



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 01:54 am BST

PDB ID : 3FVL  
Title : Crystallogic studies on the Complex of Carboxypeptidase A with inhibitors using alpha-hydroxy ketone as zinc-binding group  
Authors : Wang, S.F.; Jin, J.-Y.; Tian, G.R.  
Deposited on : 2009-01-16  
Resolution : 1.85 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.

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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 : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

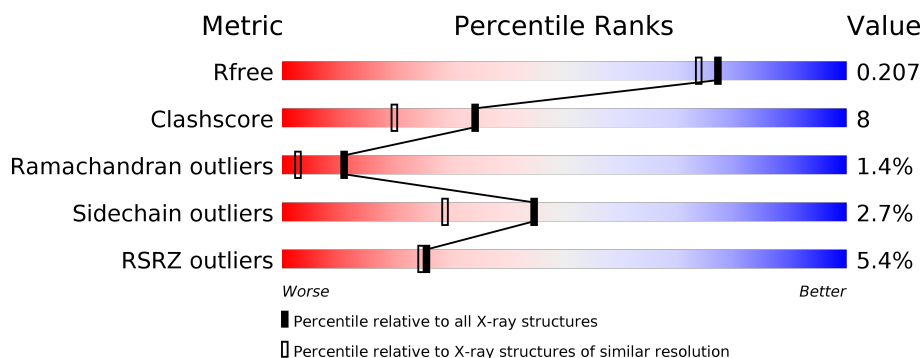
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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 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	307	<div> <div>5%</div> <div> <div></div> <div>83%</div> <div>15%</div> <div>..</div> </div> </div>
1	C	307	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>..</div> </div> </div>
1	E	307	<div> <div>5%</div> <div> <div></div> <div>82%</div> <div>16%</div> <div>..</div> </div> </div>

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
3	BHK	A	311	-	-	-	X
3	BHK	C	311	-	-	-	X
3	BHK	E	311	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7907 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 Carboxypeptidase A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2441	1563	403	470	5			
1	C	307	Total	C	N	O	S	0	0	0
			2441	1563	403	470	5			
1	E	307	Total	C	N	O	S	0	0	0
			2441	1563	403	470	5			

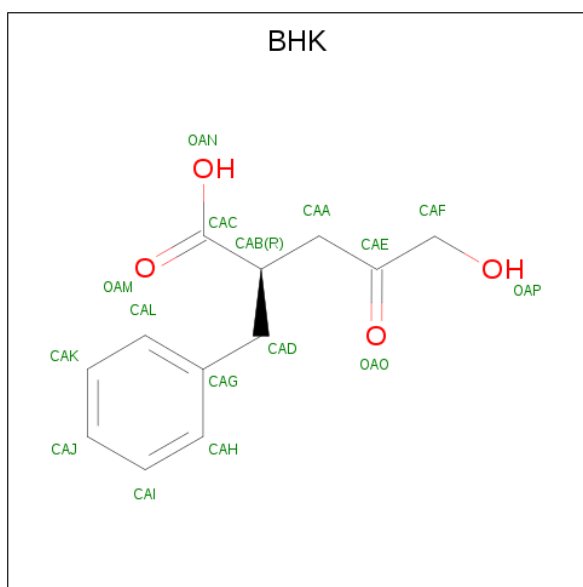
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	305	VAL	LEU	SEE REMARK 999	UNP P00730
C	305	VAL	LEU	SEE REMARK 999	UNP P00730
E	305	VAL	LEU	SEE REMARK 999	UNP P00730

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

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

- Molecule 3 is (2R)-2-benzyl-5-hydroxy-4-oxopentanoic acid (three-letter code: BHK) (formula: C<sub>12</sub>H<sub>14</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			16	12	4		
3	C	1	Total	C	O	0	0
			16	12	4		
3	E	1	Total	C	O	0	0
			16	12	4		

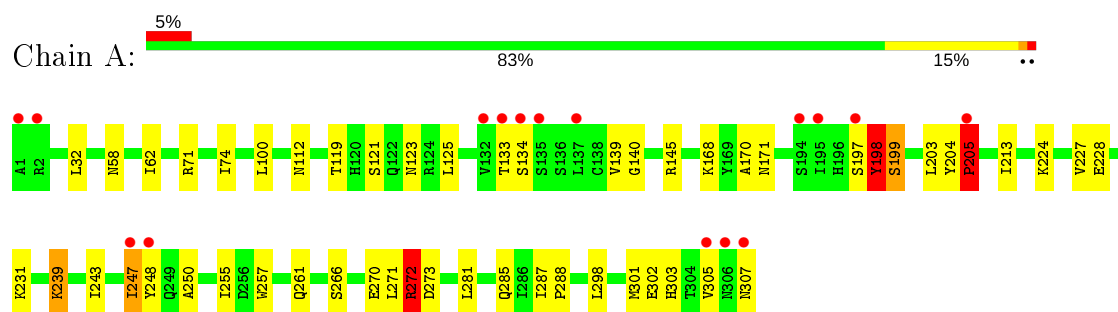
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	177	Total	O	0	0
			177	177		
4	C	177	Total	O	0	0
			177	177		
4	E	179	Total	O	0	0
			179	179		

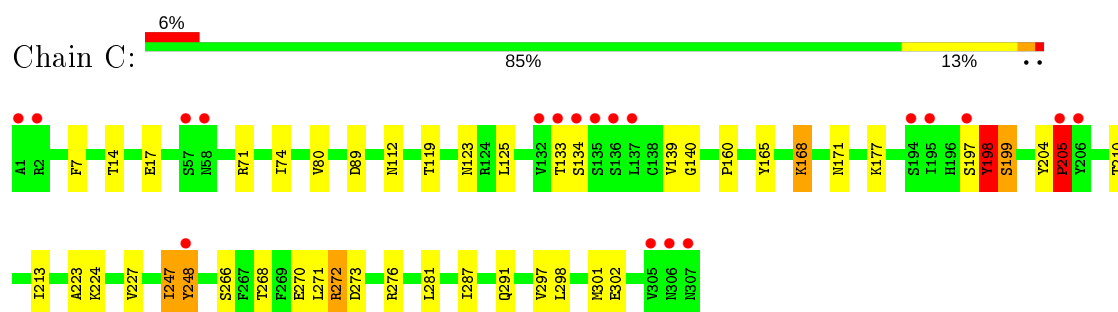
### 3 Residue-property plots [i](#)

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

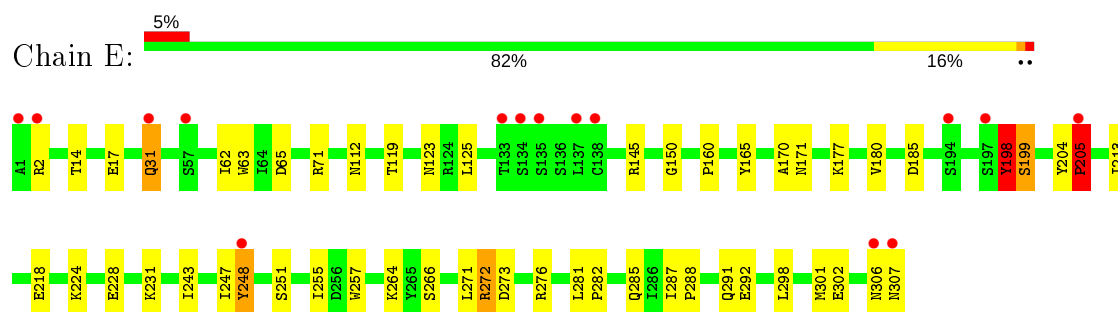
- Molecule 1: Carboxypeptidase A1



- Molecule 1: Carboxypeptidase A1



- Molecule 1: Carboxypeptidase A1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.35Å 59.79Å 99.36Å 90.00° 103.94° 90.00°	Depositor
Resolution (Å)	50.00 – 1.85 35.59 – 1.85	Depositor EDS
% Data completeness (in resolution range)	95.5 (50.00-1.85) 95.5 (35.59-1.85)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	20.68 (at 1.85Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.209 , 0.238 0.205 , 0.207	Depositor DCC
$R_{free}$ test set	3464 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	16.3	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 45.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7907	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 69.34 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.7614e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: ZN, BHK

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	0.34	0/2507	0.65	2/3407 (0.1%)
1	C	0.33	0/2507	0.63	1/3407 (0.0%)
1	E	0.34	0/2507	0.63	0/3407
All	All	0.34	0/7521	0.64	3/10221 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	272	ARG	N-CA-C	5.89	126.91	111.00
1	A	205	PRO	N-CA-C	5.43	126.22	112.10
1	A	272	ARG	N-CA-C	5.24	125.15	111.00

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	2441	0	2347	49	0
1	C	2441	0	2347	35	0
1	E	2441	0	2347	38	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	1	0	0	0	0
3	A	16	0	13	6	0
3	C	16	0	13	1	0
3	E	16	0	13	2	0
4	A	177	0	0	2	0
4	C	177	0	0	1	0
4	E	179	0	0	0	0
All	All	7907	0	7080	122	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 8.

All (122) 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:168:LYS:HE3	1:C:168:LYS:H	1.09	1.17
1:A:168:LYS:HD2	1:A:168:LYS:H	1.04	1.08
1:C:168:LYS:H	1:C:168:LYS:CE	1.82	0.93
1:A:133:THR:HG21	1:A:168:LYS:NZ	1.86	0.91
1:A:168:LYS:HD2	1:A:168:LYS:N	1.87	0.88
1:A:203:LEU:HD13	1:A:247:ILE:HD11	1.56	0.87
1:A:243:ILE:O	1:A:247:ILE:HD12	1.82	0.79
1:A:168:LYS:CD	1:A:168:LYS:H	1.90	0.78
1:E:31:GLN:H	1:E:31:GLN:NE2	1.82	0.78
1:E:198:TYR:HA	1:E:271:LEU:O	1.87	0.74
1:C:168:LYS:N	1:C:168:LYS:HE3	1.95	0.73
1:E:224:LYS:O	1:E:228:GLU:HG3	1.92	0.69
1:E:204:TYR:HB2	1:E:205:PRO:HD2	1.76	0.67
1:C:204:TYR:HB2	1:C:205:PRO:HD2	1.77	0.67
1:E:243:ILE:HD11	1:E:255:ILE:HD11	1.76	0.67
1:C:205:PRO:HD3	1:C:266:SER:O	1.95	0.67
1:A:204:TYR:HB2	1:A:205:PRO:HD2	1.78	0.66
1:A:133:THR:HG21	1:A:168:LYS:CE	2.27	0.64
1:A:203:LEU:CD1	1:A:247:ILE:HD11	2.27	0.64
1:C:205:PRO:HB2	1:C:213:ILE:HG21	1.79	0.64
1:A:224:LYS:O	1:A:228:GLU:HG3	1.98	0.63
1:A:205:PRO:HB2	1:A:213:ILE:HG21	1.81	0.62
1:C:198:TYR:HA	1:C:271:LEU:O	2.01	0.61
1:A:133:THR:HG21	1:A:168:LYS:HZ1	1.64	0.60
1:E:298:LEU:O	1:E:302:GLU:HG3	2.02	0.60
1:C:270:GLU:OE2	3:C:311:BHK:HAA	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:THR:HA	1:A:123:ASN:O	2.02	0.58
1:A:62:ILE:HD13	1:A:301:MET:HG2	1.85	0.58
1:A:198:TYR:HA	1:A:271:LEU:O	2.02	0.58
1:E:288:PRO:O	1:E:292:GLU:HG2	2.04	0.58
1:A:74:ILE:HD13	1:A:281:LEU:HD23	1.85	0.58
1:A:243:ILE:HD11	1:A:255:ILE:HD11	1.85	0.57
1:E:205:PRO:HD3	1:E:266:SER:O	2.05	0.57
1:C:210:THR:HG22	1:E:185:ASP:OD1	2.05	0.56
1:E:71:ARG:HD3	1:E:125:LEU:O	2.04	0.56
1:A:145:ARG:HG2	3:A:311:BHK:HAI	1.86	0.56
1:A:257:TRP:O	1:A:261:GLN:HG2	2.07	0.55
1:E:31:GLN:H	1:E:31:GLN:CD	2.11	0.53
1:A:224:LYS:O	1:A:227:VAL:HG22	2.08	0.53
3:A:311:BHK:HAH	3:A:311:BHK:CAC	2.39	0.53
1:A:205:PRO:HD3	1:A:266:SER:O	2.08	0.53
1:A:133:THR:HG21	1:A:168:LYS:HZ2	1.70	0.53
1:A:231:LYS:HZ2	1:A:231:LYS:HB2	1.73	0.53
1:A:273:ASP:HB2	1:A:281:LEU:HD13	1.92	0.52
1:C:272:ARG:O	1:C:273:ASP:HB3	2.10	0.52
1:C:74:ILE:HD13	1:C:281:LEU:HD23	1.90	0.52
1:A:231:LYS:NZ	1:A:231:LYS:HB2	2.25	0.51
1:C:119:THR:HA	1:C:123:ASN:O	2.10	0.51
1:C:247:ILE:HG22	1:C:248:TYR:CG	2.46	0.51
1:A:272:ARG:O	1:A:273:ASP:HB3	2.11	0.51
1:A:71:ARG:HD3	1:A:125:LEU:O	2.11	0.50
1:C:204:TYR:HB2	1:C:205:PRO:CD	2.41	0.50
1:A:250:ALA:HB1	3:A:311:BHK:CAI	2.42	0.50
1:A:298:LEU:O	1:A:302:GLU:HG3	2.11	0.49
1:E:62:ILE:HD13	1:E:301:MET:HG2	1.93	0.49
1:A:139:VAL:HG22	1:A:140:GLY:N	2.27	0.49
1:C:133:THR:O	1:C:134:SER:HB3	2.13	0.49
3:E:311:BHK:HAH	3:E:311:BHK:CAC	2.43	0.49
1:A:198:TYR:O	1:A:199:SER:HB3	2.12	0.48
1:E:150:GLY:O	1:E:251:SER:HB2	2.14	0.48
1:C:198:TYR:O	1:C:199:SER:HB3	2.14	0.48
1:A:303:HIS:O	1:A:307:ASN:HB2	2.13	0.48
1:C:74:ILE:CG2	1:C:197:SER:HB2	2.42	0.48
1:C:71:ARG:HD3	1:C:125:LEU:O	2.14	0.48
1:C:168:LYS:N	1:C:168:LYS:CE	2.64	0.48
1:E:198:TYR:O	1:E:199:SER:HB3	2.14	0.47
1:C:171:ASN:HB3	1:C:177:LYS:HD3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:287:ILE:O	1:C:291:GLN:HG3	2.14	0.47
1:E:14:THR:OG1	1:E:17:GLU:HG3	2.15	0.47
1:C:224:LYS:O	1:C:227:VAL:HG22	2.14	0.47
1:C:273:ASP:HB2	1:C:281:LEU:HD13	1.96	0.47
1:E:145:ARG:HG2	3:E:311:BHK:HAI	1.97	0.47
1:E:119:THR:HA	1:E:123:ASN:O	2.15	0.47
1:A:272:ARG:HH11	1:A:285:GLN:HE21	1.63	0.47
1:E:272:ARG:O	1:E:273:ASP:HB3	2.15	0.46
1:C:297:VAL:O	1:C:301:MET:HG3	2.16	0.46
1:E:287:ILE:O	1:E:291:GLN:HG3	2.15	0.46
1:C:160:PRO:HA	1:C:165:TYR:CD2	2.51	0.46
1:C:198:TYR:O	1:C:199:SER:CB	2.63	0.46
1:E:63:TRP:HE1	1:E:65:ASP:HB3	1.80	0.46
1:E:231:LYS:HB2	1:E:231:LYS:NZ	2.31	0.46
1:C:268:THR:HG21	4:C:480:HOH:O	2.16	0.45
1:A:133:THR:HG21	1:A:168:LYS:HE3	1.97	0.45
1:E:160:PRO:HA	1:E:165:TYR:CD2	2.52	0.45
1:A:303:HIS:HD2	4:A:404:HOH:O	2.00	0.44
1:E:170:ALA:O	1:E:171:ASN:HB2	2.18	0.44
1:E:287:ILE:HB	1:E:288:PRO:HD3	2.00	0.44
1:A:204:TYR:HB2	1:A:205:PRO:CD	2.46	0.44
1:E:204:TYR:HB2	1:E:205:PRO:CD	2.45	0.43
1:C:14:THR:OG1	1:C:17:GLU:HG3	2.19	0.43
1:E:171:ASN:HB3	1:E:177:LYS:HD3	2.01	0.43
1:A:198:TYR:O	1:A:199:SER:CB	2.66	0.43
1:E:205:PRO:HB2	1:E:213:ILE:HG21	2.00	0.43
1:E:218:GLU:OE1	1:E:264:LYS:HE2	2.18	0.43
1:E:247:ILE:HG22	1:E:248:TYR:CG	2.55	0.42
1:E:31:GLN:N	1:E:31:GLN:NE2	2.61	0.42
1:C:139:VAL:HG12	1:C:140:GLY:N	2.35	0.42
1:C:298:LEU:O	1:C:302:GLU:HG3	2.19	0.42
1:E:247:ILE:HG22	1:E:248:TYR:CD2	2.55	0.42
1:A:287:ILE:HB	1:A:288:PRO:HD3	2.02	0.42
1:E:198:TYR:O	1:E:199:SER:CB	2.68	0.42
1:E:306:ASN:O	1:E:307:ASN:CB	2.68	0.42
1:E:281:LEU:HA	1:E:282:PRO:HD3	1.96	0.42
1:A:272:ARG:O	1:A:273:ASP:CB	2.68	0.41
1:A:170:ALA:O	1:A:171:ASN:HB2	2.20	0.41
1:C:7:PHE:CE2	1:C:80:VAL:HG11	2.55	0.41
1:E:273:ASP:HB2	1:E:281:LEU:HD13	2.02	0.41
1:A:302:GLU:O	1:A:305:VAL:HG12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:ARG:CG	3:A:311:BHK:HAI	2.49	0.41
1:C:272:ARG:O	1:C:273:ASP:CB	2.68	0.41
1:E:180:VAL:HG13	1:E:257:TRP:CZ2	2.56	0.41
3:A:311:BHK:CAH	3:A:311:BHK:CAC	2.99	0.41
1:A:239:LYS:HG3	4:A:318:HOH:O	2.20	0.41
1:C:223:ALA:O	1:C:227:VAL:HG13	2.20	0.41
1:E:285:GLN:HB3	1:E:285:GLN:HE21	1.72	0.40
1:A:139:VAL:CG2	1:A:140:GLY:N	2.84	0.40
1:A:32:LEU:HD11	1:A:100:LEU:HD21	2.03	0.40
1:A:121:SER:HB2	1:C:247:ILE:O	2.21	0.40
1:A:121:SER:CB	1:C:247:ILE:O	2.70	0.40
1:A:270:GLU:OE2	3:A:311:BHK:HAA	2.22	0.40
1:A:74:ILE:CG2	1:A:197:SER:HB2	2.51	0.40
1:E:63:TRP:NE1	1:E:65:ASP:HB3	2.36	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	305/307 (99%)	290 (95%)	10 (3%)	5 (2%)	9	2
1	C	305/307 (99%)	287 (94%)	14 (5%)	4 (1%)	12	3
1	E	305/307 (99%)	290 (95%)	11 (4%)	4 (1%)	12	3
All	All	915/921 (99%)	867 (95%)	35 (4%)	13 (1%)	11	3

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	SER
1	C	199	SER
1	A	134	SER

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Mol	Chain	Res	Type
1	A	205	PRO
1	C	205	PRO
1	E	198	TYR
1	E	199	SER
1	E	205	PRO
1	A	272	ARG
1	C	198	TYR
1	E	272	ARG
1	A	198	TYR
1	C	247	ILE

### 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	264/264 (100%)	257 (97%)	7 (3%)	44	29
1	C	264/264 (100%)	257 (97%)	7 (3%)	44	29
1	E	264/264 (100%)	257 (97%)	7 (3%)	44	29
All	All	792/792 (100%)	771 (97%)	21 (3%)	44	29

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	ASN
1	A	112	ASN
1	A	198	TYR
1	A	205	PRO
1	A	239	LYS
1	A	247	ILE
1	A	248	TYR
1	C	89	ASP
1	C	112	ASN
1	C	168	LYS
1	C	198	TYR
1	C	205	PRO

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Mol	Chain	Res	Type
1	C	248	TYR
1	C	276	ARG
1	E	2	ARG
1	E	31	GLN
1	E	112	ASN
1	E	198	TYR
1	E	205	PRO
1	E	248	TYR
1	E	276	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	37	GLN
1	A	92	GLN
1	A	112	ASN
1	A	171	ASN
1	A	220	ASN
1	A	285	GLN
1	A	303	HIS
1	C	5	ASN
1	C	37	GLN
1	C	92	GLN
1	C	112	ASN
1	C	171	ASN
1	C	221	GLN
1	C	285	GLN
1	E	5	ASN
1	E	31	GLN
1	E	37	GLN
1	E	92	GLN
1	E	112	ASN
1	E	171	ASN
1	E	285	GLN
1	E	307	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	BHK	A	311	2	13,16,16	0.78	0	15,20,20	0.61	0
3	BHK	C	311	2	13,16,16	0.77	0	15,20,20	0.73	0
3	BHK	E	311	2	13,16,16	0.76	0	15,20,20	0.63	0

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	BHK	A	311	2	-	4/10/14/14	0/1/1/1
3	BHK	C	311	2	-	2/10/14/14	0/1/1/1
3	BHK	E	311	2	-	4/10/14/14	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	311	BHK	CAA-CAE-CAF-OAP
3	E	311	BHK	OAQ-CAE-CAF-OAP
3	A	311	BHK	CAA-CAE-CAF-OAP
3	C	311	BHK	CAA-CAE-CAF-OAP
3	C	311	BHK	OAQ-CAE-CAF-OAP
3	A	311	BHK	OAQ-CAE-CAF-OAP
3	E	311	BHK	CAB-CAA-CAE-CAF
3	A	311	BHK	CAB-CAA-CAE-OAQ
3	A	311	BHK	CAB-CAA-CAE-CAF
3	E	311	BHK	CAB-CAA-CAE-OAQ

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	311	BHK	6	0
3	C	311	BHK	1	0
3	E	311	BHK	2	0

## 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	307/307 (100%)	0.03	16 (5%)	27 26	8, 14, 33, 90	3 (0%)
1	C	307/307 (100%)	0.28	19 (6%)	20 20	10, 16, 36, 109	3 (0%)
1	E	307/307 (100%)	0.23	15 (4%)	29 28	9, 16, 33, 94	3 (0%)
All	All	921/921 (100%)	0.18	50 (5%)	25 24	8, 15, 34, 109	9 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	1	ALA	11.8
1	C	307	ASN	8.7
1	E	135	SER	7.7
1	C	137	LEU	7.1
1	E	307	ASN	6.7
1	C	133	THR	6.6
1	A	134	SER	6.1
1	A	307	ASN	6.0
1	C	134	SER	5.9
1	A	137	LEU	5.7
1	E	2	ARG	5.5
1	E	133	THR	5.4
1	A	135	SER	5.1
1	A	2	ARG	5.0
1	C	2	ARG	5.0
1	C	248	TYR	4.9
1	A	306	ASN	4.8
1	E	137	LEU	4.7
1	E	306	ASN	4.2
1	A	133	THR	3.9
1	C	1	ALA	3.7
1	C	132	VAL	3.6
1	E	134	SER	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	57	SER	3.5
1	C	306	ASN	3.2
1	C	58	ASN	3.0
1	C	135	SER	2.9
1	A	197	SER	2.8
1	A	247	ILE	2.8
1	C	197	SER	2.7
1	C	305	VAL	2.6
1	E	205	PRO	2.6
1	E	194	SER	2.6
1	C	194	SER	2.5
1	C	205	PRO	2.5
1	E	31	GLN	2.5
1	A	305	VAL	2.5
1	C	195	ILE	2.4
1	E	197	SER	2.4
1	A	1	ALA	2.4
1	A	194	SER	2.3
1	E	248	TYR	2.3
1	A	205	PRO	2.3
1	E	57	SER	2.3
1	E	138	CYS	2.3
1	A	132	VAL	2.2
1	A	248	TYR	2.2
1	C	136	SER	2.2
1	A	195	ILE	2.1
1	C	206	TYR	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 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( $\text{\AA}^2$ )	Q<0.9
3	BHK	E	311	16/16	0.21	0.65	39,54,56,56	0
3	BHK	A	311	16/16	0.36	0.63	28,37,45,46	0
3	BHK	C	311	16/16	0.52	0.55	28,41,43,43	0
2	ZN	A	1309	1/1	1.00	0.07	12,12,12,12	0
2	ZN	E	1309	1/1	1.00	0.06	13,13,13,13	0
2	ZN	C	1309	1/1	1.00	0.06	15,15,15,15	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.