



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:01 PM GMT

PDB ID : 4ZJZ  
Title : Crystal structure of a benzoate coenzyme A ligase with Benzoyl-AMP  
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Deposited on : 2015-04-29  
Resolution : 1.70 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
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) : rb-20026982

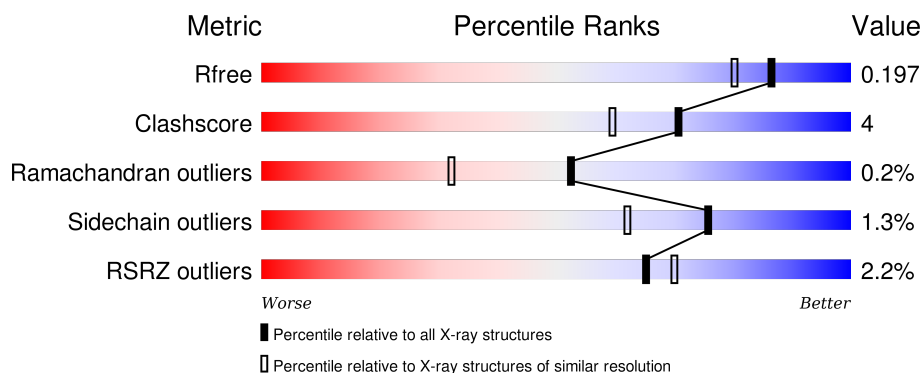
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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	524	<div> <div>2%</div> <div>88%</div> <div>10%</div> <div>..</div> </div>
1	B	524	<div> <div>3%</div> <div>87%</div> <div>11%</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	BEZ	B	601	-	-	-	X
4	GOL	A	1004	-	-	-	X
4	GOL	A	1005	-	-	-	X
4	GOL	B	604	-	-	X	X
4	GOL	B	605	-	-	-	X
4	GOL	B	606	-	-	-	X
4	GOL	B	608	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8658 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.

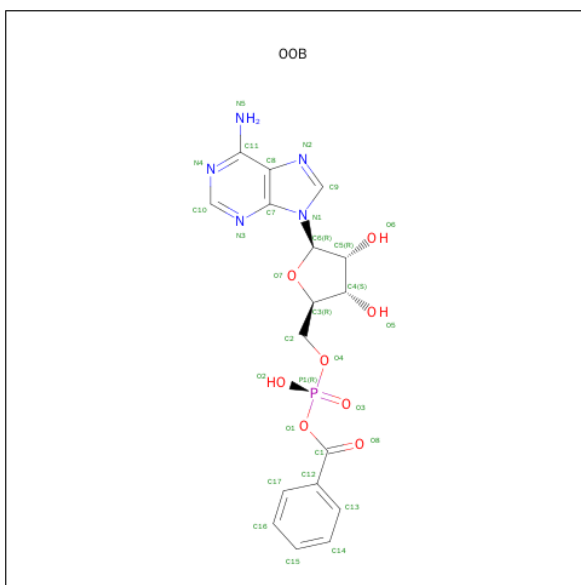
- Molecule 1 is a protein called Benzoate-coenzyme A ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	517	Total	C	N	O	S	0	2	0
			3921	2511	678	722	10			
1	B	519	Total	C	N	O	S	0	5	0
			3950	2531	680	728	11			

There are 8 discrepancies between the modelled and reference sequences:

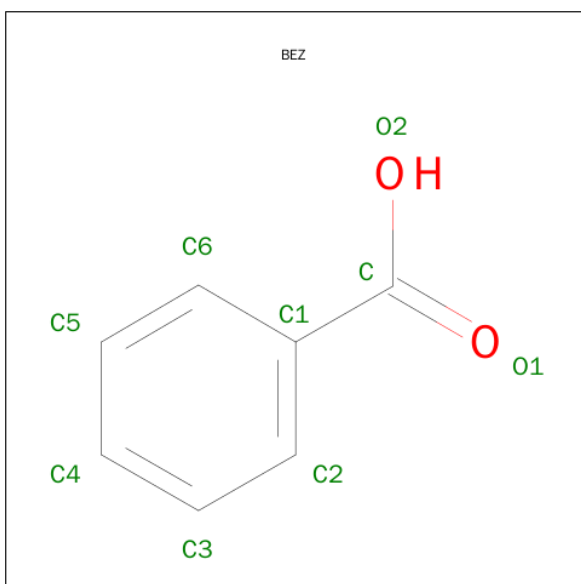
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q93TK0
A	83	ALA	THR	conflict	UNP Q93TK0
A	341	ASP	GLY	conflict	UNP Q93TK0
A	524	GLY	SER	conflict	UNP Q93TK0
B	1	MET	-	initiating methionine	UNP Q93TK0
B	83	ALA	THR	conflict	UNP Q93TK0
B	341	ASP	GLY	conflict	UNP Q93TK0
B	524	GLY	SER	conflict	UNP Q93TK0

- Molecule 2 is 5'-O-[(R)-(benzoyloxy)(hydroxy)phosphoryl]adenosine (three-letter code: OOB) (formula: C<sub>17</sub>H<sub>18</sub>N<sub>5</sub>O<sub>8</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 17	N 5	O 8	P 1	0	0
2	B	1	Total 31	C 17	N 5	O 8	P 1	0	0

- Molecule 3 is BENZOIC ACID (three-letter code: BEZ) (formula:  $C_7H_6O_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 9 7 2	0	0
3	B	1	Total C O 9 7 2	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	316	Total	O	0	0
			316	316		

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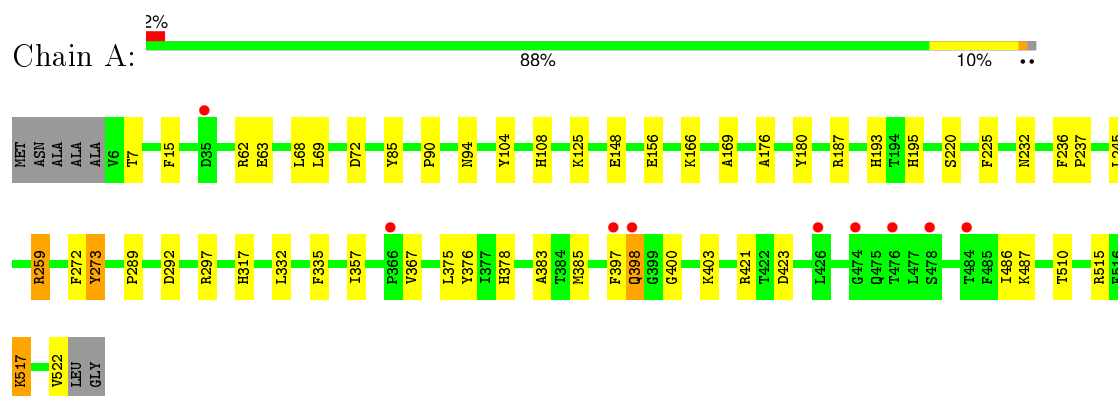
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	331	Total	O	0	0
			331	331		

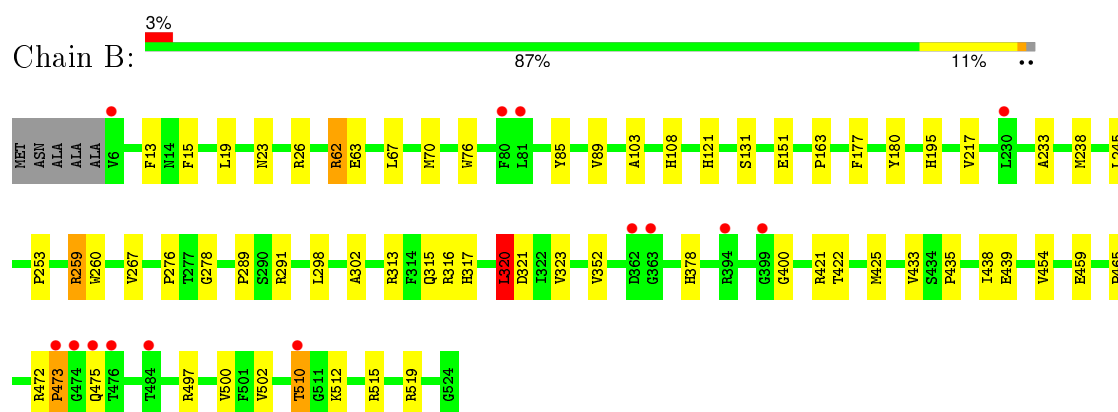
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Benzoate-coenzyme A ligase



- Molecule 1: Benzoate-coenzyme A ligase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.62Å 95.61Å 95.71Å 90.00° 104.53° 90.00°	Depositor
Resolution (Å)	29.39 – 1.70 29.39 – 1.70	Depositor EDS
% Data completeness (in resolution range)	96.3 (29.39-1.70) 96.3 (29.39-1.70)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.35 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.152 , 0.192 0.161 , 0.197	Depositor DCC
$R_{free}$ test set	5429 reflections (5.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	15.2	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 52.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 108571 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8658	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BEZ, GOL, OOB

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	1.32	5/4032 (0.1%)	1.12	11/5500 (0.2%)
1	B	1.31	8/4070 (0.2%)	1.10	4/5552 (0.1%)
All	All	1.31	13/8102 (0.2%)	1.11	15/11052 (0.1%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	13	PHE	CG-CD2	5.86	1.47	1.38
1	A	85	TYR	CD1-CE1	5.61	1.47	1.39
1	B	421	ARG	CG-CD	-5.47	1.38	1.51
1	B	233	ALA	CA-CB	5.46	1.64	1.52
1	A	335	PHE	CE1-CZ	5.34	1.47	1.37
1	B	260	TRP	CZ3-CH2	5.34	1.48	1.40
1	B	302	ALA	CA-CB	5.33	1.63	1.52
1	A	259	ARG	CG-CD	-5.28	1.38	1.51
1	B	259	ARG	CG-CD	-5.21	1.39	1.51
1	A	169	ALA	CA-CB	5.11	1.63	1.52
1	B	85	TYR	CD1-CE1	5.08	1.47	1.39
1	B	352	VAL	CB-CG1	5.04	1.63	1.52
1	A	104	TYR	CE1-CZ	-5.02	1.32	1.38

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	273	TYR	O-C-N	-10.90	104.68	123.20
1	A	273	TYR	CA-C-N	9.87	135.94	116.20
1	A	272	PHE	O-C-N	6.15	132.54	122.70
1	A	272	PHE	CA-C-N	-6.08	103.83	117.20
1	A	69	LEU	CB-CG-CD2	-6.02	100.76	111.00
1	A	292	ASP	CB-CG-OD2	5.61	123.35	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	187	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	A	517	LYS	CD-CE-NZ	-5.48	99.09	111.70
1	A	72	ASP	CB-CG-OD1	5.45	123.21	118.30
1	A	166	LYS	CD-CE-NZ	-5.34	99.42	111.70
1	A	297	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	B	425	MET	O-C-N	5.27	131.13	122.70
1	B	320	LEU	CA-CB-CG	5.24	127.35	115.30
1	B	62	ARG	NE-CZ-NH2	5.14	122.87	120.30
1	B	67	LEU	CB-CG-CD2	-5.08	102.37	111.00

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3921	0	3847	32	0
1	B	3950	0	3884	38	0
2	A	31	0	17	0	0
2	B	31	0	17	0	0
3	A	9	0	5	0	0
3	B	9	0	5	1	0
4	A	24	0	32	0	0
4	B	36	0	48	5	0
5	A	316	0	0	2	0
5	B	331	0	0	4	0
All	All	8658	0	7855	70	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 4.

All (70) 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:B:510:THR:HG21	5:B:964:HOH:O	1.42	1.16

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:385:MET:HA	1:A:397:PHE:CE2	2.06	0.90
1:B:121:HIS:HE1	1:B:151:GLU:OE2	1.56	0.88
1:A:15:PHE:H	1:A:195:HIS:HD2	1.23	0.84
1:B:15:PHE:H	1:B:195:HIS:HD2	1.27	0.83
1:A:62:ARG:O	1:A:63:GLU:HB2	1.84	0.78
1:A:385:MET:CA	1:A:397:PHE:CE2	2.72	0.72
1:A:245:LEU:HD21	1:A:259:ARG:HG2	1.72	0.71
1:B:321:ASP:HB2	4:B:604:GOL:H31	1.74	0.70
1:A:398:GLN:CB	5:A:1386:HOH:O	2.40	0.69
1:A:108:HIS:HD2	1:A:180:TYR:OH	1.76	0.68
1:B:298:LEU:HD12	4:B:604:GOL:H12	1.75	0.68
1:B:62:ARG:O	1:B:63:GLU:HB2	1.94	0.67
1:B:108:HIS:HD2	1:B:180:TYR:OH	1.79	0.66
1:B:15:PHE:H	1:B:195:HIS:CD2	2.14	0.66
1:B:435:PRO:O	1:B:439:GLU:HG3	1.97	0.65
1:B:163:PRO:HG3	5:B:921:HOH:O	1.97	0.63
1:A:515:ARG:HG2	1:A:515:ARG:HH11	1.62	0.63
1:A:289:PRO:O	1:A:317:HIS:HE1	1.81	0.63
1:B:454:VAL:HB	1:B:519:ARG:HG3	1.84	0.59
1:A:15:PHE:H	1:A:195:HIS:CD2	2.12	0.59
1:B:23:ASN:HD22	1:B:26:ARG:HH11	1.50	0.59
1:A:195:HIS:HE1	5:A:1311:HOH:O	1.85	0.59
1:B:121:HIS:HD2	5:B:1012:HOH:O	1.86	0.58
1:B:378:HIS:HD2	1:B:400:GLY:O	1.87	0.58
1:A:385:MET:HA	1:A:397:PHE:HE2	1.69	0.57
1:B:217:VAL:HG21	1:B:267:VAL:CG2	2.34	0.56
1:B:289:PRO:O	1:B:317:HIS:HE1	1.87	0.56
1:A:62:ARG:O	1:A:63:GLU:CB	2.53	0.54
1:B:23:ASN:ND2	1:B:26:ARG:HH11	2.06	0.54
1:B:313:ARG:HG2	1:B:316:ARG:NH2	2.24	0.53
1:B:323:VAL:HG23	4:B:604:GOL:H11	1.90	0.52
1:A:421:ARG:NH2	1:A:423:ASP:OD2	2.34	0.52
1:B:195:HIS:HE1	5:B:902:HOH:O	1.92	0.52
1:B:439:GLU:OE2	1:B:515:ARG:NH1	2.37	0.52
1:B:19:LEU:HB3	1:B:238[A]:MET:HE2	1.92	0.52
1:B:298:LEU:CD1	4:B:604:GOL:H12	2.38	0.52
1:B:62:ARG:O	1:B:63:GLU:CB	2.56	0.52
1:A:383:ALA:O	1:A:397:PHE:CE1	2.63	0.51
1:B:245:LEU:HD21	1:B:259:ARG:HG2	1.93	0.51
1:B:465:PRO:HD2	1:B:497:ARG:HD3	1.95	0.48
1:A:125:LYS:HE3	1:A:148:GLU:OE2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:515:ARG:NH1	1:A:515:ARG:HG2	2.26	0.48
1:B:510:THR:HG22	1:B:512:LYS:H	1.80	0.47
1:B:323:VAL:CG2	4:B:604:GOL:H11	2.45	0.47
1:A:357:ILE:HG23	1:A:375:LEU:CD1	2.45	0.47
1:A:517:LYS:HB3	1:A:522:VAL:HG11	1.95	0.46
1:B:103:ALA:HB1	1:B:131:SER:HB3	1.98	0.46
1:B:459:GLU:CD	1:B:459:GLU:H	2.18	0.46
1:B:121:HIS:CE1	1:B:151:GLU:OE2	2.49	0.46
1:A:236:PHE:HB2	1:A:237:PRO:CD	2.46	0.46
1:A:193:HIS:CE1	1:A:332:LEU:HB2	2.50	0.46
1:A:220:SER:HA	1:A:273:TYR:HB2	1.98	0.45
1:A:385:MET:CA	1:A:397:PHE:HE2	2.25	0.44
1:A:378:HIS:HD2	1:A:400:GLY:O	2.00	0.44
1:A:232:ASN:HB3	1:A:273:TYR:CD2	2.53	0.43
1:B:472:ARG:O	1:B:473:PRO:C	2.57	0.43
1:A:90:PRO:HD2	1:A:176:ALA:O	2.18	0.43
1:A:357:ILE:HG23	1:A:375:LEU:HD11	2.00	0.43
1:B:500:VAL:HG12	1:B:502:VAL:HG13	2.01	0.42
1:B:89:VAL:HG13	1:B:177:PHE:HA	2.02	0.42
1:B:278:GLY:HA3	3:B:601:BEZ:C6	2.51	0.41
1:A:486:ILE:HG13	1:A:487:LYS:N	2.35	0.41
1:B:315:GLN:HA	1:B:320:LEU:O	2.20	0.41
1:A:376:TYR:CE1	1:A:403:LYS:HB2	2.55	0.41
1:A:94:ASN:HA	1:A:225:PHE:CE1	2.55	0.41
1:B:70[B]:MET:HB3	1:B:76:TRP:CD1	2.55	0.41
1:A:68:LEU:C	1:A:68:LEU:HD23	2.41	0.41
1:A:367:VAL:O	1:A:367:VAL:HG23	2.21	0.40
1:B:433:VAL:CG1	1:B:438:ILE:HD11	2.52	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	517/524 (99%)	504 (98%)	12 (2%)	1 (0%)	52	32
1	B	522/524 (100%)	516 (99%)	5 (1%)	1 (0%)	52	32
All	All	1039/1048 (99%)	1020 (98%)	17 (2%)	2 (0%)	52	32

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	398	GLN
1	B	473	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	399/414 (96%)	396 (99%)	3 (1%)	86	79
1	B	404/414 (98%)	397 (98%)	7 (2%)	68	51
All	All	803/828 (97%)	793 (99%)	10 (1%)	76	65

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	THR
1	A	156	GLU
1	A	510	THR
1	B	253	PRO
1	B	276	PRO
1	B	291	ARG
1	B	320	LEU
1	B	422	THR
1	B	475	GLN
1	B	510	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	108	HIS
1	A	195	HIS
1	A	317	HIS
1	A	378	HIS
1	B	23	ASN
1	B	50	GLN
1	B	108	HIS
1	B	112	GLN
1	B	121	HIS
1	B	161	HIS
1	B	195	HIS
1	B	317	HIS
1	B	378	HIS
1	B	411	ASN
1	B	475	GLN

### 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 ⓘ

14 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
2	OOB	A	1000	-	31,34,34	2.10	6 (19%)	33,50,50	1.98	10 (30%)
3	BEZ	A	1001	-	6,9,9	1.61	1 (16%)	8,11,11	1.67	1 (12%)
4	GOL	A	1002	-	5,5,5	0.88	0	5,5,5	1.24	0
4	GOL	A	1003	-	5,5,5	0.19	0	5,5,5	0.64	0
4	GOL	A	1004	-	5,5,5	0.74	0	5,5,5	1.59	1 (20%)
4	GOL	A	1005	-	5,5,5	0.37	0	5,5,5	1.03	0
3	BEZ	B	601	-	6,9,9	1.65	1 (16%)	8,11,11	1.10	0
4	GOL	B	602	-	5,5,5	0.37	0	5,5,5	1.77	2 (40%)
2	OOB	B	603	-	31,34,34	2.16	5 (16%)	33,50,50	2.26	8 (24%)
4	GOL	B	604	-	5,5,5	0.46	0	5,5,5	1.66	2 (40%)
4	GOL	B	605	-	5,5,5	0.59	0	5,5,5	0.50	0
4	GOL	B	606	-	5,5,5	0.60	0	5,5,5	0.99	0
4	GOL	B	607	-	5,5,5	0.33	0	5,5,5	1.30	1 (20%)
4	GOL	B	608	-	5,5,5	0.54	0	5,5,5	1.41	1 (20%)

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
2	OOB	A	1000	-	-	0/15/35/35	0/4/4/4
3	BEZ	A	1001	-	-	0/0/4/4	0/1/1/1
4	GOL	A	1002	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1003	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1004	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1005	-	-	0/4/4/4	0/0/0/0
3	BEZ	B	601	-	-	0/0/4/4	0/1/1/1
4	GOL	B	602	-	-	0/4/4/4	0/0/0/0
2	OOB	B	603	-	-	0/15/35/35	0/4/4/4
4	GOL	B	604	-	-	0/4/4/4	0/0/0/0
4	GOL	B	605	-	-	0/4/4/4	0/0/0/0
4	GOL	B	606	-	-	0/4/4/4	0/0/0/0
4	GOL	B	607	-	-	0/4/4/4	0/0/0/0
4	GOL	B	608	-	-	0/4/4/4	0/0/0/0

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1000	OOB	P1-O2	-3.04	1.42	1.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	603	OOB	C13-C12	2.01	1.42	1.39
3	B	601	BEZ	C3-C2	2.38	1.43	1.38
2	A	1000	OOB	C8-C7	2.38	1.45	1.40
2	A	1000	OOB	C16-C15	2.38	1.44	1.38
2	B	603	OOB	O7-C6	2.42	1.44	1.41
3	A	1001	BEZ	C6-C1	2.68	1.45	1.39
2	A	1000	OOB	P1-O1	2.78	1.66	1.60
2	B	603	OOB	C10-N3	3.08	1.37	1.32
2	B	603	OOB	C9-N2	3.73	1.41	1.34
2	A	1000	OOB	C7-N3	4.41	1.42	1.35
2	A	1000	OOB	O1-C1	8.13	1.45	1.37
2	B	603	OOB	O1-C1	9.28	1.47	1.37

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	603	OOB	N3-C10-N4	-7.03	123.35	128.87
2	B	603	OOB	C6-N1-C7	-6.13	119.96	126.81
2	A	1000	OOB	C6-N1-C7	-6.06	120.04	126.81
2	B	603	OOB	C16-C17-C12	-3.59	116.06	120.35
2	A	1000	OOB	N3-C10-N4	-3.57	126.06	128.87
2	B	603	OOB	O6-C5-C6	-3.11	101.89	111.61
2	A	1000	OOB	C14-C13-C12	-3.07	116.69	120.35
2	A	1000	OOB	C5-C4-C3	-2.99	96.51	102.64
3	A	1001	BEZ	C2-C1-C	-2.84	116.68	120.43
4	B	602	GOL	C3-C2-C1	-2.22	101.71	111.06
2	A	1000	OOB	O1-P1-O3	-2.14	101.33	108.38
4	B	608	GOL	O1-C1-C2	-2.13	99.19	109.97
2	A	1000	OOB	C3-O7-C6	-2.10	107.42	109.64
2	B	603	OOB	C10-N4-C11	2.05	122.42	118.77
2	B	603	OOB	C15-C16-C17	2.06	123.07	120.20
2	A	1000	OOB	N5-C11-N4	2.07	121.98	118.52
4	B	607	GOL	O2-C2-C3	2.09	118.50	108.47
4	B	604	GOL	O3-C3-C2	2.12	120.71	109.97
4	A	1004	GOL	O2-C2-C3	2.32	119.61	108.47
4	B	604	GOL	O2-C2-C3	2.52	120.58	108.47
4	B	602	GOL	O2-C2-C3	2.55	120.71	108.47
2	A	1000	OOB	O1-C1-C12	2.65	115.21	112.39
2	A	1000	OOB	O2-P1-O3	2.77	126.95	112.56
2	B	603	OOB	C17-C12-C13	3.11	122.86	118.61
2	A	1000	OOB	C17-C12-C13	3.35	123.19	118.61
2	B	603	OOB	N5-C11-N4	3.38	124.19	118.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	601	BEZ	1	0
4	B	604	GOL	5	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	517/524 (98%)	-0.27	9 (1%) 73 77	8, 16, 32, 43	0
1	B	519/524 (99%)	-0.22	14 (2%) 58 62	8, 17, 31, 51	0
All	All	1036/1048 (98%)	-0.24	23 (2%) 65 70	8, 16, 32, 51	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	397	PHE	8.8
1	B	476	THR	4.7
1	B	474	GLY	4.5
1	A	426	LEU	4.0
1	B	362	ASP	3.8
1	A	398	GLN	3.7
1	B	399	GLY	3.6
1	A	474	GLY	3.4
1	B	363	GLY	3.3
1	B	473	PRO	3.0
1	B	230	LEU	2.7
1	B	510	THR	2.7
1	A	476	THR	2.7
1	A	478	SER	2.6
1	B	81	LEU	2.5
1	A	35	ASP	2.4
1	B	475	GLN	2.3
1	B	484	THR	2.3
1	B	6	VAL	2.2
1	B	80	PHE	2.1
1	B	394	ARG	2.0
1	A	366	PRO	2.0
1	A	484	THR	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. 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
4	GOL	A	1005	6/6	0.94	0.17	13.90	22,33,43,51	0
4	GOL	B	608	6/6	0.95	0.14	8.36	24,38,40,46	0
4	GOL	A	1004	6/6	0.87	0.15	5.94	36,37,41,42	0
4	GOL	B	604	6/6	0.94	0.13	5.38	13,24,38,38	0
3	BEZ	B	601	9/9	0.91	0.14	3.92	21,28,31,32	0
4	GOL	B	606	6/6	0.91	0.21	3.20	27,37,41,46	0
4	GOL	B	605	6/6	0.95	0.10	2.20	19,23,24,24	0
4	GOL	B	602	6/6	0.95	0.09	0.72	19,26,32,37	0
3	BEZ	A	1001	9/9	0.91	0.11	0.64	20,26,28,28	0
4	GOL	A	1003	6/6	0.91	0.10	0.13	31,38,41,42	0
2	OOB	B	603	31/31	0.97	0.08	-0.13	7,13,17,20	0
4	GOL	A	1002	6/6	0.97	0.09	-0.17	14,20,24,26	0
2	OOB	A	1000	31/31	0.98	0.08	-0.30	8,13,19,19	0
4	GOL	B	607	6/6	0.96	0.07	-0.40	19,26,31,36	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.