



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:26 AM GMT

PDB ID : 3OZ1  
Title : cIAP1-BIR3 domain in complex with the Smac-mimetic compound Smac066  
Authors : Cossu, F.; Malvezzi, F.; Mastrangelo, E.; Canevari, G.; Bolognesi, M.; Milani, M.  
Deposited on : 2010-09-24  
Resolution : 3.00 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

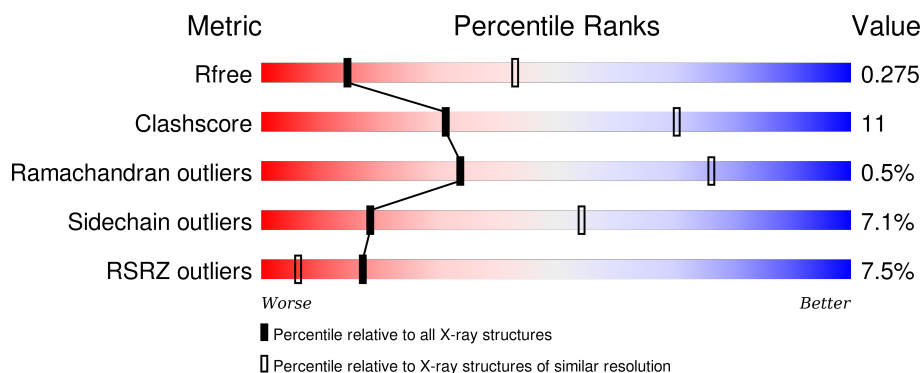
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	122	<div> <div>72%</div> <div>11% • 16%</div> </div>
1	B	122	<div> <div>78%</div> <div>5% • 15%</div> </div>
1	C	122	<div> <div>25%</div> <div>69%</div> <div>10% 6% 16%</div> </div>
1	D	122	<div> <div>25%</div> <div>58%</div> <div>22% • 16%</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	BMB	D	600	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3579 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 Baculoviral IAP repeat-containing protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	103	Total	C	N	O	S	0	0	0
			845	540	145	151	9			
1	B	104	Total	C	N	O	S	0	0	0
			852	544	146	153	9			
1	C	103	Total	C	N	O	S	0	0	0
			850	543	148	150	9			
1	D	102	Total	C	N	O	S	0	0	0
			839	537	144	149	9			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	244	MET	-	EXPRESSION TAG	UNP Q13490
A	358	LEU	-	EXPRESSION TAG	UNP Q13490
A	359	GLU	-	EXPRESSION TAG	UNP Q13490
A	360	HIS	-	EXPRESSION TAG	UNP Q13490
A	361	HIS	-	EXPRESSION TAG	UNP Q13490
A	362	HIS	-	EXPRESSION TAG	UNP Q13490
A	363	HIS	-	EXPRESSION TAG	UNP Q13490
A	364	HIS	-	EXPRESSION TAG	UNP Q13490
A	365	HIS	-	EXPRESSION TAG	UNP Q13490
B	244	MET	-	EXPRESSION TAG	UNP Q13490
B	358	LEU	-	EXPRESSION TAG	UNP Q13490
B	359	GLU	-	EXPRESSION TAG	UNP Q13490
B	360	HIS	-	EXPRESSION TAG	UNP Q13490
B	361	HIS	-	EXPRESSION TAG	UNP Q13490
B	362	HIS	-	EXPRESSION TAG	UNP Q13490
B	363	HIS	-	EXPRESSION TAG	UNP Q13490
B	364	HIS	-	EXPRESSION TAG	UNP Q13490
B	365	HIS	-	EXPRESSION TAG	UNP Q13490
C	244	MET	-	EXPRESSION TAG	UNP Q13490
C	358	LEU	-	EXPRESSION TAG	UNP Q13490
C	359	GLU	-	EXPRESSION TAG	UNP Q13490

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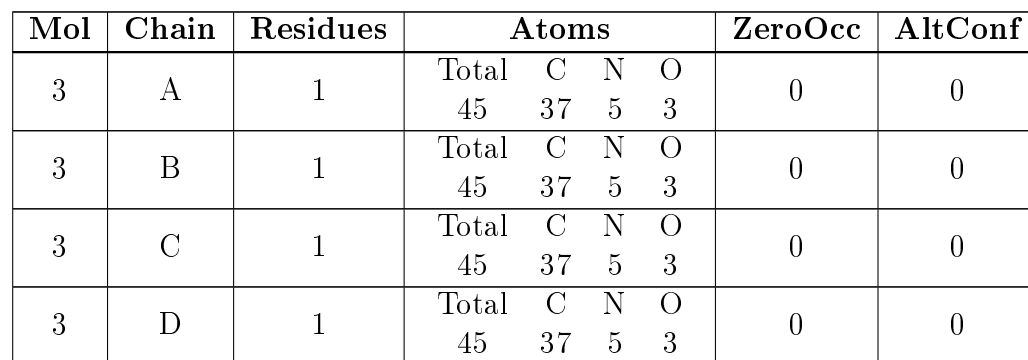
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Chain	Residue	Modelled	Actual	Comment	Reference
C	360	HIS	-	EXPRESSION TAG	UNP Q13490
C	361	HIS	-	EXPRESSION TAG	UNP Q13490
C	362	HIS	-	EXPRESSION TAG	UNP Q13490
C	363	HIS	-	EXPRESSION TAG	UNP Q13490
C	364	HIS	-	EXPRESSION TAG	UNP Q13490
C	365	HIS	-	EXPRESSION TAG	UNP Q13490
D	244	MET	-	EXPRESSION TAG	UNP Q13490
D	358	LEU	-	EXPRESSION TAG	UNP Q13490
D	359	GLU	-	EXPRESSION TAG	UNP Q13490
D	360	HIS	-	EXPRESSION TAG	UNP Q13490
D	361	HIS	-	EXPRESSION TAG	UNP Q13490
D	362	HIS	-	EXPRESSION TAG	UNP Q13490
D	363	HIS	-	EXPRESSION TAG	UNP Q13490
D	364	HIS	-	EXPRESSION TAG	UNP Q13490
D	365	HIS	-	EXPRESSION TAG	UNP Q13490

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

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

- Molecule 3 is (3S,6S,7R,9AS)-7-[2-(BENZYLAMINO)ETHYL]-N-(DIPHENYLMETHYL)-6-{|(2S)-2-(METHYLAMINO)BUTANOYL|AMINO}-5-OXOOCTAHYDRO-1H-PYRROL O[1,2-A]AZEPINE-3-CARBOXAMIDE (three-letter code: BMB) (formula: C<sub>37</sub>H<sub>47</sub>N<sub>5</sub>O<sub>3</sub>).



- | Mol | Chain | Residues | Atoms          | ZeroOcc | AltConf |
|-----|-------|----------|----------------|---------|---------|
| 4   | A     | 1        | Total O<br>1 1 | 0       | 0       |
| 4   | B     | 3        | Total O<br>3 3 | 0       | 0       |
| 4   | C     | 4        | Total O<br>4 4 | 0       | 0       |
| 4   | D     | 1        | Total O<br>1 1 | 0       | 0       |

### 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: Baculoviral IAP repeat-containing protein 2

Chain A: 



- Molecule 1: Baculoviral IAP repeat-containing protein 2

Chain B: 



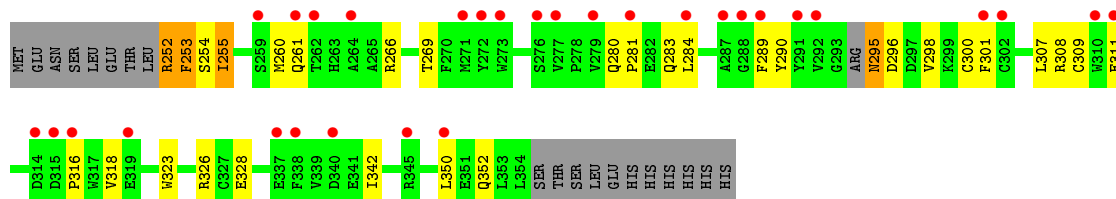
- Molecule 1: Baculoviral IAP repeat-containing protein 2

Chain C: 



- Molecule 1: Baculoviral IAP repeat-containing protein 2

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	113.59Å 114.63Å 92.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.93 – 3.00 60.93 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (60.93-3.00) 95.3 (60.93-3.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.01 (at 3.01Å)	Xtriage
Refinement program	BUSTER 2.9.2	Depositor
R, $R_{free}$	0.192 , 0.257 0.210 , 0.275	Depositor DCC
$R_{free}$ test set	575 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	69.4	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 102.3	EDS
Estimated twinning fraction	0.048 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 11908 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	3579	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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: BMB, 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	0.51	0/872	0.69	0/1180
1	B	0.53	0/879	0.69	0/1190
1	C	0.58	0/877	0.72	0/1186
1	D	0.50	0/865	0.70	0/1169
All	All	0.53	0/3493	0.70	0/4725

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	845	0	790	8	0
1	B	852	0	797	8	0
1	C	850	0	798	34	0
1	D	839	0	784	25	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	45	0	47	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	45	0	47	0	0
3	C	45	0	47	1	0
3	D	45	0	47	4	0
4	A	1	0	0	0	0
4	B	3	0	0	0	0
4	C	4	0	0	0	0
4	D	1	0	0	0	0
All	All	3579	0	3357	74	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 11.

All (74) 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:350:LEU:HD12	1:C:350:LEU:C	1.57	1.18
1:C:252:ARG:HH11	1:C:252:ARG:CG	1.57	1.15
1:D:254:SER:C	1:D:255:ILE:HD13	1.69	1.11
1:D:255:ILE:HD13	1:D:255:ILE:N	1.66	1.10
1:C:252:ARG:HG2	1:C:252:ARG:HH11	1.17	1.07
1:C:252:ARG:HH11	1:C:252:ARG:CB	1.69	1.04
1:C:252:ARG:HB3	1:C:252:ARG:NH1	1.72	1.03
1:C:350:LEU:HD12	1:C:350:LEU:O	1.60	1.01
1:C:350:LEU:C	1:C:350:LEU:CD1	2.30	1.00
1:D:254:SER:C	1:D:255:ILE:CD1	2.30	1.00
1:C:252:ARG:CB	1:C:252:ARG:NH1	2.30	0.93
1:D:255:ILE:CD1	1:D:255:ILE:N	2.30	0.88
1:C:252:ARG:HG2	1:C:252:ARG:NH1	1.88	0.87
1:B:345:ARG:HG3	1:B:345:ARG:HH11	1.42	0.85
1:C:252:ARG:O	1:C:253:PHE:CD2	2.30	0.84
1:D:253:PHE:HB3	1:D:255:ILE:HD11	1.62	0.80
1:C:350:LEU:CD1	1:C:350:LEU:O	2.30	0.77
1:C:350:LEU:HD11	1:C:354:LEU:HD13	1.70	0.73
1:C:252:ARG:O	1:C:253:PHE:HD2	1.72	0.72
1:A:332:ARG:HH11	1:B:331:ILE:HG21	1.55	0.72
1:C:350:LEU:CD1	1:C:354:LEU:HD13	2.21	0.71
1:D:254:SER:O	1:D:255:ILE:HD12	1.92	0.70
1:C:296:ASP:C	1:C:296:ASP:OD1	2.30	0.68
1:D:254:SER:C	1:D:255:ILE:HD12	2.14	0.67
1:D:283:GLN:HB3	1:D:316:PRO:HG2	1.79	0.64
1:D:253:PHE:CZ	1:D:328:GLU:HB3	2.33	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:LYS:HG2	1:A:349:LEU:HD21	1.82	0.62
1:B:345:ARG:NH1	1:B:345:ARG:HG3	2.12	0.62
1:C:295:ASN:O	1:C:296:ASP:CB	2.48	0.62
1:B:322:LYS:HG2	1:B:349:LEU:HD21	1.85	0.58
1:C:349:LEU:O	1:C:352:GLN:HG2	2.04	0.57
1:C:352:GLN:HG3	1:C:353:LEU:N	2.19	0.57
1:D:284:LEU:HD22	1:D:298:VAL:HG11	1.87	0.57
1:D:254:SER:CA	1:D:255:ILE:HD13	2.35	0.57
1:C:252:ARG:HB3	1:C:252:ARG:CZ	2.34	0.57
1:D:308:ARG:HB3	3:D:600:BMB:HAK	1.86	0.56
1:C:252:ARG:C	1:C:253:PHE:CD2	2.80	0.55
1:C:258:LEU:O	1:C:261:GLN:HB2	2.09	0.52
1:C:350:LEU:HD12	1:C:351:GLU:N	2.21	0.52
1:C:350:LEU:HB2	1:D:350:LEU:HD13	1.91	0.52
1:A:348:HIS:HB3	1:A:351:GLU:HG2	1.92	0.52
1:C:295:ASN:O	1:C:296:ASP:OD1	2.29	0.50
1:D:323:TRP:CE2	3:D:600:BMB:HBO	2.46	0.49
1:D:307:LEU:HD22	3:D:600:BMB:HAAB	1.93	0.48
1:D:308:ARG:HH12	3:D:600:BMB:CAT	2.27	0.48
1:B:354:LEU:HD12	1:B:354:LEU:HA	1.38	0.48
1:C:349:LEU:HA	1:C:349:LEU:HD12	1.70	0.48
1:D:295:ASN:HB3	1:D:296:ASP:H	1.56	0.48
1:C:253:PHE:O	1:C:254:SER:OG	2.30	0.47
1:D:318:VAL:HG13	1:D:342:ILE:HD12	1.97	0.47
3:A:600:BMB:HAR	3:A:600:BMB:CAT	2.44	0.46
1:A:258:LEU:O	1:A:261:GLN:HB2	2.15	0.46
1:B:318:VAL:HG13	1:B:342:ILE:HD12	1.97	0.45
1:B:349:LEU:HD12	1:B:349:LEU:HA	1.76	0.45
1:D:252:ARG:HB2	1:D:253:PHE:H	1.51	0.45
1:C:295:ASN:O	1:C:296:ASP:HB3	2.16	0.45
1:C:349:LEU:HD12	1:C:352:GLN:NE2	2.30	0.45
1:C:351:GLU:OE2	1:D:326:ARG:NH1	2.50	0.45
1:C:294:ARG:HG3	1:C:295:ASN:HB2	1.98	0.45
1:C:295:ASN:O	1:C:296:ASP:CG	2.55	0.45
1:D:260:MET:HB3	1:D:266:ARG:HG2	1.98	0.44
1:D:280:GLN:HG2	1:D:283:GLN:HE22	1.82	0.44
1:A:282:GLU:H	1:A:282:GLU:HG3	1.46	0.44
1:A:284:LEU:O	1:A:289:PHE:HB2	2.17	0.44
1:A:318:VAL:HG13	1:A:342:ILE:HD12	1.99	0.44
1:D:290:TYR:HE2	1:D:301:PHE:HA	1.84	0.43
3:C:600:BMB:CAT	3:C:600:BMB:HAR	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:289:PHE:HB3	1:D:298:VAL:HB	2.01	0.42
1:C:296:ASP:CG	1:C:296:ASP:O	2.58	0.42
1:C:318:VAL:HG13	1:C:342:ILE:HD12	2.01	0.42
1:D:289:PHE:HA	1:D:300:CYS:HA	2.03	0.41
1:B:322:LYS:NZ	1:B:349:LEU:HD11	2.34	0.41
1:A:349:LEU:HA	1:A:352:GLN:HB2	2.02	0.40
1:C:332:ARG:HH11	1:C:332:ARG:HB3	1.86	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	101/122 (83%)	96 (95%)	5 (5%)	0	100	100
1	B	102/122 (84%)	99 (97%)	3 (3%)	0	100	100
1	C	101/122 (83%)	95 (94%)	5 (5%)	1 (1%)	19	61
1	D	98/122 (80%)	89 (91%)	8 (8%)	1 (1%)	19	61
All	All	402/488 (82%)	379 (94%)	21 (5%)	2 (0%)	34	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	281	PRO
1	C	255	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	91/110 (83%)	87 (96%)	4 (4%)	35	74
1	B	92/110 (84%)	87 (95%)	5 (5%)	27	66
1	C	91/110 (83%)	83 (91%)	8 (9%)	12	42
1	D	90/110 (82%)	81 (90%)	9 (10%)	9	34
All	All	364/440 (83%)	338 (93%)	26 (7%)	18	54

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	253	PHE
1	A	282	GLU
1	A	332	ARG
1	A	353	LEU
1	B	254	SER
1	B	345	ARG
1	B	349	LEU
1	B	354	LEU
1	B	355	SER
1	C	252	ARG
1	C	271	MET
1	C	294	ARG
1	C	296	ASP
1	C	349	LEU
1	C	350	LEU
1	C	352	GLN
1	C	354	LEU
1	D	252	ARG
1	D	253	PHE
1	D	255	ILE
1	D	261	GLN
1	D	269	THR
1	D	295	ASN
1	D	309	CYS
1	D	311	GLU
1	D	352	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	352	GLN
1	C	352	GLN
1	D	257	ASN
1	D	280	GLN
1	D	283	GLN

### 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](#)

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	BMB	A	600	-	47,49,49	0.65	0	53,66,66	1.02	3 (5%)
3	BMB	B	600	-	47,49,49	0.70	0	53,66,66	1.16	4 (7%)
3	BMB	C	600	-	47,49,49	0.73	0	53,66,66	1.07	3 (5%)
3	BMB	D	600	-	47,49,49	0.76	0	53,66,66	1.53	9 (16%)

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	BMB	A	600	-	-	0/35/65/65	0/5/5/5
3	BMB	B	600	-	-	0/35/65/65	0/5/5/5
3	BMB	C	600	-	-	0/35/65/65	0/5/5/5
3	BMB	D	600	-	-	0/35/65/65	0/5/5/5

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	600	BMB	OAE-CBI-NBS	-3.53	117.28	121.67
3	B	600	BMB	OAE-CBI-NBS	-3.34	117.53	121.67
3	D	600	BMB	OAE-CBI-NBS	-2.83	118.15	121.67
3	A	600	BMB	OAE-CBI-NBS	-2.78	118.22	121.67
3	B	600	BMB	CBB-NBD-CAV	-2.03	106.00	113.48
3	D	600	BMB	CBH-CBP-NBS	2.13	115.55	111.84
3	A	600	BMB	CBM-CBQ-NBE	2.14	113.42	110.70
3	D	600	BMB	CBI-CBQ-NBE	2.34	113.84	108.47
3	C	600	BMB	CBM-CBQ-NBE	2.54	113.92	110.70
3	A	600	BMB	CBP-NBS-CBI	2.63	121.58	118.27
3	B	600	BMB	CBP-NBS-CBI	2.70	121.67	118.27
3	D	600	BMB	CBR-NBF-CBH	2.77	127.03	121.18
3	C	600	BMB	CBP-NBS-CBI	2.94	121.97	118.27
3	B	600	BMB	CBM-CBQ-NBE	2.97	114.46	110.70
3	D	600	BMB	CAZ-CBO-NBS	2.97	105.07	101.93
3	D	600	BMB	CBA-CBP-NBS	3.66	106.37	102.92
3	D	600	BMB	CBL-CBR-NBF	3.86	118.39	111.18
3	D	600	BMB	CBP-NBS-CBI	4.21	123.57	118.27
3	D	600	BMB	CBM-CBQ-NBE	4.57	116.48	110.70

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	600	BMB	1	0
3	C	600	BMB	1	0
3	D	600	BMB	4	0

## 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	103/122 (84%)	0.18	0 100 100	45, 71, 114, 131	4 (3%)
1	B	104/122 (85%)	-0.06	0 100 100	41, 64, 97, 122	3 (2%)
1	C	103/122 (84%)	-0.02	1 (0%) 84 60	41, 61, 96, 120	3 (2%)
1	D	102/122 (83%)	1.50	30 (29%) 1 0	77, 122, 167, 179	9 (8%)
All	All	412/488 (84%)	0.40	31 (7%) 17 6	41, 74, 149, 179	19 (4%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	271	MET	6.8
1	D	314	ASP	5.1
1	D	279	VAL	4.8
1	D	284	LEU	4.8
1	D	273	TRP	4.1
1	D	281	PRO	4.0
1	D	264	ALA	3.8
1	D	310	TRP	3.6
1	D	311	GLU	3.5
1	D	291	TYR	3.4
1	D	277	VAL	3.3
1	D	276	SER	3.1
1	D	340	ASP	3.1
1	D	292	VAL	3.0
1	D	262	THR	2.8
1	D	315	ASP	2.6
1	D	288	GLY	2.6
1	D	289	PHE	2.5
1	D	259	SER	2.5
1	D	302	CYS	2.4
1	D	316	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	301	PHE	2.4
1	D	345	ARG	2.3
1	D	272	TYR	2.3
1	D	319	GLU	2.3
1	D	338	PHE	2.2
1	D	261	GLN	2.2
1	D	337	GLU	2.2
1	D	350	LEU	2.1
1	D	287	ALA	2.1
1	C	350	LEU	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
3	BMB	A	600	45/45	0.95	0.34	1.33	16,74,197,262	0
3	BMB	B	600	45/45	0.96	0.26	1.26	16,76,239,281	0
3	BMB	C	600	45/45	0.92	0.27	0.55	19,75,216,277	0
3	BMB	D	600	45/45	0.89	0.45	0.36	62,118,264,284	0
2	ZN	C	501	1/1	1.00	0.19	-0.39	53,53,53,53	0
2	ZN	A	501	1/1	0.99	0.17	-0.78	59,59,59,59	0
2	ZN	B	501	1/1	0.99	0.17	-1.05	55,55,55,55	0
2	ZN	D	501	1/1	0.99	0.13	-1.41	100,100,100,100	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.