



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 08:26 AM GMT

PDB ID : 3ENH  
Title : Crystal structure of Cgi121/Bud32/Kae1 complex  
Authors : Neculai, D.  
Deposited on : 2008-09-25  
Resolution : 3.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

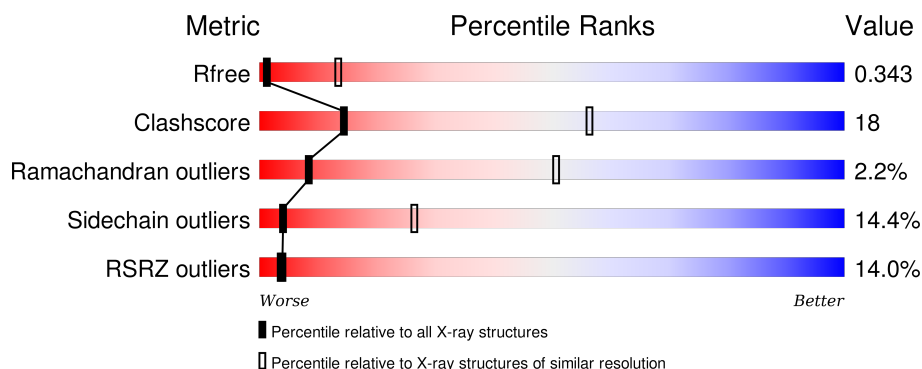
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.60 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))
$R_{free}$	91344	1408 (3.80-3.40)
Clashscore	102246	1010 (3.74-3.46)
Ramachandran outliers	100387	1007 (3.76-3.44)
Sidechain outliers	100360	1007 (3.76-3.44)
RSRZ outliers	91569	1003 (3.78-3.42)

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	540	 7% 50% 34% 7% 8%
1	B	540	 16% 52% 32% 5% 10%
2	C	150	 28% 56% 30% 5% 8%
2	D	150	 5% 53% 33% 5% 8%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	TBR	A	4769	-	-	-	X
3	TBR	A	5769	-	-	X	-
3	TBR	B	2769	-	-	-	X
3	TBR	B	7769	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10142 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 Putative O-sialoglycoprotein endopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	495	Total	C	N	O	S	0	0	0
			3926	2509	661	736	20			
1	B	485	Total	C	N	O	S	0	0	0
			3856	2466	643	726	21			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP Q58530
A	-3	ALA	-	EXPRESSION TAG	UNP Q58530
A	-2	MET	-	EXPRESSION TAG	UNP Q58530
A	-1	ASP	-	EXPRESSION TAG	UNP Q58530
A	0	PRO	-	EXPRESSION TAG	UNP Q58530
B	-4	GLY	-	EXPRESSION TAG	UNP Q58530
B	-3	ALA	-	EXPRESSION TAG	UNP Q58530
B	-2	MET	-	EXPRESSION TAG	UNP Q58530
B	-1	ASP	-	EXPRESSION TAG	UNP Q58530
B	0	PRO	-	EXPRESSION TAG	UNP Q58530

- Molecule 2 is a protein called Uncharacterized protein MJ0187.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	138	Total	C	N	O	S	0	0	0
			1108	708	196	200	4			
2	D	138	Total	C	N	O	S	0	0	0
			1108	708	196	200	4			

There are 10 discrepancies between the modelled and reference sequences:

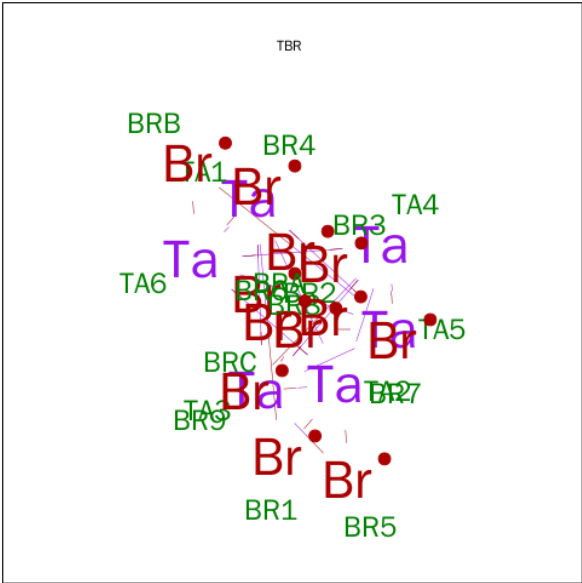
Chain	Residue	Modelled	Actual	Comment	Reference
C	1	GLY	-	EXPRESSION TAG	UNP Q57646
C	2	ALA	-	EXPRESSION TAG	UNP Q57646

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Chain	Residue	Modelled	Actual	Comment	Reference
C	3	MET	-	EXPRESSION TAG	UNP Q57646
C	4	ASP	-	EXPRESSION TAG	UNP Q57646
C	5	PRO	-	EXPRESSION TAG	UNP Q57646
D	1	GLY	-	EXPRESSION TAG	UNP Q57646
D	2	ALA	-	EXPRESSION TAG	UNP Q57646
D	3	MET	-	EXPRESSION TAG	UNP Q57646
D	4	ASP	-	EXPRESSION TAG	UNP Q57646
D	5	PRO	-	EXPRESSION TAG	UNP Q57646

- Molecule 3 is HEXATANTALUM DODECABROMIDE (three-letter code: TBR) (formula: Br<sub>12</sub>Ta<sub>6</sub>).



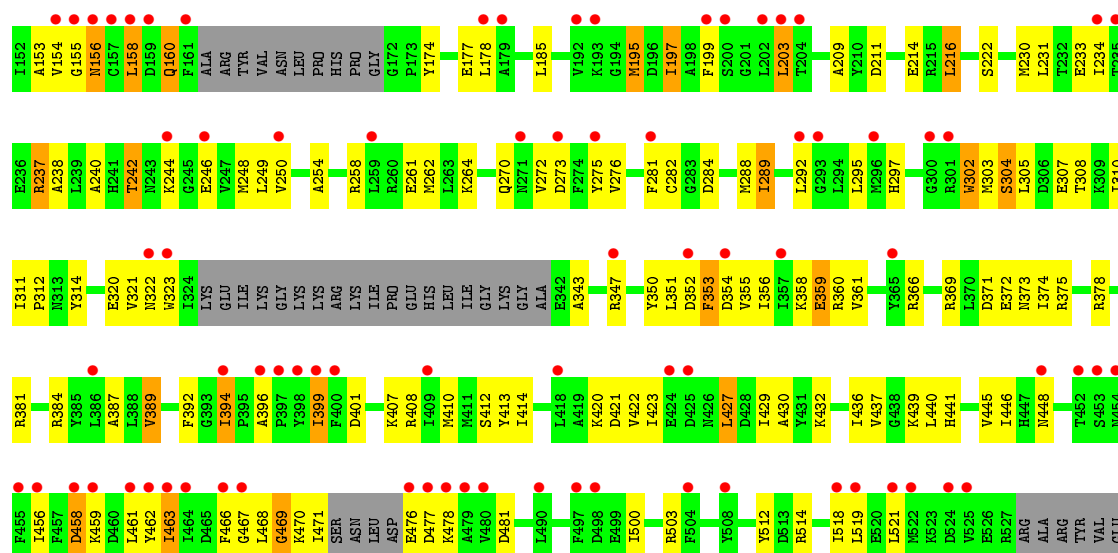
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Br	Ta	0	0
			18	12	6		
3	B	1	Total	Br	Ta	0	0
			18	12	6		
3	B	1	Total	Br	Ta	0	0
			18	12	6		
3	A	1	Total	Br	Ta	0	0
			18	12	6		
3	A	1	Total	Br	Ta	0	0
			18	12	6		
3	A	1	Total	Br	Ta	0	0
			18	12	6		
3	B	1	Total	Br	Ta	0	0
			18	12	6		

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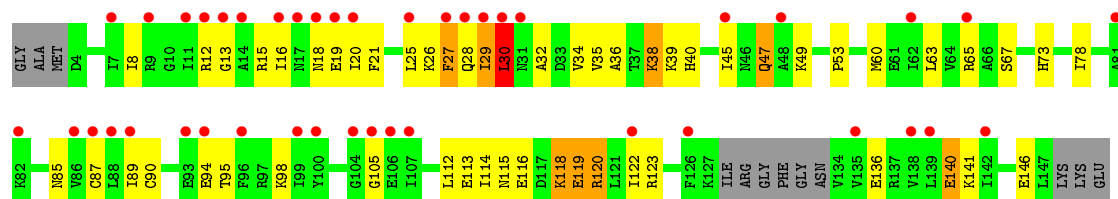
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	Br	Ta		
3	B	1	18	12	6	0	0

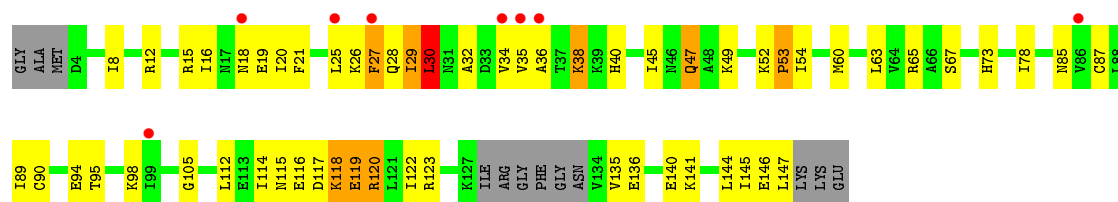




• Molecule 2: Uncharacterized protein MJ0187



• Molecule 2: Uncharacterized protein MJ0187





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	76.99Å 106.91Å 209.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	94.92 – 3.60 72.26 – 3.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (94.92-3.60) 100.0 (72.26-3.60)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.89 (at 3.58Å)	Xtriage
Refinement program	REFMAC 5.5	Depositor
R, $R_{free}$	0.271 , 0.324 0.287 , 0.343	Depositor DCC
$R_{free}$ test set	1052 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	146.1	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 144.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 20739 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	10142	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	104.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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: TBR

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	0.35	0/3994	0.56	0/5387
1	B	0.35	0/3921	0.54	0/5284
2	C	0.36	0/1119	0.59	1/1499 (0.1%)
2	D	0.37	0/1119	0.59	1/1499 (0.1%)
All	All	0.35	0/10153	0.56	2/13669 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	30	LEU	CA-CB-CG	5.80	128.65	115.30
2	C	30	LEU	CA-CB-CG	5.72	128.47	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	3926	0	3959	168	0
1	B	3856	0	3887	140	0
2	C	1108	0	1168	31	0
2	D	1108	0	1168	34	0
3	A	72	0	0	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	72	0	0	5	0
All	All	10142	0	10182	365	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 18.

The worst 5 of 365 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:11:GLU:OE2	3:A:5769:TBR:BR5	2.09	1.25
2:C:114:ILE:HD11	2:C:119:GLU:HB3	1.33	1.09
2:D:114:ILE:HD11	2:D:119:GLU:HB3	1.35	1.08
1:A:394:ILE:HD12	1:A:436:ILE:HB	1.47	0.96
1:B:394:ILE:HD12	1:B:436:ILE:HB	1.47	0.95

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	487/540 (90%)	423 (87%)	57 (12%)	7 (1%)	14	59
1	B	475/540 (88%)	421 (89%)	50 (10%)	4 (1%)	24	69
2	C	134/150 (89%)	115 (86%)	11 (8%)	8 (6%)	2	24
2	D	134/150 (89%)	115 (86%)	11 (8%)	8 (6%)	2	24
All	All	1230/1380 (89%)	1074 (87%)	129 (10%)	27 (2%)	8	51

5 of 27 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	316	THR

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Mol	Chain	Res	Type
1	B	240	ALA
2	D	21	PHE
1	A	467	GLY
2	C	21	PHE

### 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	421/460 (92%)	357 (85%)	64 (15%)	3	24
1	B	415/460 (90%)	360 (87%)	55 (13%)	5	30
2	C	120/128 (94%)	102 (85%)	18 (15%)	3	25
2	D	120/128 (94%)	102 (85%)	18 (15%)	3	25
All	All	1076/1176 (92%)	921 (86%)	155 (14%)	4	26

5 of 155 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	93	LEU
1	B	261	GLU
2	D	47	GLN
1	B	133	THR
1	B	195	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 24 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	297	HIS
1	B	27	ASN
2	D	46	ASN
1	A	313	ASN
1	A	441	HIS

### 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 ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	TBR	A	1769	-	0,36,36	0.00	-	0,180,180	0.00	-
3	TBR	A	4769	-	0,36,36	0.00	-	0,180,180	0.00	-
3	TBR	A	5769	-	0,36,36	0.00	-	0,180,180	0.00	-
3	TBR	A	6769	-	0,36,36	0.00	-	0,180,180	0.00	-
3	TBR	B	2769	-	0,36,36	0.00	-	0,180,180	0.00	-
3	TBR	B	3769	-	0,36,36	0.00	-	0,180,180	0.00	-
3	TBR	B	7769	-	0,36,36	0.00	-	0,180,180	0.00	-
3	TBR	B	8769	-	0,36,36	0.00	-	0,180,180	0.00	-

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	TBR	A	1769	-	-	0/0/696/696	0/0/19/19
3	TBR	A	4769	-	-	0/0/696/696	0/0/19/19

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TBR	A	5769	-	-	0/0/696/696	0/0/19/19
3	TBR	A	6769	-	-	0/0/696/696	0/0/19/19
3	TBR	B	2769	-	-	0/0/696/696	0/0/19/19
3	TBR	B	3769	-	-	0/0/696/696	0/0/19/19
3	TBR	B	7769	-	-	0/0/696/696	0/0/19/19
3	TBR	B	8769	-	-	0/0/696/696	0/0/19/19

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	5769	TBR	12	0
3	A	6769	TBR	3	0
3	B	2769	TBR	1	0
3	B	3769	TBR	2	0
3	B	7769	TBR	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	495/540 (91%)	0.56	38 (7%) 16 11	70, 102, 104, 115	2 (0%)
1	B	485/540 (89%)	0.94	88 (18%) 2 2	99, 103, 120, 126	6 (1%)
2	C	138/150 (92%)	1.22	42 (30%) 1 1	101, 106, 114, 116	1 (0%)
2	D	138/150 (92%)	0.48	8 (5%) 26 18	101, 106, 114, 116	2 (1%)
All	All	1256/1380 (91%)	0.77	176 (14%) 4 3	70, 103, 114, 126	11 (0%)

The worst 5 of 176 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	GLY	9.6
1	B	461	LEU	6.4
1	B	455	PHE	6.1
1	B	463	ILE	5.9
1	A	9	THR	5.5

### 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	TBR	A	4769	18/18	0.96	0.43	1.84	151,222,500,500	18
3	TBR	A	6769	18/18	0.97	0.39	0.96	83,119,191,196	18
3	TBR	B	7769	18/18	0.84	0.45	0.90	79,159,374,500	18
3	TBR	B	2769	18/18	0.98	0.43	0.08	25,81,98,244	18
3	TBR	B	3769	18/18	0.98	0.39	-0.50	122,181,265,500	18
3	TBR	A	5769	18/18	0.90	0.28	-0.98	72,127,269,500	18
3	TBR	A	1769	18/18	0.99	0.33	-1.08	67,105,143,170	18
3	TBR	B	8769	18/18	0.87	0.63	-	51,146,180,230	18

## 6.5 Other polymers [i](#)

There are no such residues in this entry.