



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 10:54 AM UTC

PDB ID : 3VAL / pdb\_00003val  
Title : Structure of U2AF65 variant with BrU5C1 DNA  
Authors : Jenkins, J.L.; Frato, K.H.; Kielkopf, C.L.  
Deposited on : 2011-12-29  
Resolution : 2.50 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

<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 : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

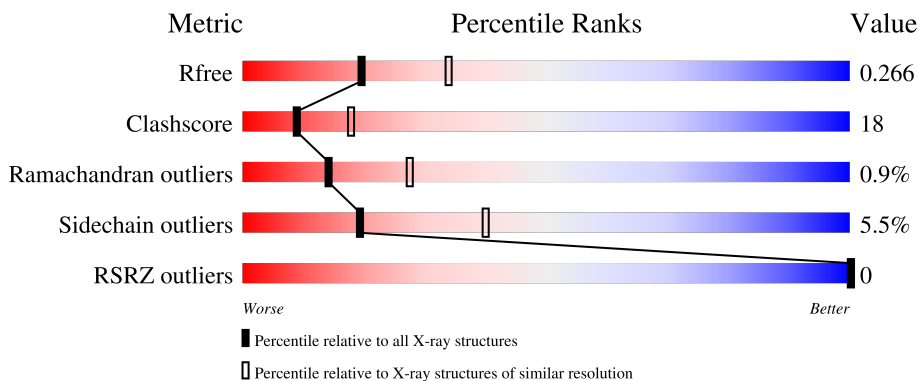
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	174	
1	B	174	
1	D	174	
1	I	174	
2	E	7	

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Mol	Chain	Length	Quality of chain
2	H	7	 86% 14%
2	K	7	 71% 29%
2	P	7	 29% 29% 29% 14%

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6102 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 Splicing factor U2AF 65 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	174	1343	857	230	250	6	0	0	0
1	B	173	1339	855	229	249	6	0	0	0
1	D	174	1343	857	230	250	6	0	0	0
1	I	172	1332	850	228	248	6	0	0	0

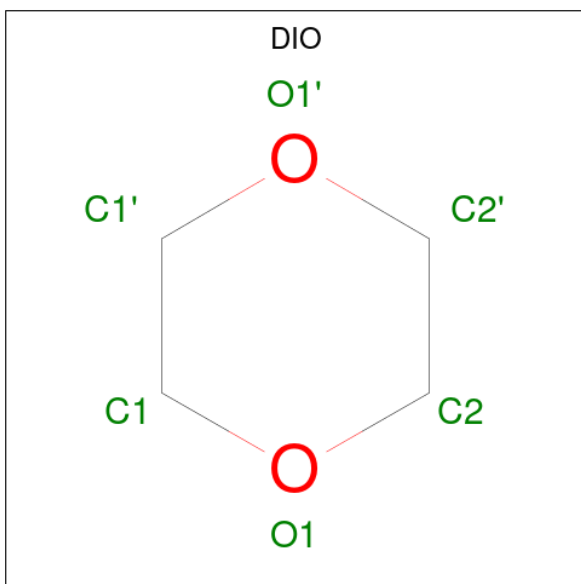
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	GLY	-	expression tag	UNP P26368
A	144	PRO	-	expression tag	UNP P26368
A	145	LEU	-	expression tag	UNP P26368
A	146	GLY	-	expression tag	UNP P26368
A	147	SER	-	expression tag	UNP P26368
B	143	GLY	-	expression tag	UNP P26368
B	144	PRO	-	expression tag	UNP P26368
B	145	LEU	-	expression tag	UNP P26368
B	146	GLY	-	expression tag	UNP P26368
B	147	SER	-	expression tag	UNP P26368
D	143	GLY	-	expression tag	UNP P26368
D	144	PRO	-	expression tag	UNP P26368
D	145	LEU	-	expression tag	UNP P26368
D	146	GLY	-	expression tag	UNP P26368
D	147	SER	-	expression tag	UNP P26368
I	143	GLY	-	expression tag	UNP P26368
I	144	PRO	-	expression tag	UNP P26368
I	145	LEU	-	expression tag	UNP P26368
I	146	GLY	-	expression tag	UNP P26368
I	147	SER	-	expression tag	UNP P26368

- Molecule 2 is a DNA chain called DNA (5'-D(\*C\*UP\*UP\*UP\*(BRU)P\*UP\*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	Br	C	N	O				P
2	E	6	Total 112	Br 1	C 54	N 12	O 40	P 5	0	0	0
2	H	7	Total 116	Br 1	C 54	N 12	O 43	P 6	0	0	1
2	K	7	Total 116	Br 1	C 54	N 12	O 43	P 6	0	0	1
2	P	6	Total 97	Br 1	C 45	N 10	O 36	P 5	0	0	1

- Molecule 3 is 1,4-DIETHYLENE DIOXIDE (CCD ID: DIO) (formula: C<sub>4</sub>H<sub>8</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	C O		
3	A	1	Total 6	C 4 O 2	0	0
3	B	1	Total 6	C 4 O 2	0	0
3	D	1	Total 6	C 4 O 2	0	0
3	I	1	Total 6	C 4 O 2	0	0

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	68	Total	O	0	0
			68	68		

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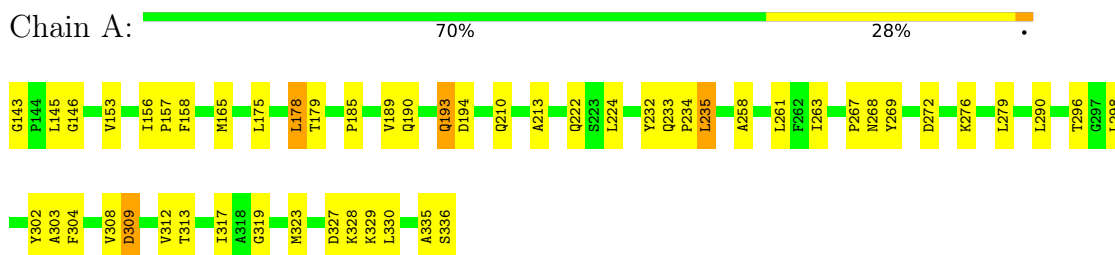
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
5	B	29	Total O 29 29	0	0
5	D	49	Total O 49 49	0	0
5	I	43	Total O 43 43	0	0
5	E	8	Total O 8 8	0	0
5	H	12	Total O 12 12	0	0
5	K	10	Total O 10 10	0	0
5	P	6	Total O 6 6	0	0

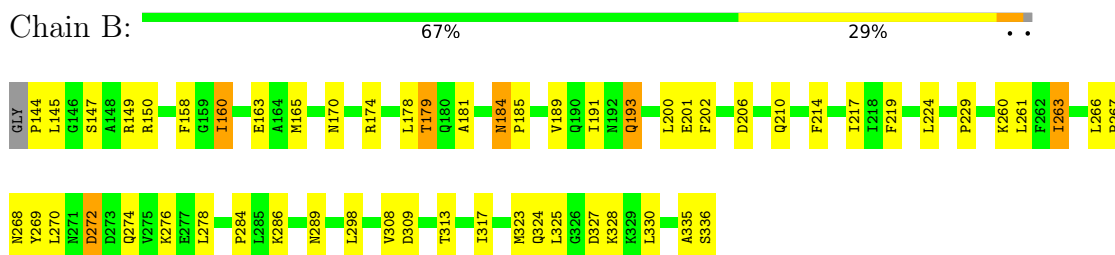
### 3 Residue-property plots [i](#)

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.

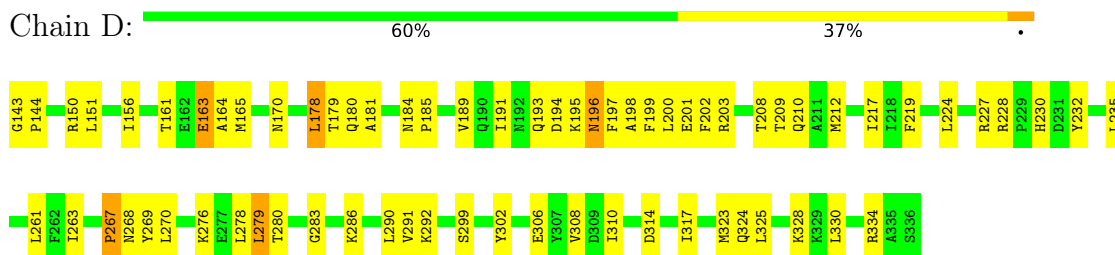
- Molecule 1: Splicing factor U2AF 65 kDa subunit



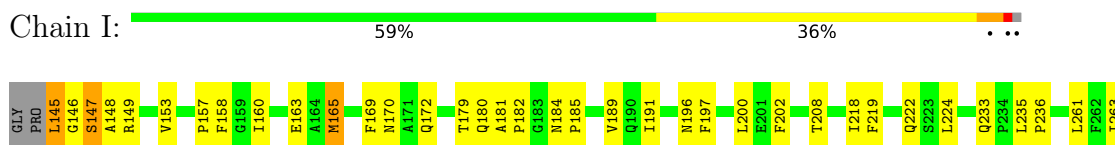
- Molecule 1: Splicing factor U2AF 65 kDa subunit

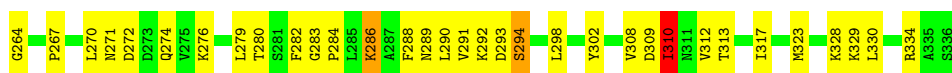


- Molecule 1: Splicing factor U2AF 65 kDa subunit

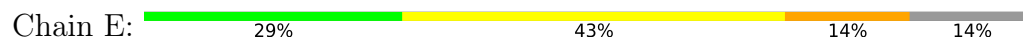


- Molecule 1: Splicing factor U2AF 65 kDa subunit

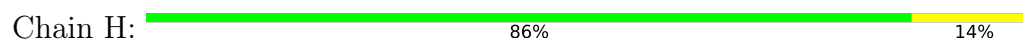




• Molecule 2: DNA (5'-D(\*C\*UP\*UP\*UP\*(BRU)P\*UP\*U)-3')



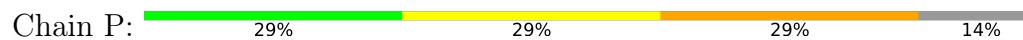
• Molecule 2: DNA (5'-D(\*C\*UP\*UP\*UP\*(BRU)P\*UP\*U)-3')



• Molecule 2: DNA (5'-D(\*C\*UP\*UP\*UP\*(BRU)P\*UP\*U)-3')



• Molecule 2: DNA (5'-D(\*C\*UP\*UP\*UP\*(BRU)P\*UP\*U)-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	37.58Å 137.08Å 83.96Å 90.00° 102.97° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.1 (50.00-2.50) 92.0 (50.00-2.50)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.88 (at 2.51Å)	Xtrriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.208 , 0.271 0.203 , 0.266	Depositor DCC
$R_{free}$ test set	5066 reflections (9.68%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.7	Xtrriage
Anisotropy	0.667	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 30.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	0.306 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	6102	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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: DIO, BRU, SO4

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	0/1370	0.89	2/1849 (0.1%)
1	B	0.43	0/1366	1.00	7/1843 (0.4%)
1	D	0.43	0/1370	0.94	2/1849 (0.1%)
1	I	0.45	0/1358	0.93	2/1832 (0.1%)
2	E	0.94	2/100 (2.0%)	0.63	0/148
2	H	1.05	1/104 (1.0%)	0.72	0/155
2	K	1.04	0/104	0.77	0/155
2	P	1.02	1/83 (1.2%)	0.84	0/123
All	All	0.50	4/5855 (0.1%)	0.93	13/7954 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	E	0	1
2	P	0	1
All	All	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	2	DU	O3'-P	5.35	1.61	1.56
2	H	2	DU	O3'-P	5.28	1.61	1.56
2	E	3	DU	O3'-P	5.12	1.61	1.56
2	P	3	DU	O3'-P	5.08	1.61	1.56

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	SER	N-CA-C	-10.28	100.74	113.18
1	B	286	LYS	N-CA-C	-6.58	105.55	113.97
1	I	191	ILE	N-CA-C	6.48	117.25	108.17
1	B	184	ASN	CA-C-N	5.98	125.52	119.19
1	B	184	ASN	C-N-CA	5.98	125.52	119.19

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	E	6	DU	Sidechain
2	P	6	DU	Sidechain

## 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	1343	0	1336	50	0
1	B	1339	0	1334	50	0
1	D	1343	0	1336	62	0
1	I	1332	0	1326	50	0
2	E	112	0	60	1	0
2	H	116	0	60	0	0
2	K	116	0	60	2	0
2	P	97	0	50	8	0
3	A	6	0	8	2	0
3	B	6	0	8	0	0
3	D	6	0	8	0	0
3	I	6	0	8	0	0
4	A	10	0	0	0	0
4	B	10	0	0	0	0
4	D	15	0	0	0	0
4	H	5	0	0	0	0
4	I	15	0	0	0	0
5	A	68	0	0	0	0
5	B	29	0	0	2	0
5	D	49	0	0	5	0
5	E	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	H	12	0	0	0	0
5	I	43	0	0	4	0
5	K	10	0	0	1	0
5	P	6	0	0	1	0
All	All	6102	0	5594	212	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 212 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:190:GLN:HE22	2:P:7:DU:H4'	1.35	0.91
1:A:143:GLY:HA2	1:A:146:GLY:H	1.36	0.88
1:B:144:PRO:HD2	1:B:149:ARG:HH22	1.38	0.86
1:I:145:LEU:HD22	1:I:147:SER:H	1.41	0.85
1:D:161:THR:OG1	1:D:163:GLU:HG3	1.77	0.84

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	172/174 (99%)	164 (95%)	8 (5%)	0	100	100
1	B	171/174 (98%)	158 (92%)	12 (7%)	1 (1%)	21	38
1	D	172/174 (99%)	168 (98%)	2 (1%)	2 (1%)	10	20
1	I	170/174 (98%)	152 (89%)	15 (9%)	3 (2%)	6	12
All	All	685/696 (98%)	642 (94%)	37 (5%)	6 (1%)	14	27

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	268	ASN
1	I	294	SER
1	I	236	PRO
1	D	267	PRO
1	I	310	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	141/141 (100%)	133 (94%)	8 (6%)	18	39
1	B	141/141 (100%)	134 (95%)	7 (5%)	22	44
1	D	141/141 (100%)	135 (96%)	6 (4%)	26	51
1	I	140/141 (99%)	130 (93%)	10 (7%)	13	29
All	All	563/564 (100%)	532 (94%)	31 (6%)	19	40

5 of 31 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	298	LEU
1	I	279	LEU
1	D	178	LEU
1	I	298	LEU
1	I	179	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	324	GLN
1	I	315	GLN
1	I	271	ASN
1	B	220	GLN
1	D	315	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](#)

4 non-standard protein/DNA/RNA residues 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	BRU	H	5	2	18,21,22	0.36	0	25,30,33	0.46	0
2	BRU	P	5	2	18,21,22	0.45	0	25,30,33	0.59	0
2	BRU	E	5	2	18,21,22	0.39	0	25,30,33	0.53	0
2	BRU	K	5	2	18,21,22	0.36	0	25,30,33	0.54	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
2	BRU	H	5	2	-	0/7/21/22	0/2/2/2
2	BRU	P	5	2	-	0/7/21/22	0/2/2/2
2	BRU	E	5	2	-	0/7/21/22	0/2/2/2
2	BRU	K	5	2	-	0/7/21/22	0/2/2/2

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.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

15 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
4	SO4	A	402	-	4,4,4	0.35	0	6,6,6	0.06	0
4	SO4	I	402	-	4,4,4	0.35	0	6,6,6	0.10	0
3	DIO	A	401	-	6,6,6	0.60	0	6,6,6	2.00	2 (33%)
4	SO4	I	404	-	4,4,4	0.35	0	6,6,6	0.10	0
3	DIO	I	401	-	6,6,6	0.65	0	6,6,6	1.99	2 (33%)
4	SO4	H	101	-	4,4,4	0.34	0	6,6,6	0.11	0
3	DIO	B	401	-	6,6,6	0.64	0	6,6,6	2.00	2 (33%)
4	SO4	A	403	-	4,4,4	0.36	0	6,6,6	0.10	0
4	SO4	D	404	-	4,4,4	0.36	0	6,6,6	0.08	0
4	SO4	D	403	-	4,4,4	0.38	0	6,6,6	0.04	0
4	SO4	I	403	-	4,4,4	0.38	0	6,6,6	0.08	0
3	DIO	D	401	-	6,6,6	0.61	0	6,6,6	2.00	2 (33%)
4	SO4	D	402	-	4,4,4	0.35	0	6,6,6	0.11	0
4	SO4	B	403	-	4,4,4	0.36	0	6,6,6	0.10	0
4	SO4	B	402	-	4,4,4	0.37	0	6,6,6	0.09	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	DIO	B	401	-	-	-	0/1/1/1
3	DIO	A	401	-	-	-	0/1/1/1
3	DIO	D	401	-	-	-	0/1/1/1
3	DIO	I	401	-	-	-	0/1/1/1

There are no bond length outliers.

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	DIO	C2-O1-C1	3.56	121.40	109.88
3	D	401	DIO	C2-O1-C1	3.53	121.30	109.88
3	I	401	DIO	C2-O1-C1	3.52	121.26	109.88
3	A	401	DIO	C2-O1-C1	3.50	121.19	109.88
3	A	401	DIO	C2'-O1'-C1'	3.39	120.84	109.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	DIO	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	174/174 (100%)	-1.54	0 100 100	21, 35, 54, 69	0
1	B	173/174 (99%)	-1.50	0 100 100	20, 40, 65, 84	0
1	D	174/174 (100%)	-1.54	0 100 100	20, 35, 57, 64	0
1	I	172/174 (98%)	-1.48	0 100 100	21, 40, 66, 82	0
2	E	0/7	-	-	-	-
2	H	1/7 (14%)	-2.15	0 100 100	66, 66, 66, 66	0
2	K	1/7 (14%)	-1.55	0 100 100	68, 68, 68, 68	0
2	P	0/7	-	-	-	-
All	All	695/724 (95%)	-1.52	0 100 100	20, 37, 62, 84	0

There are no RSRZ outliers to report.

### 6.2 Non-standard residues in protein, DNA, RNA chains [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
2	BRU	P	5	20/21	0.99	0.03	28,40,49,59	0
2	BRU	H	5	20/21	1.00	0.02	20,30,36,52	0
2	BRU	K	5	20/21	1.00	0.02	22,32,36,46	0
2	BRU	E	5	20/21	1.00	0.03	31,35,43,60	0

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides 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	DIO	B	401	6/6	0.98	0.08	64,66,67,67	0
3	DIO	I	401	6/6	0.98	0.08	65,66,67,68	0
4	SO4	A	403	5/5	0.98	0.04	99,100,101,101	0
4	SO4	B	402	5/5	0.98	0.04	92,92,93,93	0
4	SO4	D	403	5/5	0.98	0.05	87,87,88,89	0
4	SO4	D	404	5/5	0.98	0.04	105,106,106,107	0
4	SO4	I	403	5/5	0.98	0.06	93,93,94,94	0
4	SO4	I	404	5/5	0.98	0.07	122,122,123,123	0
4	SO4	D	402	5/5	0.99	0.04	68,69,70,71	0
4	SO4	A	402	5/5	0.99	0.03	71,71,72,72	0
3	DIO	D	401	6/6	0.99	0.07	95,96,96,96	0
4	SO4	I	402	5/5	0.99	0.04	70,72,74,74	0
3	DIO	A	401	6/6	0.99	0.06	84,85,85,85	0
4	SO4	B	403	5/5	0.99	0.04	72,73,75,75	0
4	SO4	H	101	5/5	0.99	0.04	104,104,105,106	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.