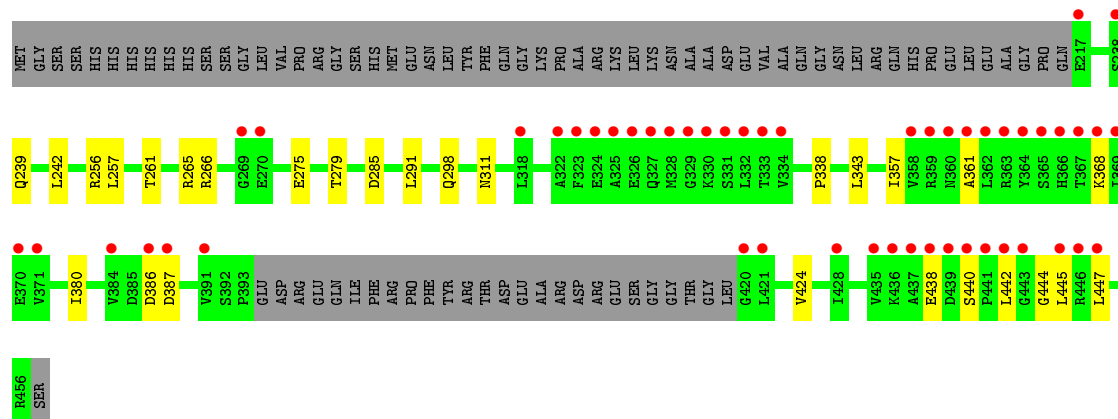
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	46	Total O 46 46	0	0
4	B	36	Total O 36 36	0	0
4	C	57	Total O 57 57	0	0
4	D	23	Total O 23 23	0	0
4	E	28	Total O 28 28	0	0
4	F	28	Total O 28 28	0	0

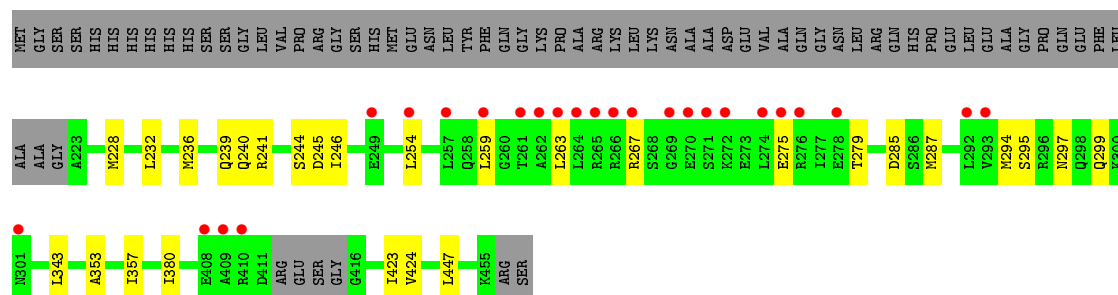


ARG  
SER

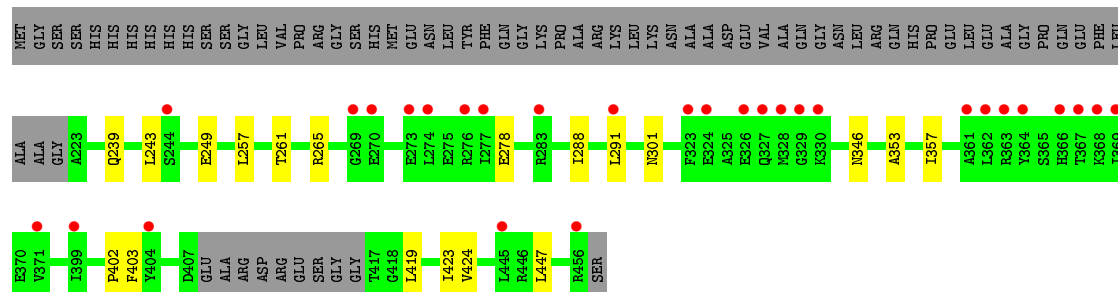
• Molecule 1: SENSOR PROTEIN CPXA



• Molecule 1: SENSOR PROTEIN CPXA



• Molecule 1: SENSOR PROTEIN CPXA



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	120.45Å 125.11Å 159.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.01 – 2.65 49.01 – 2.65	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.01-2.65) 100.0 (49.01-2.65)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.28 (at 2.65Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, $R_{free}$	0.206 , 0.235 0.212 , 0.237	Depositor DCC
$R_{free}$ test set	3571 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	71.2	Xtriage
Anisotropy	0.185	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 69.4	EDS
Estimated twinning fraction	0.014 for k,h,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 70756 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	11128	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, ADP

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	0.50	0/1867	0.67	0/2526
1	B	0.48	0/1816	0.66	0/2459
1	C	0.49	0/1912	0.66	0/2588
1	D	0.49	0/1700	0.67	0/2302
1	E	0.47	0/1838	0.65	0/2488
1	F	0.45	0/1812	0.64	0/2453
All	All	0.48	0/10945	0.66	0/14816

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1838	0	1832	38	0
1	B	1787	0	1791	36	0
1	C	1881	0	1872	25	0
1	D	1674	0	1686	25	0
1	E	1809	0	1805	25	0
1	F	1783	0	1787	20	0
2	A	27	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	27	0	12	0	0
2	E	27	0	12	0	0
2	F	27	0	12	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	E	5	0	0	0	0
3	F	5	0	0	0	0
4	A	46	0	0	0	0
4	B	36	0	0	0	0
4	C	57	0	0	1	0
4	D	23	0	0	0	0
4	E	28	0	0	0	0
4	F	28	0	0	0	0
All	All	11128	0	10821	128	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 6.

All (128) 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:C:357:ILE:HD12	1:C:447:LEU:CD2	1.52	1.37
1:C:357:ILE:HD12	1:C:447:LEU:HD23	1.28	1.14
1:A:251:ARG:NH2	1:A:289:ASN:OD1	1.83	1.10
1:A:246:ILE:HB	1:B:291:LEU:HD21	1.39	1.02
1:D:368:LYS:HG2	1:D:387:ASP:OD2	1.61	1.00
1:C:357:ILE:CD1	1:C:447:LEU:CD2	2.41	0.98
1:E:246:ILE:HD13	1:F:291:LEU:HB3	1.48	0.95
1:C:357:ILE:CD1	1:C:447:LEU:HD23	2.07	0.84
1:A:302:ALA:HB1	1:A:303:LEU:HA	1.57	0.83
1:A:246:ILE:CG2	1:B:403:PHE:HE2	1.90	0.83
1:C:426:THR:HG22	4:C:2049:HOH:O	1.79	0.83
1:C:357:ILE:HD12	1:C:447:LEU:HD21	1.59	0.81
1:A:426:THR:HG21	1:B:242:LEU:HB2	1.65	0.79
1:E:245:ASP:HB3	1:F:402:PRO:HD2	1.63	0.79
1:C:357:ILE:HD12	1:C:447:LEU:HD22	1.65	0.78
1:E:246:ILE:HG21	1:F:291:LEU:HD23	1.63	0.78
1:A:302:ALA:HB1	1:A:303:LEU:CA	2.14	0.77
1:E:357:ILE:HD11	1:E:424:VAL:HG11	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:ARG:NH1	1:A:285:ASP:OD1	2.17	0.75
1:C:357:ILE:HD11	1:C:424:VAL:HG11	1.72	0.72
1:D:361:ALA:HA	1:D:386:ASP:OD2	1.90	0.72
1:A:357:ILE:HD11	1:A:424:VAL:HG11	1.73	0.71
1:A:246:ILE:HG21	1:B:403:PHE:HE2	1.55	0.71
1:A:343:LEU:HD11	1:A:380:ILE:HD12	1.72	0.71
1:A:246:ILE:CB	1:B:291:LEU:HD21	2.17	0.70
1:E:343:LEU:HD11	1:E:380:ILE:HD12	1.74	0.70
1:F:357:ILE:HD11	1:F:424:VAL:HG11	1.74	0.69
1:A:241:ARG:NH1	1:B:426:THR:HG22	2.07	0.69
1:E:246:ILE:HG12	1:F:403:PHE:HE2	1.57	0.68
1:F:243:LEU:HD22	1:F:291:LEU:HD13	1.75	0.68
1:D:440:SER:HB3	1:D:444:GLY:O	1.94	0.67
1:B:343:LEU:HD11	1:B:380:ILE:HD12	1.78	0.66
1:C:265:ARG:HA	1:C:274:LEU:HD21	1.79	0.64
1:E:263:LEU:HD22	1:E:267:ARG:CZ	2.27	0.64
1:A:302:ALA:HA	1:A:346:ASN:HD21	1.62	0.63
1:E:259:LEU:O	1:E:263:LEU:HG	1.99	0.63
1:C:357:ILE:CD1	1:C:447:LEU:HD21	2.22	0.61
1:D:343:LEU:HD11	1:D:380:ILE:HD12	1.83	0.61
1:C:426:THR:HG21	1:D:242:LEU:HB2	1.84	0.60
1:A:288:ILE:O	1:A:292:LEU:HG	2.01	0.60
1:A:246:ILE:HG23	1:B:403:PHE:HE2	1.64	0.59
1:B:357:ILE:HD11	1:B:424:VAL:HG11	1.82	0.59
1:A:227:GLN:HB3	1:D:266:ARG:NH1	2.19	0.58
1:B:265:ARG:HD3	1:B:274:LEU:HB3	1.87	0.57
1:C:254:LEU:HD21	1:C:285:ASP:HA	1.86	0.57
1:B:230:THR:HG22	1:D:256:ARG:HH12	1.70	0.56
1:C:436:LYS:O	1:C:447:LEU:HD12	2.05	0.55
1:B:365:SER:HA	1:B:388:GLY:HA3	1.88	0.55
1:E:246:ILE:HG12	1:F:403:PHE:CE2	2.38	0.55
1:E:287:MET:HE3	1:F:249:GLU:HB3	1.89	0.55
1:A:302:ALA:HB1	1:A:303:LEU:CB	2.37	0.54
1:A:362:LEU:HD23	1:A:369:ILE:HD13	1.89	0.54
1:A:254:LEU:HD21	1:A:285:ASP:HA	1.90	0.54
1:C:402:PRO:HG3	1:D:242:LEU:HD12	1.89	0.54
1:D:311:ASN:HA	1:D:338:PRO:HB2	1.88	0.54
1:A:264:LEU:HD23	1:A:274:LEU:HD13	1.89	0.53
1:C:246:ILE:HD13	1:D:291:LEU:HB2	1.90	0.53
1:A:273:GLU:HG2	1:B:264:LEU:HD21	1.90	0.53
1:B:266:ARG:HB2	1:C:224:SER:HB2	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:357:ILE:HD12	1:B:447:LEU:HD13	1.91	0.53
1:A:240:GLN:HE21	1:A:298:GLN:HB3	1.73	0.53
1:F:243:LEU:CD2	1:F:291:LEU:HD13	2.39	0.53
1:E:275:GLU:O	1:E:279:THR:HG23	2.09	0.53
1:D:357:ILE:HD12	1:D:447:LEU:HD13	1.90	0.52
1:A:357:ILE:HD12	1:A:447:LEU:HD13	1.91	0.52
1:B:305:SER:O	1:D:265:ARG:NH2	2.41	0.52
1:E:254:LEU:HD21	1:E:285:ASP:HA	1.89	0.52
1:E:241:ARG:HD2	1:E:299:GLN:HE22	1.74	0.52
1:B:275:GLU:O	1:B:279:THR:HG23	2.10	0.51
1:A:246:ILE:CG2	1:B:403:PHE:CE2	2.82	0.51
1:C:353:ALA:O	1:C:357:ILE:HG12	2.12	0.50
1:C:430:GLN:OE1	1:D:239:GLN:HG2	2.12	0.50
1:B:257:LEU:O	1:B:261:THR:HG23	2.12	0.50
1:E:357:ILE:HG23	1:E:447:LEU:HD12	1.93	0.49
1:E:240:GLN:HG3	1:F:239:GLN:HE21	1.76	0.49
1:E:244:SER:HA	1:E:295:SER:HB2	1.94	0.49
1:C:357:ILE:CG1	1:C:447:LEU:HD23	2.42	0.49
1:C:275:GLU:O	1:C:279:THR:HG23	2.12	0.49
1:A:302:ALA:HA	1:A:346:ASN:ND2	2.27	0.49
1:B:261:THR:O	1:B:265:ARG:HG2	2.13	0.48
1:B:276:ARG:O	1:B:280:GLU:HG2	2.14	0.48
1:A:275:GLU:O	1:A:279:THR:HG23	2.14	0.48
1:A:246:ILE:HG21	1:B:403:PHE:CE2	2.43	0.48
1:D:357:ILE:HD11	1:D:424:VAL:HG11	1.96	0.47
1:C:357:ILE:CG2	1:C:447:LEU:HD23	2.44	0.47
1:F:265:ARG:NH1	1:F:278:GLU:OE1	2.47	0.47
1:A:246:ILE:HD12	1:B:291:LEU:HD23	1.97	0.47
1:A:273:GLU:HB3	1:B:264:LEU:HD11	1.95	0.47
1:E:245:ASP:CB	1:F:402:PRO:HD2	2.40	0.47
1:B:353:ALA:O	1:B:357:ILE:HG12	2.15	0.47
1:D:275:GLU:O	1:D:279:THR:HG23	2.15	0.47
1:E:232:LEU:O	1:E:236:MET:HG3	2.15	0.47
1:B:361:ALA:O	1:B:365:SER:HB3	2.15	0.47
1:C:359:ARG:HH21	1:C:363:ARG:NH2	2.13	0.47
1:E:357:ILE:HD12	1:E:447:LEU:HD13	1.95	0.47
1:B:345:GLY:HA3	1:B:431:HIS:CD2	2.50	0.46
1:A:239:GLN:HG3	1:B:298:GLN:OE1	2.15	0.46
1:A:246:ILE:HG23	1:B:403:PHE:CE2	2.46	0.45
1:A:353:ALA:O	1:A:357:ILE:HG12	2.17	0.45
1:B:230:THR:CG2	1:D:256:ARG:HH12	2.28	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:TYR:CE2	1:B:219:LEU:HD21	2.51	0.45
1:F:288:ILE:O	1:F:291:LEU:HG	2.17	0.45
1:E:294:MET:HG3	1:E:423:ILE:HD11	1.99	0.44
1:F:353:ALA:O	1:F:357:ILE:HG12	2.17	0.44
1:B:296:ARG:NH1	1:D:285:ASP:OD2	2.50	0.44
1:A:224:SER:HB2	1:D:266:ARG:HB2	1.99	0.44
1:A:227:GLN:HB3	1:D:266:ARG:HH11	1.82	0.44
1:D:438:GLU:O	1:D:445:LEU:HD12	2.17	0.44
1:E:357:ILE:HG23	1:E:447:LEU:CD1	2.47	0.44
1:B:301:ASN:HA	1:B:346:ASN:HD22	1.83	0.44
1:C:238:SER:HB2	1:D:298:GLN:HG2	2.00	0.43
1:E:246:ILE:CD1	1:F:291:LEU:HB3	2.34	0.43
1:E:353:ALA:O	1:E:357:ILE:HG12	2.19	0.43
1:A:241:ARG:HH11	1:B:426:THR:HG22	1.82	0.42
1:A:246:ILE:HD12	1:B:291:LEU:CD2	2.49	0.42
1:C:435:VAL:HG12	1:C:447:LEU:HD11	2.01	0.42
1:E:236:MET:HB3	1:F:239:GLN:HE22	1.84	0.42
1:D:257:LEU:O	1:D:261:THR:HG23	2.19	0.42
1:D:368:LYS:CG	1:D:387:ASP:OD2	2.49	0.41
1:F:301:ASN:HA	1:F:346:ASN:HD22	1.85	0.41
1:A:359:ARG:CZ	1:A:359:ARG:HB3	2.51	0.41
1:F:257:LEU:O	1:F:261:THR:HG23	2.21	0.41
1:F:357:ILE:HD12	1:F:447:LEU:HD13	2.02	0.40
1:C:235:MET:HG3	1:D:298:GLN:HB3	2.03	0.40
1:F:419:LEU:O	1:F:423:ILE:HG12	2.21	0.40
1:B:301:ASN:HD22	1:B:430:GLN:NE2	2.19	0.40
1:E:228:MET:HA	1:E:228:MET:CE	2.51	0.40
1:D:387:ASP:OD1	1:D:442:LEU:HB3	2.21	0.40

There are no symmetry-related clashes.

## 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	229/298 (77%)	222 (97%)	7 (3%)	0	100	100
1	B	223/298 (75%)	215 (96%)	8 (4%)	0	100	100
1	C	237/298 (80%)	234 (99%)	3 (1%)	0	100	100
1	D	210/298 (70%)	203 (97%)	7 (3%)	0	100	100
1	E	225/298 (76%)	220 (98%)	5 (2%)	0	100	100
1	F	221/298 (74%)	215 (97%)	6 (3%)	0	100	100
All	All	1345/1788 (75%)	1309 (97%)	36 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	199/251 (79%)	199 (100%)	0	100	100
1	B	194/251 (77%)	193 (100%)	1 (0%)	92	98
1	C	203/251 (81%)	202 (100%)	1 (0%)	92	98
1	D	182/251 (72%)	182 (100%)	0	100	100
1	E	197/251 (78%)	195 (99%)	2 (1%)	82	94
1	F	195/251 (78%)	195 (100%)	0	100	100
All	All	1170/1506 (78%)	1166 (100%)	4 (0%)	94	99

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	279	THR
1	C	359	ARG
1	E	239	GLN
1	E	297	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	240	GLN
1	A	346	ASN
1	B	346	ASN
1	B	430	GLN
1	D	356	ASN
1	E	299	GLN
1	F	239	GLN
1	F	301	ASN
1	F	346	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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

10 ligands 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	$\# Z  > 2$
3	SO4	A	1456	-	4,4,4	0.15	0	6,6,6	0.26	0
2	ADP	A	501	-	22,29,29	0.61	0	27,45,45	0.79	1 (3%)
3	SO4	B	1456	-	4,4,4	0.15	0	6,6,6	0.08	0
3	SO4	C	1456	-	4,4,4	0.22	0	6,6,6	0.14	0
2	ADP	C	501	-	22,29,29	0.76	0	27,45,45	0.63	0
3	SO4	D	1457	-	4,4,4	0.21	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	SO4	E	1456	-	4,4,4	0.27	0	6,6,6	0.22	0
2	ADP	E	501	-	22,29,29	0.74	0	27,45,45	0.63	0
3	SO4	F	1457	-	4,4,4	0.27	0	6,6,6	0.29	0
2	ADP	F	501	-	22,29,29	0.63	0	27,45,45	0.82	2 (7%)

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
3	SO4	A	1456	-	-	0/0/0/0	0/0/0/0
2	ADP	A	501	-	-	0/12/32/32	0/3/3/3
3	SO4	B	1456	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1456	-	-	0/0/0/0	0/0/0/0
2	ADP	C	501	-	-	0/12/32/32	0/3/3/3
3	SO4	D	1457	-	-	0/0/0/0	0/0/0/0
3	SO4	E	1456	-	-	0/0/0/0	0/0/0/0
2	ADP	E	501	-	-	0/12/32/32	0/3/3/3
3	SO4	F	1457	-	-	0/0/0/0	0/0/0/0
2	ADP	F	501	-	-	0/12/32/32	0/3/3/3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ADP	C4-C5-N7	2.08	111.39	109.48
2	F	501	ADP	C4-C5-N7	2.09	111.40	109.48
2	F	501	ADP	O3A-PA-O5'	2.82	110.42	102.94

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 ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	233/298 (78%)	0.70	22 (9%) 11 8	51, 76, 127, 158	0
1	B	227/298 (76%)	1.16	39 (17%) 2 1	48, 91, 161, 193	0
1	C	239/298 (80%)	0.61	17 (7%) 19 16	52, 71, 126, 173	0
1	D	214/298 (71%)	1.56	51 (23%) 1 1	51, 96, 163, 173	0
1	E	229/298 (76%)	0.78	25 (10%) 7 5	56, 81, 137, 172	0
1	F	225/298 (75%)	0.94	29 (12%) 5 3	58, 88, 150, 186	0
All	All	1367/1788 (76%)	0.95	183 (13%) 4 3	48, 82, 151, 193	0

All (183) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	301	ASN	10.0
1	D	441	PRO	9.7
1	D	332	LEU	8.5
1	B	329	GLY	8.2
1	C	266	ARG	7.5
1	B	396	ARG	7.3
1	F	270	GLU	7.2
1	D	366	HIS	7.1
1	D	322	ALA	7.0
1	D	439	ASP	6.9
1	D	329	GLY	6.4
1	D	437	ALA	6.3
1	D	442	LEU	5.9
1	D	323	PHE	5.9
1	D	369	ILE	5.8
1	B	328	MET	5.8
1	E	259	LEU	5.7
1	D	327	GLN	5.7
1	E	266	ARG	5.7

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Mol	Chain	Res	Type	RSRZ
1	D	361	ALA	5.5
1	D	325	ALA	5.5
1	D	328	MET	5.4
1	B	368	LYS	5.3
1	D	447	LEU	5.2
1	E	267	ARG	5.2
1	A	298	GLN	5.1
1	D	363	ARG	5.1
1	D	358	VAL	5.1
1	E	274	LEU	5.1
1	C	259	LEU	5.0
1	C	218	PHE	5.0
1	D	384	VAL	5.0
1	D	270	GLU	4.9
1	B	332	LEU	4.8
1	E	265	ARG	4.8
1	D	269	GLY	4.8
1	B	358	VAL	4.8
1	C	412	ARG	4.7
1	B	367	THR	4.7
1	F	369	ILE	4.7
1	D	367	THR	4.7
1	B	441	PRO	4.6
1	D	334	VAL	4.6
1	A	299	GLN	4.5
1	A	302	ALA	4.5
1	D	359	ARG	4.5
1	D	330	LYS	4.4
1	B	324	GLU	4.4
1	E	270	GLU	4.4
1	B	325	ALA	4.4
1	B	363	ARG	4.3
1	C	413	GLU	4.3
1	D	371	VAL	4.3
1	D	326	GLU	4.2
1	D	387	ASP	4.2
1	B	331	SER	4.1
1	D	440	SER	4.1
1	A	408	GLU	4.1
1	D	364	TYR	4.1
1	D	365	SER	4.1
1	B	442	LEU	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	415	GLY	4.0
1	D	362	LEU	4.0
1	D	421	LEU	4.0
1	E	264	LEU	4.0
1	B	327	GLN	4.0
1	C	219	LEU	3.9
1	F	323	PHE	3.9
1	B	400	PHE	3.9
1	E	271	SER	3.9
1	E	263	LEU	3.8
1	F	276	ARG	3.7
1	D	368	LYS	3.7
1	B	270	GLU	3.7
1	A	296	ARG	3.6
1	D	435	VAL	3.6
1	D	436	LYS	3.6
1	E	272	LYS	3.6
1	E	261	THR	3.5
1	F	362	LEU	3.5
1	B	362	LEU	3.5
1	A	245	ASP	3.5
1	B	404	TYR	3.5
1	A	246	ILE	3.4
1	D	386	ASP	3.4
1	E	262	ALA	3.4
1	D	360	ASN	3.4
1	C	274	LEU	3.4
1	F	367	THR	3.4
1	F	366	HIS	3.4
1	B	365	SER	3.4
1	E	408	GLU	3.4
1	A	409	ALA	3.3
1	D	445	LEU	3.2
1	A	300	LYS	3.2
1	F	277	ILE	3.2
1	B	369	ILE	3.2
1	A	325	ALA	3.2
1	E	276	ARG	3.1
1	B	366	HIS	3.1
1	D	331	SER	3.1
1	E	293	VAL	3.0
1	C	414	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	416	GLY	3.0
1	B	445	LEU	3.0
1	A	411	ASP	3.0
1	B	269	GLY	3.0
1	F	274	LEU	3.0
1	F	328	MET	3.0
1	A	292	LEU	2.9
1	A	412	ARG	2.9
1	D	443	GLY	2.9
1	D	318	LEU	2.9
1	D	438	GLU	2.9
1	F	327	GLN	2.9
1	B	390	GLY	2.9
1	B	386	ASP	2.9
1	B	330	LYS	2.9
1	B	322	ALA	2.8
1	B	399	ILE	2.8
1	A	295	SER	2.8
1	B	395	ASP	2.8
1	C	267	ARG	2.8
1	B	394	GLU	2.8
1	E	292	LEU	2.8
1	B	391	VAL	2.8
1	B	405	ARG	2.8
1	B	439	ASP	2.8
1	D	324	GLU	2.7
1	F	329	GLY	2.7
1	D	446	ARG	2.7
1	E	410	ARG	2.7
1	B	323	PHE	2.7
1	D	333	THR	2.7
1	A	410	ARG	2.7
1	D	217	GLU	2.6
1	F	399	ILE	2.6
1	C	264	LEU	2.6
1	F	456	ARG	2.6
1	E	301	ASN	2.6
1	B	398	GLN	2.6
1	B	291	LEU	2.6
1	D	370	GLU	2.6
1	F	445	LEU	2.5
1	D	391	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	268	SER	2.5
1	C	265	ARG	2.5
1	F	361	ALA	2.5
1	A	363	ARG	2.5
1	C	411	ASP	2.4
1	B	389	PRO	2.4
1	E	278	GLU	2.4
1	E	269	GLY	2.4
1	B	320	ASN	2.4
1	E	257	LEU	2.4
1	A	241	ARG	2.4
1	C	270	GLU	2.4
1	E	409	ALA	2.4
1	F	330	LYS	2.3
1	D	238	SER	2.3
1	F	244	SER	2.3
1	A	369	ILE	2.3
1	F	291	LEU	2.3
1	C	262	ALA	2.3
1	E	254	LEU	2.2
1	F	269	GLY	2.2
1	F	326	GLU	2.2
1	A	303	LEU	2.2
1	F	364	TYR	2.2
1	C	408	GLU	2.2
1	F	324	GLU	2.2
1	F	368	LYS	2.2
1	E	275	GLU	2.2
1	D	428	ILE	2.1
1	F	283	ARG	2.1
1	E	249	GLU	2.1
1	F	273	GLU	2.1
1	F	404	TYR	2.1
1	F	363	ARG	2.1
1	D	420	GLY	2.1
1	F	371	VAL	2.1
1	C	269	GLY	2.1
1	A	328	MET	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 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. 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, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ADP	F	501	27/27	0.86	0.26	-0.25	118,131,137,139	0
2	ADP	E	501	27/27	0.98	0.19	-0.42	62,76,89,92	0
2	ADP	C	501	27/27	0.97	0.18	-0.57	54,66,80,82	0
2	ADP	A	501	27/27	0.97	0.19	-0.58	60,75,87,91	0
3	SO4	A	1456	5/5	0.97	0.15	-	98,101,101,103	0
3	SO4	D	1457	5/5	0.81	0.20	-	147,148,149,150	0
3	SO4	F	1457	5/5	0.97	0.13	-	91,94,95,96	0
3	SO4	C	1456	5/5	0.89	0.19	-	119,120,121,121	0
3	SO4	E	1456	5/5	0.92	0.17	-	114,114,115,117	0
3	SO4	B	1456	5/5	0.84	0.21	-	157,159,159,159	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.