



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 17, 2023 – 12:30 PM EDT

PDB ID : 4XZ5  
Title : Structure of the thermostable alpha-Carbonic Anydrase from *Thiomicrospira crunogena* XCL-2 gammaproteobacterium  
Authors : Mahon, B.P.; Diaz-Torres, N.A.; Pinard, M.A.; McKenna, R.  
Deposited on : 2015-02-03  
Resolution : 2.60 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

<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>.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35.1

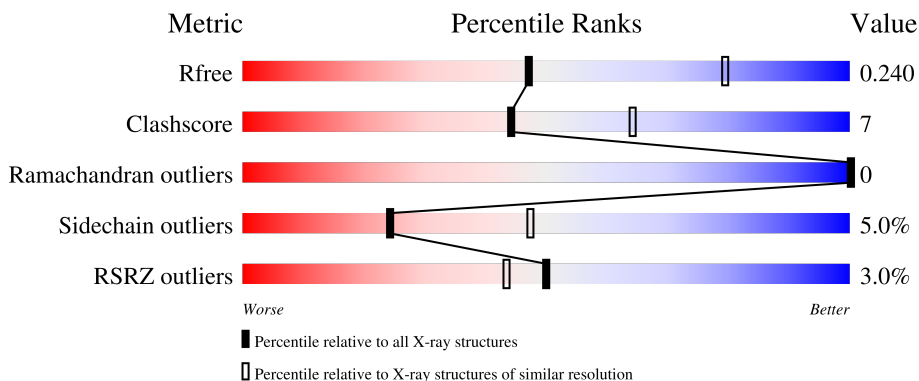
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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  $\geq 3$ , 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	233	 3% 79% 17% ..
1	B	233	 3% 80% 15% ..
1	C	233	 3% 78% 20% ..
1	D	233	 3% 78% 18% ..

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7497 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 Carbonic anhydrase, alpha family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	230	1860	1197	316	340	7	0	0	0
1	B	230	1860	1197	316	340	7	0	0	0
1	C	230	1860	1197	316	340	7	0	0	0
1	D	230	1860	1197	316	340	7	0	0	0

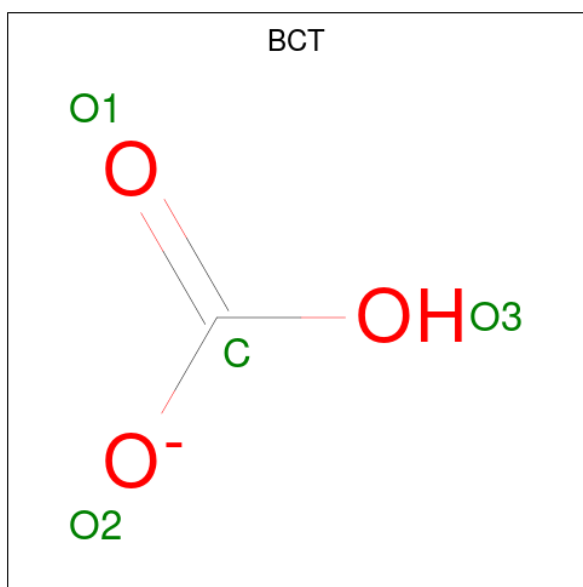
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	307	ALA	-	expression tag	UNP Q31FD6
B	307	ALA	-	expression tag	UNP Q31FD6
C	307	ALA	-	expression tag	UNP Q31FD6
D	307	ALA	-	expression tag	UNP Q31FD6

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

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

- Molecule 3 is BICARBONATE ION (three-letter code: BCT) (formula: CHO<sub>3</sub>).



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

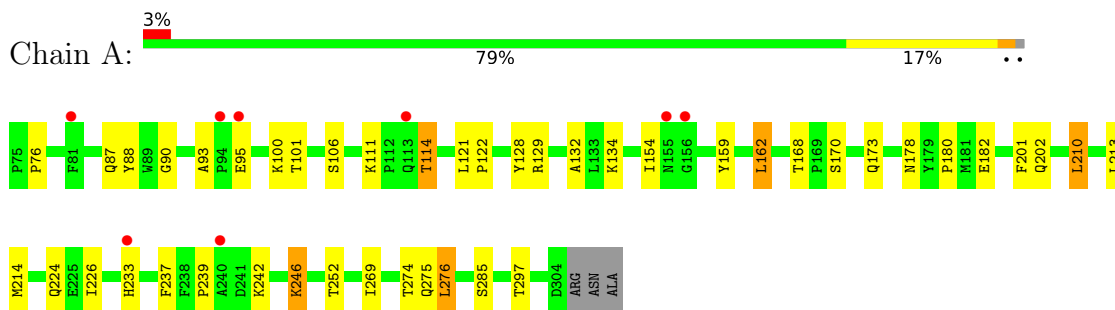
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	6	Total O 6 6	0	0
4	B	17	Total O 17 17	0	0
4	C	7	Total O 7 7	0	0
4	D	7	Total O 7 7	0	0

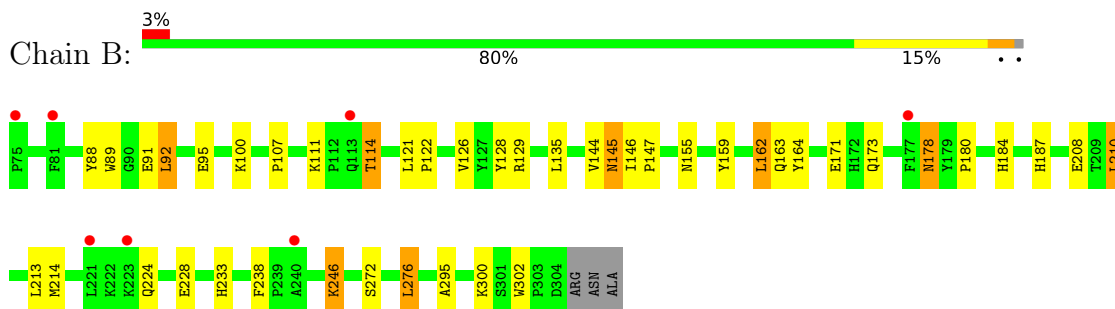
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

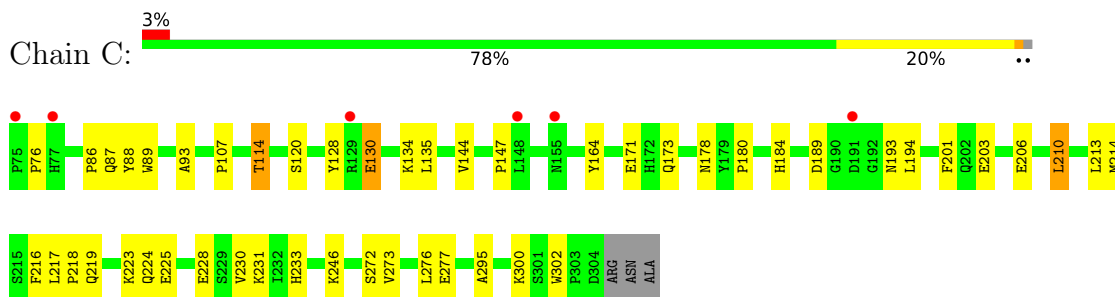
- Molecule 1: Carbonic anhydrase, alpha family



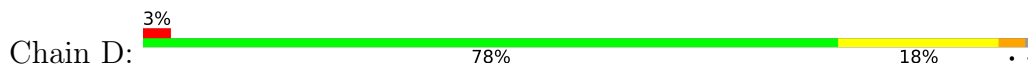
- Molecule 1: Carbonic anhydrase, alpha family

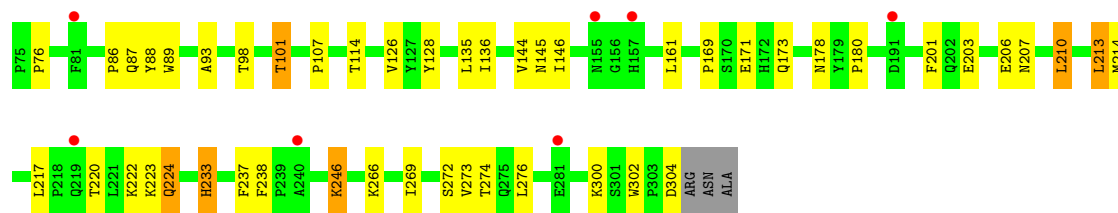


- Molecule 1: Carbonic anhydrase, alpha family



- Molecule 1: Carbonic anhydrase, alpha family





## 4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.06Å 102.23Å 105.02Å 90.00° 127.26° 90.00°	Depositor
Resolution (Å)	19.91 – 2.60 19.91 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.1 (19.91-2.60) 94.8 (19.91-2.60)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.02 (at 2.59Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: dev_1839)	Depositor
R, $R_{free}$	0.171 , 0.240 0.172 , 0.240	Depositor DCC
$R_{free}$ test set	1643 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.7	Xtrriage
Anisotropy	0.495	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 55.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.022 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7497	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: BCT, 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	0.45	0/1919	0.61	1/2604 (0.0%)
1	B	0.47	0/1919	0.66	1/2604 (0.0%)
1	C	0.46	0/1919	0.64	0/2604
1	D	0.45	0/1919	0.62	0/2604
All	All	0.46	0/7676	0.63	2/10416 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	162	LEU	CA-CB-CG	5.48	127.91	115.30
1	A	162	LEU	CA-CB-CG	5.27	127.43	115.30

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	1860	0	1806	25	1
1	B	1860	0	1806	24	2
1	C	1860	0	1806	27	0
1	D	1860	0	1806	28	1
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	4	0	0	0	0
3	B	4	0	1	0	0
3	C	4	0	1	0	0
3	D	4	0	1	0	0
4	A	6	0	0	0	0
4	B	17	0	0	0	0
4	C	7	0	0	2	0
4	D	7	0	0	0	0
All	All	7497	0	7227	97	2

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 7.

All (97) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:ARG:NH2	1:C:203:GLU:O	2.13	0.81
1:B:173:GLN:HE21	1:B:178:ASN:HD21	1.35	0.73
1:D:107:PRO:HG3	1:D:171:GLU:HB3	1.80	0.64
1:B:91:GLU:OE2	1:B:100:LYS:NZ	2.28	0.63
1:C:300:LYS:HD2	1:C:302:TRP:CZ2	2.32	0.63
1:D:98:THR:HA	1:D:101:THR:HG22	1.80	0.63
1:C:76:PRO:HG3	1:C:93:ALA:HB2	1.85	0.59
1:C:135:LEU:HD13	1:C:144:VAL:HG22	1.85	0.59
1:A:132:ALA:HB3	1:C:273:VAL:HG21	1.85	0.58
1:C:107:PRO:HG3	1:C:171:GLU:HB3	1.85	0.58
1:C:128:TYR:HB2	1:C:233:HIS:CE1	2.38	0.58
1:B:88:TYR:O	1:B:92:LEU:HD13	2.06	0.56
1:C:130:GLU:HG2	1:C:231:LYS:HB3	1.87	0.56
1:B:111:LYS:O	1:B:114:THR:HG22	2.06	0.56
1:D:173:GLN:HG2	1:D:178:ASN:OD1	2.06	0.55
1:C:210:LEU:HD13	1:C:214:MET:HE1	1.90	0.53
1:B:300:LYS:HD2	1:B:302:TRP:CZ2	2.43	0.53
1:A:121:LEU:HD12	1:A:122:PRO:HD2	1.90	0.53
1:A:90:GLY:HA3	1:A:100:LYS:HB3	1.90	0.52
1:C:218:PRO:HB3	1:C:225:GLU:HG3	1.92	0.52
1:D:300:LYS:HD2	1:D:302:TRP:CZ2	2.44	0.52
1:B:180:PRO:HB2	1:B:276:LEU:HG	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:135:LEU:HD13	1:D:144:VAL:HG22	1.91	0.52
1:B:128:TYR:HB2	1:B:233:HIS:CE1	2.46	0.51
1:A:173:GLN:HG2	1:A:178:ASN:OD1	2.10	0.51
1:B:159:TYR:HB3	1:B:187:HIS:HB3	1.93	0.50
1:D:87:GLN:HG3	1:D:88:TYR:CD1	2.47	0.50
1:D:128:TYR:HB2	1:D:233:HIS:HE1	1.76	0.49
1:D:203:GLU:OE1	1:D:273:VAL:HG22	2.11	0.49
1:C:114:THR:HG23	4:C:504:HOH:O	2.12	0.49
1:A:76:PRO:HG3	1:A:93:ALA:HB2	1.94	0.49
1:C:206:GLU:HA	1:C:272:SER:HB3	1.95	0.49
1:B:173:GLN:HE21	1:B:178:ASN:ND2	2.09	0.49
1:A:210:LEU:O	1:A:214:MET:HG3	2.13	0.49
1:B:178:ASN:N	1:B:178:ASN:HD22	2.12	0.48
1:C:120:SER:OG	1:D:304:ASP:OD1	2.18	0.48
1:B:107:PRO:HG3	1:B:171:GLU:HB3	1.96	0.48
1:D:126:VAL:HG21	1:D:238:PHE:HE1	1.79	0.48
1:D:76:PRO:HG3	1:D:93:ALA:HB2	1.94	0.48
1:B:145:ASN:OD1	1:B:145:ASN:N	2.46	0.47
1:D:128:TYR:HB2	1:D:233:HIS:CE1	2.49	0.47
1:D:86:PRO:HA	1:D:89:TRP:CD2	2.49	0.47
1:D:146:ILE:HG13	1:D:161:LEU:HG	1.97	0.47
1:D:214:MET:HA	1:D:217:LEU:HG	1.96	0.47
1:A:111:LYS:O	1:A:114:THR:HG22	2.15	0.47
1:C:173:GLN:HG2	1:C:178:ASN:ND2	2.29	0.46
1:C:216:PHE:CZ	1:C:230:VAL:HG21	2.50	0.46
1:B:135:LEU:HD13	1:B:144:VAL:HG22	1.98	0.46
1:D:180:PRO:HD2	1:D:201:PHE:O	2.16	0.46
1:D:206:GLU:HG3	1:D:207:ASN:N	2.31	0.46
1:A:128:TYR:HB2	1:A:233:HIS:CE1	2.51	0.45
1:A:297:THR:OG1	1:B:246:LYS:HE2	2.17	0.45
1:C:214:MET:HA	1:C:217:LEU:HG	1.97	0.44
1:B:121:LEU:HD12	1:B:122:PRO:HD2	1.98	0.44
1:D:266:LYS:HD3	1:D:302:TRP:CH2	2.52	0.44
1:B:272:SER:O	1:B:276:LEU:HD22	2.16	0.44
1:D:220:THR:HB	1:D:223:LYS:HG3	2.00	0.44
1:A:180:PRO:HB2	1:A:276:LEU:HG	1.99	0.44
1:A:210:LEU:HD13	1:A:214:MET:CE	2.47	0.44
1:C:87:GLN:HG3	1:C:88:TYR:CD1	2.52	0.44
1:D:169:PRO:HA	1:D:180:PRO:O	2.17	0.44
1:C:164:TYR:HA	1:C:184:HIS:O	2.18	0.43
1:A:180:PRO:HD2	1:A:201:PHE:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:VAL:HG21	1:B:238:PHE:HE1	1.83	0.43
1:C:295:ALA:O	1:D:246:LYS:NZ	2.52	0.43
1:A:134:LYS:HG2	1:A:226:ILE:HG12	2.01	0.42
1:B:210:LEU:O	1:B:214:MET:HG3	2.19	0.42
1:A:226:ILE:HD12	1:C:277:GLU:OE2	2.19	0.42
1:B:122:PRO:HB3	1:B:155:ASN:HB2	2.00	0.42
1:C:180:PRO:HD2	1:C:201:PHE:O	2.20	0.42
1:D:210:LEU:HD13	1:D:214:MET:HE1	2.02	0.42
1:A:239:PRO:O	1:A:242:LYS:NZ	2.37	0.42
1:B:89:TRP:HA	1:B:92:LEU:HD22	2.02	0.42
1:B:146:ILE:HA	1:B:147:PRO:HD3	1.88	0.42
1:A:100:LYS:HG2	1:A:101:THR:HG23	2.02	0.42
1:C:86:PRO:HA	1:C:89:TRP:CD2	2.55	0.42
1:B:164:TYR:HA	1:B:184:HIS:O	2.20	0.41
1:D:210:LEU:HD13	1:D:214:MET:CE	2.50	0.41
1:C:189:ASP:CG	1:C:193:ASN:HB2	2.40	0.41
1:C:273:VAL:HG23	4:C:503:HOH:O	2.19	0.41
1:A:168:THR:O	1:A:285:SER:HA	2.19	0.41
1:A:170:SER:HB2	1:A:182:GLU:HB2	2.03	0.41
1:D:206:GLU:HB2	1:D:272:SER:HB3	2.02	0.41
1:D:237:PHE:O	1:D:269:ILE:HD13	2.20	0.41
1:D:145:ASN:HD22	1:D:224:GLN:NE2	2.18	0.41
1:A:106:SER:C	1:A:252:THR:HG22	2.41	0.41
1:A:154:ILE:HD12	1:A:159:TYR:CE2	2.55	0.41
1:A:246:LYS:NZ	1:B:295:ALA:O	2.54	0.41
1:A:237:PHE:O	1:A:269:ILE:HD13	2.20	0.41
1:C:216:PHE:CE2	1:C:230:VAL:HG21	2.56	0.41
1:C:223:LYS:HG2	1:C:225:GLU:OE2	2.20	0.41
1:D:136:ILE:HD11	1:D:222:LYS:HA	2.03	0.41
1:A:87:GLN:HG3	1:A:88:TYR:CD1	2.56	0.41
1:D:213:LEU:HD23	1:D:213:LEU:HA	1.94	0.41
1:B:95:GLU:H	1:B:95:GLU:CD	2.23	0.40
1:C:134:LYS:HE2	1:C:147:PRO:HG3	2.03	0.40
1:A:214:MET:HE1	1:A:275:GLN:HB3	2.02	0.40

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	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:ARG:NH2	1:D:203:GLU:O[3_545]	2.03	0.17
1:A:202:GLN:NE2	1:B:208:GLU:OE1[4_7410]	2.06	0.14

## 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	228/233 (98%)	223 (98%)	5 (2%)	0	100	100
1	B	228/233 (98%)	223 (98%)	5 (2%)	0	100	100
1	C	228/233 (98%)	223 (98%)	5 (2%)	0	100	100
1	D	228/233 (98%)	222 (97%)	6 (3%)	0	100	100
All	All	912/932 (98%)	891 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	202/204 (99%)	193 (96%)	9 (4%)	27	52
1	B	202/204 (99%)	190 (94%)	12 (6%)	19	39
1	C	202/204 (99%)	192 (95%)	10 (5%)	24	47
1	D	202/204 (99%)	193 (96%)	9 (4%)	27	52
All	All	808/816 (99%)	768 (95%)	40 (5%)	24	47

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	GLU
1	A	114	THR

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	162	LEU
1	A	210	LEU
1	A	213	LEU
1	A	224	GLN
1	A	246	LYS
1	A	274	THR
1	A	276	LEU
1	B	92	LEU
1	B	114	THR
1	B	145	ASN
1	B	162	LEU
1	B	163	GLN
1	B	178	ASN
1	B	210	LEU
1	B	213	LEU
1	B	224	GLN
1	B	228	GLU
1	B	246	LYS
1	B	276	LEU
1	C	114	THR
1	C	130	GLU
1	C	194	LEU
1	C	210	LEU
1	C	213	LEU
1	C	219	GLN
1	C	224	GLN
1	C	228	GLU
1	C	246	LYS
1	C	276	LEU
1	D	101	THR
1	D	114	THR
1	D	210	LEU
1	D	213	LEU
1	D	224	GLN
1	D	233	HIS
1	D	246	LYS
1	D	274	THR
1	D	276	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	143	GLN
1	B	178	ASN
1	C	178	ASN
1	D	145	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

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

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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
3	BCT	B	402[B]	2	2,3,3	1.12	0	2,3,3	4.29	2 (100%)
3	BCT	C	402[B]	2	2,3,3	1.17	0	2,3,3	4.01	2 (100%)
3	BCT	D	402[B]	2	2,3,3	1.08	0	2,3,3	4.36	2 (100%)
3	BCT	A	402[B]	2	2,3,3	1.23	0	2,3,3	3.92	2 (100%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	B	402[B]	BCT	O2-C-O1	5.56	133.96	119.55
3	D	402[B]	BCT	O2-C-O1	5.24	133.15	119.55
3	A	402[B]	BCT	O2-C-O1	5.09	132.76	119.55
3	C	402[B]	BCT	O2-C-O1	4.82	132.06	119.55
3	D	402[B]	BCT	O3-C-O1	-3.24	111.13	119.55
3	C	402[B]	BCT	O3-C-O1	-2.99	111.80	119.55
3	B	402[B]	BCT	O3-C-O1	-2.43	113.25	119.55
3	A	402[B]	BCT	O3-C-O1	-2.18	113.88	119.55

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<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	230/233 (98%)	-0.16	8 (3%) 44 36	14, 29, 59, 76	0
1	B	230/233 (98%)	-0.15	7 (3%) 50 43	14, 28, 58, 79	0
1	C	230/233 (98%)	-0.09	6 (2%) 56 50	17, 33, 71, 95	0
1	D	230/233 (98%)	-0.00	7 (3%) 50 43	17, 34, 66, 84	0
All	All	920/932 (98%)	-0.10	28 (3%) 50 43	14, 31, 64, 95	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	81	PHE	4.8
1	D	155	ASN	3.9
1	B	81	PHE	3.6
1	C	75	PRO	3.3
1	B	75	PRO	3.2
1	C	191	ASP	3.1
1	B	240	ALA	2.8
1	C	155	ASN	2.7
1	B	177	PHE	2.7
1	D	281	GLU	2.6
1	C	148	LEU	2.6
1	D	240	ALA	2.5
1	A	156	GLY	2.5
1	A	81	PHE	2.4
1	C	77	HIS	2.4
1	C	129	ARG	2.3
1	A	240	ALA	2.2
1	A	94	PRO	2.2
1	A	113	GLN	2.2
1	A	95	GLU	2.2
1	B	221	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	233	HIS	2.1
1	B	223	LYS	2.1
1	D	191	ASP	2.1
1	B	113	GLN	2.1
1	A	155	ASN	2.0
1	D	219	GLN	2.0
1	D	157	HIS	2.0

## 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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<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	B-factors(Å <sup>2</sup> )	Q<0.9
3	BCT	B	402[B]	4/4	0.98	0.15	28,31,34,37	0
3	BCT	C	402[B]	4/4	0.98	0.11	24,24,29,38	0
3	BCT	A	402[B]	4/4	0.99	0.12	27,33,34,42	0
3	BCT	D	402[B]	4/4	0.99	0.11	19,20,25,28	0
2	ZN	A	401	1/1	1.00	0.12	25,25,25,25	0
2	ZN	B	401	1/1	1.00	0.10	24,24,24,24	0
2	ZN	C	401	1/1	1.00	0.07	26,26,26,26	0
2	ZN	D	401	1/1	1.00	0.09	24,24,24,24	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.