



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:52 PM GMT

PDB ID : 3VAL
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 Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at 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.

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

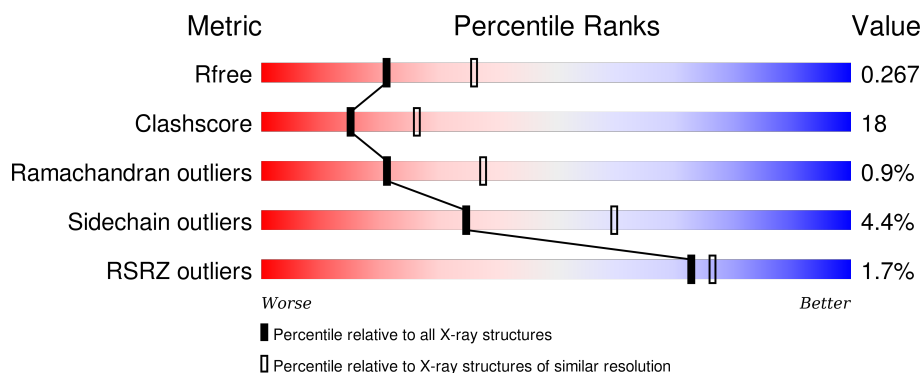
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	174	<div> <div>2%</div> <div>72%</div> <div>26%</div> <div>.</div> </div>
1	B	174	<div> <div>%</div> <div>70%</div> <div>29%</div> <div>..</div> </div>
1	D	174	<div> <div>%</div> <div>62%</div> <div>36%</div> <div>.</div> </div>
1	I	174	<div> <div>3%</div> <div>62%</div> <div>34%</div> <div>...</div> </div>
2	E	7	<div> <div>43%</div> <div>43%</div> <div>14%</div> </div>

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Mol	Chain	Length	Quality of chain
2	H	7	 86%14%
2	K	7	 57%43%
2	P	7	 14%57%14%14%

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	DIO	A	401	-	-	-	X
3	DIO	B	401	-	-	-	X
3	DIO	D	401	-	-	-	X
4	SO4	D	404	-	-	-	X

2 Entry composition

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
1	A	174	Total	C	N	O	S	0	0	0
			1343	857	230	250	6			
1	B	173	Total	C	N	O	S	0	0	0
			1339	855	229	249	6			
1	D	174	Total	C	N	O	S	0	0	0
			1343	857	230	250	6			
1	I	172	Total	C	N	O	S	0	0	0
			1332	850	228	248	6			

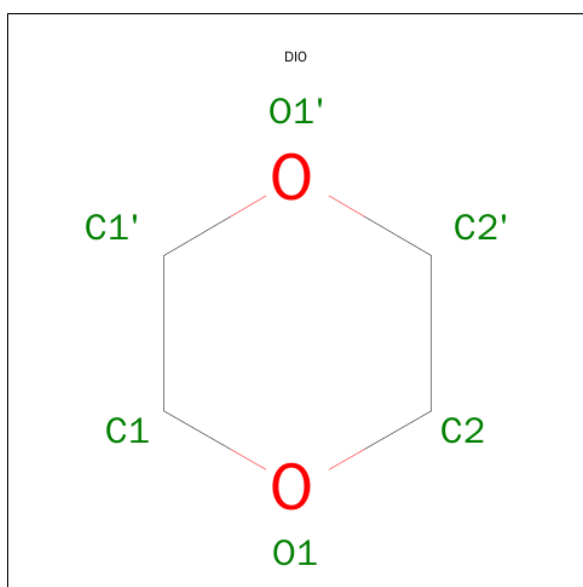
There are 20 discrepancies between the modelled and reference sequences:

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
2	E	6	Total	Br	C	N	O	P	0	0	0
			112	1	54	12	40	5			
2	H	7	Total	Br	C	N	O	P	0	0	1
			116	1	54	12	43	6			
2	K	7	Total	Br	C	N	O	P	0	0	1
			116	1	54	12	43	6			
2	P	6	Total	Br	C	N	O	P	0	0	1
			97	1	45	10	36	5			

- Molecule 3 is 1,4-DIETHYLENE DIOXIDE (three-letter code: DIO) (formula: C₄H₈O₂).



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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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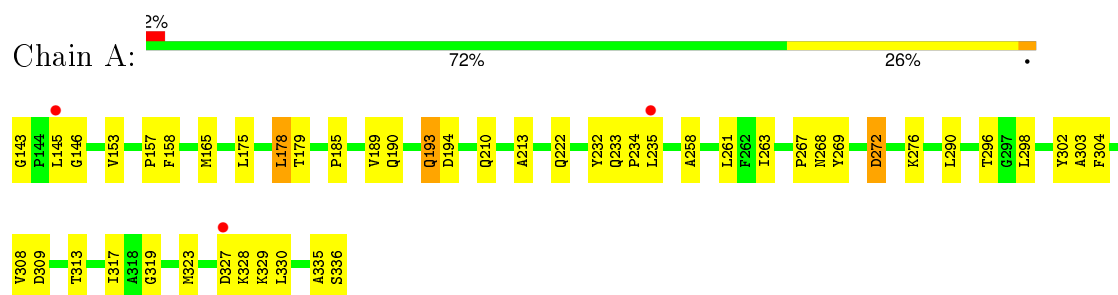
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	29	Total 29	O 29	0	0
5	D	49	Total 49	O 49	0	0
5	I	43	Total 43	O 43	0	0
5	E	8	Total 8	O 8	0	0
5	H	12	Total 12	O 12	0	0
5	K	10	Total 10	O 10	0	0
5	P	6	Total 6	O 6	0	0

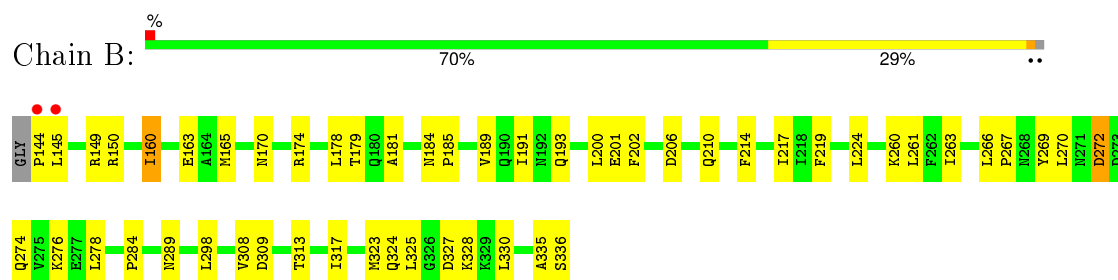
3 Residue-property plots [i](#)

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.

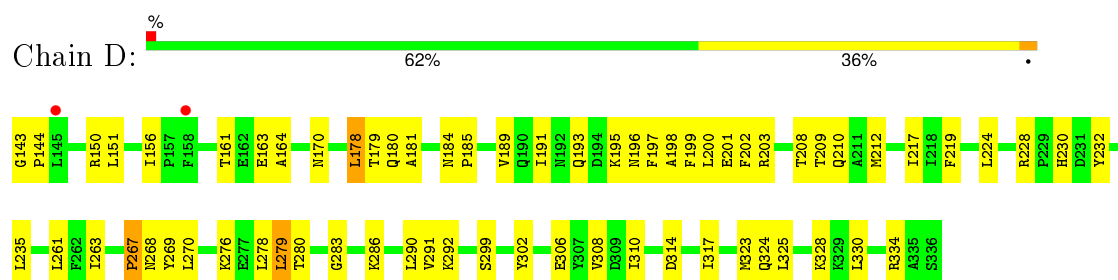
- Molecule 1: Splicing factor U2AF 65 kDa subunit



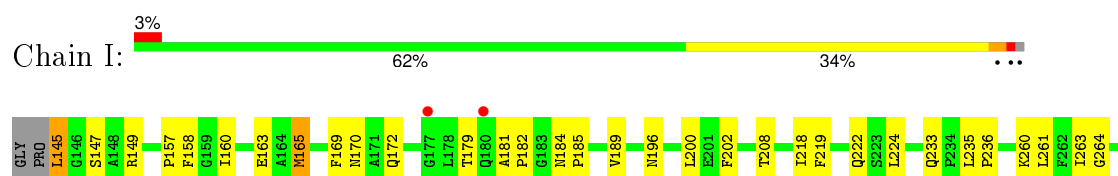
- Molecule 1: Splicing factor U2AF 65 kDa subunit

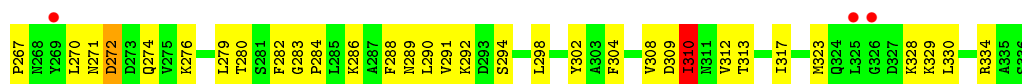


- Molecule 1: Splicing factor U2AF 65 kDa subunit

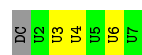


- Molecule 1: Splicing factor U2AF 65 kDa subunit

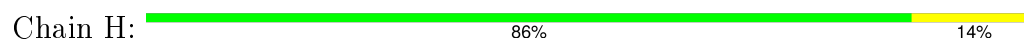




- Molecule 2: DNA (5'-D(*C*UP*UP*UP*(BRU)P*UP*U)-3')



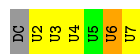
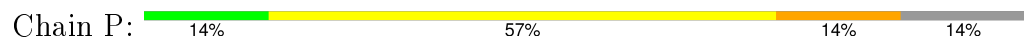
- Molecule 2: DNA (5'-D(*C*UP*UP*UP*(BRU)P*UP*U)-3')



- 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, α , β , γ	37.58Å 137.08Å 83.96Å 90.00° 102.97° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 28.58 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.1 (50.00-2.50) 93.4 (28.58-2.50)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.51Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.208 , 0.271 0.206 , 0.267	Depositor DCC
R_{free} test set	2602 reflections (10.93%)	DCC
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.667	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 26.1	EDS
Estimated twinning fraction	0.306 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 52319 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6102	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage'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.*

¹Intensities estimated from amplitudes.

²Theoretical 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: BRU, DIO, 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	0.39	0/1370	0.61	0/1849
1	B	0.38	0/1366	0.63	0/1843
1	D	0.38	0/1370	0.60	0/1849
1	I	0.39	0/1358	0.61	0/1832
2	E	0.31	0/100	0.63	0/148
2	H	0.91	1/104 (1.0%)	0.72	0/155
2	K	0.92	1/104 (1.0%)	0.77	0/155
2	P	0.99	1/83 (1.2%)	0.84	0/123
All	All	0.43	3/5855 (0.1%)	0.62	0/7954

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 (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	2	DU	O3'-P	-7.13	1.52	1.61
2	K	1	DC	O3'-P	-7.08	1.52	1.61
2	H	1	DC	O3'-P	-7.06	1.52	1.61

There are no bond angle outliers.

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 ⓘ

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	46	0
1	B	1339	0	1334	49	0
1	D	1343	0	1336	59	0
1	I	1332	0	1326	46	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
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	201	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.

All (201) 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
1:A:193:GLN:NE2	1:A:193:GLN:H	1.77	0.83
1:D:219:PHE:HB3	1:D:224:LEU:HD11	1.62	0.82
1:A:235:LEU:H	1:A:235:LEU:HD23	1.46	0.79
1:A:190:GLN:HE22	2:P:7:DU:C4'	1.96	0.79
1:A:308:VAL:HG23	1:A:309:ASP:H	1.47	0.78
1:A:193:GLN:HG2	1:A:194:ASP:H	1.49	0.76
1:D:228:ARG:HD3	1:D:232:TYR:CE2	2.23	0.74
1:B:335:ALA:O	1:B:336:SER:HB3	1.88	0.73
1:A:272:ASP:O	1:A:276:LYS:HB2	1.90	0.72
1:B:263:ILE:HD12	1:B:263:ILE:N	2.05	0.71
1:D:235:LEU:H	1:D:235:LEU:HD12	1.53	0.71
1:B:149:ARG:HG2	1:B:202:PHE:O	1.92	0.70
1:B:284:PRO:HG2	1:B:308:VAL:HG11	1.72	0.70
1:A:261:LEU:HD21	1:A:313:THR:HG23	1.74	0.70
1:D:219:PHE:CB	1:D:224:LEU:HD11	2.21	0.70
1:A:234:PRO:HA	3:A:401:DIO:H1'2	1.74	0.69
1:D:193:GLN:H	1:D:193:GLN:CD	1.96	0.69
1:A:308:VAL:HG23	1:A:309:ASP:N	2.07	0.68
1:A:193:GLN:HG2	1:A:194:ASP:N	2.10	0.66
1:I:145:LEU:HG	1:I:286:LYS:HE3	1.78	0.66
1:I:284:PRO:HG2	1:I:308:VAL:CG1	2.26	0.66
2:P:6:DU:H2'	2:P:7:DU:OP2	1.96	0.66
1:B:174:ARG:HG2	1:B:174:ARG:HH11	1.60	0.66
1:D:163:GLU:HG3	1:D:164:ALA:H	1.60	0.66
1:A:235:LEU:N	1:A:235:LEU:HD23	2.12	0.65
1:B:267:PRO:HG3	1:B:328:LYS:HD3	1.79	0.64
1:B:308:VAL:HG13	1:B:309:ASP:N	2.12	0.64
1:A:296:THR:HB	1:A:298:LEU:HG	1.77	0.64
1:A:235:LEU:HG	1:A:258:ALA:HB2	1.79	0.64
1:I:267:PRO:HD2	1:I:270:LEU:HD12	1.80	0.64
1:A:267:PRO:HA	1:A:328:LYS:NZ	2.13	0.64
1:B:266:LEU:HD23	1:B:330:LEU:HD11	1.81	0.63
1:I:284:PRO:HG2	1:I:308:VAL:HG11	1.81	0.63
1:I:270:LEU:HD23	1:I:274:GLN:OE1	1.98	0.63
1:D:230:HIS:HB3	5:D:538:HOH:O	1.97	0.63
1:I:290:LEU:HB3	1:I:292:LYS:NZ	2.13	0.62
1:B:145:LEU:N	1:B:145:LEU:HD12	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:217:ILE:O	1:D:224:LEU:HD13	2.00	0.61
1:A:190:GLN:NE2	2:P:7:DU:H4'	2.12	0.60
1:I:145:LEU:HD13	1:I:147:SER:OG	2.01	0.60
1:D:286:LYS:HB3	1:D:308:VAL:HG22	1.82	0.60
1:A:335:ALA:O	1:A:336:SER:HB3	2.02	0.59
1:B:179:THR:HG21	5:B:508:HOH:O	2.02	0.59
1:B:308:VAL:HG13	1:B:309:ASP:H	1.67	0.59
1:A:296:THR:CB	1:A:298:LEU:HG	2.32	0.59
1:A:235:LEU:H	1:A:235:LEU:CD2	2.15	0.58
1:D:228:ARG:HD3	1:D:232:TYR:CD2	2.37	0.58
1:A:235:LEU:O	1:A:258:ALA:HB2	2.04	0.58
1:A:179:THR:HG23	1:A:185:PRO:HG3	1.84	0.58
1:B:284:PRO:HG2	1:B:308:VAL:CG1	2.33	0.57
1:I:290:LEU:HB3	1:I:292:LYS:HZ2	1.67	0.57
1:D:328:LYS:HE2	2:K:2:DU:O4	2.04	0.56
1:D:143:GLY:HA3	1:D:144:PRO:C	2.25	0.56
1:A:175:LEU:HD23	1:I:172:GLN:NE2	2.21	0.56
1:D:278:LEU:CD2	1:D:325:LEU:HD11	2.36	0.55
1:D:224:LEU:N	1:D:224:LEU:HD12	2.20	0.55
1:D:267:PRO:O	1:D:269:TYR:N	2.41	0.54
1:I:264:GLY:O	1:I:330:LEU:HD12	2.07	0.54
1:D:163:GLU:HG3	1:D:164:ALA:N	2.22	0.54
1:A:165:MET:HG2	1:A:189:VAL:HG11	1.90	0.53
1:I:272:ASP:O	1:I:276:LYS:HB2	2.08	0.53
1:A:313:THR:O	1:A:317:ILE:HG13	2.09	0.53
1:I:271:ASN:ND2	1:I:274:GLN:HE21	2.07	0.53
1:I:271:ASN:ND2	1:I:274:GLN:NE2	2.57	0.52
1:I:196:ASN:O	1:I:196:ASN:CG	2.48	0.52
1:B:313:THR:O	1:B:317:ILE:HG13	2.09	0.52
1:D:179:THR:HG23	1:D:185:PRO:HG3	1.90	0.52
1:B:170:ASN:O	1:B:174:ARG:HG3	2.10	0.52
1:A:267:PRO:HA	1:A:328:LYS:HZ1	1.74	0.52
1:B:189:VAL:HG22	1:B:200:LEU:HD23	1.90	0.52
1:A:157:PRO:HA	1:A:222:GLN:HG3	1.93	0.51
1:D:278:LEU:HD22	1:D:325:LEU:HD11	1.93	0.51
1:I:282:PHE:O	1:I:312:VAL:HG11	2.10	0.51
1:D:191:ILE:HG12	1:D:198:ALA:HB2	1.93	0.51
1:I:157:PRO:HG3	1:I:222:GLN:HB2	1.92	0.50
1:B:174:ARG:HG2	1:B:174:ARG:NH1	2.26	0.50
1:B:160:ILE:HG23	1:B:191:ILE:HD13	1.94	0.50
1:I:260:LYS:HZ3	1:I:304:PHE:HB3	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:ASP:O	1:B:210:GLN:HG2	2.11	0.50
2:P:6:DU:H5	5:P:104:HOH:O	2.11	0.50
1:D:267:PRO:HB3	1:D:269:TYR:CE2	2.47	0.50
1:A:153:VAL:HG11	1:A:165:MET:HE1	1.92	0.50
1:D:209:THR:O	1:D:212:MET:HB2	2.12	0.49
1:B:144:PRO:HD2	1:B:149:ARG:NH2	2.17	0.49
1:D:286:LYS:HE2	1:D:306:GLU:HG2	1.94	0.49
1:B:150:ARG:HA	1:B:200:LEU:O	2.12	0.49
2:E:3:DU:H4'	2:E:4:DU:H5'	1.95	0.49
1:A:308:VAL:CG2	1:A:309:ASP:H	2.21	0.49
1:D:261:LEU:HD21	1:D:317:ILE:HD11	1.94	0.49
1:D:292:LYS:HG3	1:D:299:SER:HA	1.94	0.49
1:I:291:VAL:HG21	1:I:302:TYR:CE1	2.47	0.49
1:I:169:PHE:CE2	1:I:200:LEU:HD11	2.48	0.49
1:D:235:LEU:N	1:D:235:LEU:HD12	2.26	0.48
1:D:261:LEU:CD2	1:D:317:ILE:HD11	2.44	0.48
1:B:284:PRO:O	1:B:308:VAL:HG12	2.14	0.48
1:A:290:LEU:HD12	1:A:303:ALA:HB2	1.96	0.48
1:A:145:LEU:HD23	3:A:401:DIO:H22	1.94	0.47
1:D:151:LEU:O	1:D:199:PHE:HA	2.14	0.47
1:I:170:ASN:CG	1:I:184:ASN:HB3	2.34	0.47
1:A:308:VAL:CG2	1:A:309:ASP:N	2.77	0.47
1:B:165:MET:HE3	1:B:200:LEU:HD21	1.97	0.47
1:D:161:THR:HG23	5:D:516:HOH:O	2.14	0.47
1:B:289:ASN:ND2	5:B:511:HOH:O	2.37	0.47
1:B:325:LEU:HD12	1:B:330:LEU:HD21	1.95	0.47
1:B:179:THR:CG2	1:B:181:ALA:H	2.28	0.47
1:D:291:VAL:HG21	1:D:302:TYR:CE1	2.49	0.47
1:D:325:LEU:CD1	1:D:330:LEU:HD11	2.45	0.46
1:B:144:PRO:HB2	1:B:145:LEU:HD12	1.96	0.46
1:I:267:PRO:HD2	1:I:270:LEU:CD1	2.46	0.46
1:I:284:PRO:HG2	1:I:308:VAL:HG12	1.96	0.46
1:B:327:ASP:OD1	1:B:328:LYS:HD2	2.16	0.46
1:B:214:PHE:O	1:B:217:ILE:HB	2.16	0.46
1:I:149:ARG:NH1	5:I:527:HOH:O	2.47	0.46
1:D:280:THR:HA	1:D:283:GLY:O	2.16	0.46
1:B:267:PRO:HD2	1:B:270:LEU:HD12	1.98	0.46
1:B:179:THR:HG23	1:B:181:ALA:H	1.79	0.46
1:B:270:LEU:HA	1:B:274:GLN:OE1	2.15	0.45
2:P:6:DU:C2'	2:P:7:DU:OP2	2.62	0.45
1:B:178:LEU:HD22	1:B:214:PHE:CZ	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:224:LEU:CD1	1:D:224:LEU:N	2.80	0.45
1:B:308:VAL:CG1	1:B:309:ASP:H	2.30	0.45
1:I:323:MET:O	1:I:329:LYS:HA	2.16	0.45
1:B:184:ASN:HA	1:B:185:PRO:HD3	1.80	0.45
1:B:267:PRO:HB3	1:B:269:TYR:CE2	2.52	0.45
1:A:263:ILE:O	1:A:302:TYR:HA	2.17	0.45
1:B:219:PHE:HB3	1:B:224:LEU:HD11	1.98	0.45
1:I:271:ASN:HD21	1:I:274:GLN:HE21	1.63	0.45
1:I:149:ARG:HD3	5:I:527:HOH:O	2.17	0.45
1:D:178:LEU:HG	1:D:210:GLN:HB3	1.98	0.45
1:B:323:MET:HE3	1:B:325:LEU:HD21	1.98	0.44
1:I:165:MET:HG2	1:I:189:VAL:HG11	1.99	0.44
1:I:170:ASN:OD1	1:I:184:ASN:HB3	2.17	0.44
1:D:276:LYS:HE2	5:D:543:HOH:O	2.17	0.44
1:A:232:TYR:O	1:A:233:GLN:HG3	2.18	0.44
1:B:308:VAL:CG1	1:B:309:ASP:N	2.79	0.44
1:D:302:TYR:HB3	5:K:104:HOH:O	2.18	0.44
1:D:184:ASN:O	1:D:203:ARG:HD3	2.17	0.44
1:B:170:ASN:CG	1:B:184:ASN:HB3	2.38	0.44
1:D:179:THR:HG22	1:D:180:GLN:N	2.33	0.44
1:D:189:VAL:HG22	1:D:200:LEU:CD2	2.47	0.44
1:A:153:VAL:HG11	1:A:165:MET:CE	2.48	0.44
1:B:261:LEU:HD21	1:B:313:THR:HG23	2.00	0.44
1:I:160:ILE:HG12	1:I:219:PHE:HE2	1.83	0.44
1:B:163:GLU:N	1:B:163:GLU:OE1	2.43	0.44
1:D:170:ASN:OD1	1:D:185:PRO:HD2	2.18	0.43
1:D:202:PHE:HB2	1:D:208:THR:OG1	2.18	0.43
1:D:189:VAL:HG22	1:D:200:LEU:HD23	2.00	0.43
1:A:193:GLN:HE21	1:A:193:GLN:H	1.60	0.43
1:B:189:VAL:HG22	1:B:200:LEU:CD2	2.49	0.43
1:I:261:LEU:HD21	1:I:317:ILE:HD11	1.98	0.43
1:I:235:LEU:HB2	5:I:512:HOH:O	2.17	0.43
1:I:145:LEU:HD12	1:I:233:GLN:OE1	2.18	0.43
1:A:261:LEU:O	1:A:304:PHE:HA	2.18	0.43
1:I:308:VAL:HG13	1:I:309:ASP:N	2.33	0.43
1:D:179:THR:HG22	1:D:181:ALA:H	1.83	0.43
1:B:323:MET:HG2	1:B:324:GLN:N	2.34	0.43
1:I:328:LYS:HG2	1:I:329:LYS:N	2.33	0.43
1:A:319:GLY:O	1:A:323:MET:HG3	2.19	0.43
2:P:3:DU:H4'	2:P:4:DU:H5'	2.01	0.42
1:I:181:ALA:HB1	1:I:182:PRO:CD	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:310:ILE:HG12	1:I:310:ILE:O	2.18	0.42
1:A:193:GLN:CG	1:A:194:ASP:H	2.24	0.42
1:B:217:ILE:HA	1:B:217:ILE:HD12	1.90	0.42
1:I:263:ILE:HD12	1:I:263:ILE:N	2.34	0.42
1:D:150:ARG:NH1	1:D:201:GLU:HG2	2.34	0.42
1:D:267:PRO:C	1:D:269:TYR:H	2.23	0.42
1:A:267:PRO:HB3	1:A:269:TYR:CZ	2.54	0.42
1:D:325:LEU:HD12	1:D:330:LEU:HD11	2.00	0.42
1:A:329:LYS:HZ1	1:D:323:MET:HE2	1.85	0.42
1:D:270:LEU:HA	1:D:270:LEU:HD23	1.90	0.42
1:A:158:PHE:CZ	1:B:260:LYS:HD3	2.55	0.42
1:B:272:ASP:O	1:B:276:LYS:HB2	2.19	0.42
1:I:149:ARG:HA	1:I:208:THR:OG1	2.20	0.42
1:D:276:LYS:CE	5:D:543:HOH:O	2.68	0.42
1:B:323:MET:HG2	1:B:324:GLN:H	1.85	0.42
1:I:261:LEU:HD23	1:I:334:ARG:HA	2.02	0.42
1:I:185:PRO:O	1:I:202:PHE:HA	2.20	0.42
1:D:314:ASP:CG	1:D:334:ARG:HH22	2.23	0.41
1:D:228:ARG:NH1	5:D:507:HOH:O	2.53	0.41
1:D:178:LEU:HA	1:D:178:LEU:HD12	1.90	0.41
1:A:296:THR:OG1	1:A:298:LEU:HG	2.20	0.41
1:D:263:ILE:CD1	1:D:279:LEU:HD11	2.49	0.41
1:I:222:GLN:HG3	5:I:518:HOH:O	2.20	0.41
1:I:313:THR:O	1:I:317:ILE:HG13	2.21	0.41
1:D:156:ILE:HG12	1:D:196:ASN:O	2.20	0.41
1:A:143:GLY:HA2	1:A:145:LEU:N	2.36	0.41
1:D:195:LYS:HB2	1:D:197:PHE:CE2	2.56	0.41
1:D:302:TYR:OH	2:K:4:DU:H4'	2.21	0.41
1:I:288:PHE:CG	1:I:289:ASN:N	2.88	0.41
1:D:161:THR:OG1	1:D:163:GLU:CG	2.59	0.40
1:D:290:LEU:O	1:D:292:LYS:HD3	2.20	0.40
1:B:260:LYS:HE2	2:P:4:DU:O4	2.21	0.40
1:D:219:PHE:HB2	1:D:224:LEU:HD11	2.02	0.40
1:A:178:LEU:HG	1:A:210:GLN:HB3	2.03	0.40
1:I:280:THR:HA	1:I:283:GLY:O	2.22	0.40
1:A:213:ALA:O	1:I:218:ILE:HD11	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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%)	30	50
1	D	172/174 (99%)	168 (98%)	2 (1%)	2 (1%)	16	29
1	I	170/174 (98%)	152 (89%)	15 (9%)	3 (2%)	11	18
All	All	685/696 (98%)	642 (94%)	37 (5%)	6 (1%)	21	37

All (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
1	B	160	ILE

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	141/141 (100%)	135 (96%)	6 (4%)	35	61
1	B	141/141 (100%)	136 (96%)	5 (4%)	43	70
1	D	141/141 (100%)	137 (97%)	4 (3%)	51	78
1	I	140/141 (99%)	130 (93%)	10 (7%)	18	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	563/564 (100%)	538 (96%)	25 (4%)	35 60

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	178	LEU
1	A	193	GLN
1	A	268	ASN
1	A	272	ASP
1	A	327	ASP
1	A	330	LEU
1	B	193	GLN
1	B	201	GLU
1	B	272	ASP
1	B	278	LEU
1	B	298	LEU
1	D	178	LEU
1	D	279	LEU
1	D	310	ILE
1	D	324	GLN
1	I	145	LEU
1	I	158	PHE
1	I	163	GLU
1	I	165	MET
1	I	179	THR
1	I	224	LEU
1	I	272	ASP
1	I	279	LEU
1	I	298	LEU
1	I	310	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	GLN
1	A	193	GLN
1	A	220	GLN
1	A	315	GLN
1	B	190	GLN
1	B	196	ASN
1	B	220	GLN

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Mol	Chain	Res	Type
1	B	222	GLN
1	B	233	GLN
1	B	289	ASN
1	D	220	GLN
1	D	315	GLN
1	D	324	GLN
1	I	220	GLN
1	I	271	ASN
1	I	315	GLN
1	I	324	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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	E	5	2	13,21,22	1.50	2 (15%)	16,30,33	4.11	2 (12%)
2	BRU	H	5	2	13,21,22	1.49	2 (15%)	16,30,33	4.05	2 (12%)
2	BRU	K	5	2	13,21,22	1.61	2 (15%)	16,30,33	4.05	2 (12%)
2	BRU	P	5	2	13,21,22	1.61	3 (23%)	16,30,33	4.09	2 (12%)

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	E	5	2	-	0/3/21/22	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BRU	H	5	2	-	0/3/21/22	0/2/2/2
2	BRU	K	5	2	-	0/3/21/22	0/2/2/2
2	BRU	P	5	2	-	0/3/21/22	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	5	BRU	C6-N1	2.09	1.38	1.35
2	H	5	BRU	C4-N3	2.58	1.37	1.33
2	E	5	BRU	C4-N3	2.62	1.38	1.33
2	K	5	BRU	C4-N3	2.76	1.38	1.33
2	P	5	BRU	C4-N3	2.78	1.38	1.33
2	H	5	BRU	C4-C5	3.91	1.43	1.38
2	E	5	BRU	C4-C5	4.09	1.43	1.38
2	P	5	BRU	C4-C5	4.27	1.44	1.38
2	K	5	BRU	C4-C5	4.45	1.44	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	5	BRU	C5-C4-N3	-8.91	114.48	124.00
2	K	5	BRU	C5-C4-N3	-8.87	114.53	124.00
2	H	5	BRU	C5-C4-N3	-8.78	114.63	124.00
2	P	5	BRU	C5-C4-N3	-8.74	114.67	124.00
2	K	5	BRU	C4-N3-C2	13.44	126.87	115.25
2	H	5	BRU	C4-N3-C2	13.47	126.89	115.25
2	P	5	BRU	C4-N3-C2	13.70	127.09	115.25
2	E	5	BRU	C4-N3-C2	13.70	127.09	115.25

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

There are no carbohydrates in this entry.

5.6 Ligand geometry

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
3	DIO	A	401	-	6,6,6	0.58	0	6,6,6	1.91	2 (33%)
4	SO4	A	402	-	4,4,4	0.21	0	6,6,6	0.13	0
4	SO4	A	403	-	4,4,4	0.18	0	6,6,6	0.08	0
3	DIO	B	401	-	6,6,6	0.62	0	6,6,6	1.92	2 (33%)
4	SO4	B	402	-	4,4,4	0.22	0	6,6,6	0.08	0
4	SO4	B	403	-	4,4,4	0.24	0	6,6,6	0.14	0
3	DIO	D	401	-	6,6,6	0.59	0	6,6,6	1.92	2 (33%)
4	SO4	D	402	-	4,4,4	0.18	0	6,6,6	0.09	0
4	SO4	D	403	-	4,4,4	0.22	0	6,6,6	0.06	0
4	SO4	D	404	-	4,4,4	0.19	0	6,6,6	0.08	0
4	SO4	H	101	-	4,4,4	0.19	0	6,6,6	0.10	0
3	DIO	I	401	-	6,6,6	0.63	0	6,6,6	1.91	2 (33%)
4	SO4	I	402	-	4,4,4	0.21	0	6,6,6	0.15	0
4	SO4	I	403	-	4,4,4	0.25	0	6,6,6	0.10	0
4	SO4	I	404	-	4,4,4	0.21	0	6,6,6	0.10	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	A	401	-	-	0/0/6/6	0/1/1/1
4	SO4	A	402	-	-	0/0/0/0	0/0/0/0
4	SO4	A	403	-	-	0/0/0/0	0/0/0/0
3	DIO	B	401	-	-	0/0/6/6	0/1/1/1
4	SO4	B	402	-	-	0/0/0/0	0/0/0/0
4	SO4	B	403	-	-	0/0/0/0	0/0/0/0
3	DIO	D	401	-	-	0/0/6/6	0/1/1/1
4	SO4	D	402	-	-	0/0/0/0	0/0/0/0
4	SO4	D	403	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SO4	D	404	-	-	0/0/0/0	0/0/0/0
4	SO4	H	101	-	-	0/0/0/0	0/0/0/0
3	DIO	I	401	-	-	0/0/6/6	0/1/1/1
4	SO4	I	402	-	-	0/0/0/0	0/0/0/0
4	SO4	I	403	-	-	0/0/0/0	0/0/0/0
4	SO4	I	404	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	DIO	C2'-O1'-C1'	3.21	120.71	109.89
3	I	401	DIO	C2'-O1'-C1'	3.23	120.76	109.89
3	D	401	DIO	C2'-O1'-C1'	3.25	120.83	109.89
3	A	401	DIO	C2'-O1'-C1'	3.25	120.84	109.89
3	A	401	DIO	C2-O1-C1	3.35	121.19	109.89
3	I	401	DIO	C2-O1-C1	3.38	121.26	109.89
3	D	401	DIO	C2-O1-C1	3.39	121.30	109.89
3	B	401	DIO	C2-O1-C1	3.42	121.40	109.89

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

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, 95th 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 ²)	Q<0.9
1	A	174/174 (100%)	0.00	3 (1%) 73 76	21, 35, 54, 69	0
1	B	173/174 (99%)	0.11	2 (1%) 81 83	20, 40, 65, 84	0
1	D	174/174 (100%)	-0.07	2 (1%) 82 84	20, 35, 57, 64	0
1	I	172/174 (98%)	0.22	5 (2%) 55 60	21, 40, 66, 82	0
2	E	0/7	-	-	-	-
2	H	1/7 (14%)	-1.41	0 100 100	67, 67, 67, 67	0
2	K	1/7 (14%)	-0.31	0 100 100	68, 68, 68, 68	0
2	P	0/7	-	-	-	-
All	All	695/724 (95%)	0.06	12 (1%) 73 76	20, 37, 62, 84	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	326	GLY	4.5
1	B	144	PRO	3.5
1	I	325	LEU	2.9
1	D	145	LEU	2.7
1	A	145	LEU	2.6
1	B	145	LEU	2.6
1	A	235	LEU	2.5
1	I	180	GLN	2.5
1	I	177	GLY	2.4
1	I	269	TYR	2.4
1	D	158	PHE	2.3
1	A	327	ASP	2.3

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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, 95th 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(Å ²)	Q<0.9
2	BRU	H	5	20/21	0.98	0.11	-	20,30,36,52	0
2	BRU	E	5	20/21	0.97	0.12	-	31,35,43,60	0
2	BRU	K	5	20/21	0.98	0.11	-	22,32,36,46	0
2	BRU	P	5	20/21	0.94	0.12	-	28,40,49,59	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 95th 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(Å ²)	Q<0.9
3	DIO	A	401	6/6	0.82	0.62	6.70	84,85,85,85	0
3	DIO	D	401	6/6	0.86	0.59	5.48	95,96,96,96	0
4	SO4	D	404	5/5	0.87	0.23	2.64	105,106,106,107	0
3	DIO	B	401	6/6	0.82	0.28	2.03	64,66,67,67	0
3	DIO	I	401	6/6	0.95	0.17	-0.04	65,66,67,68	0
4	SO4	A	403	5/5	0.87	0.18	-	99,100,101,101	0
4	SO4	B	402	5/5	0.85	0.17	-	92,92,93,93	0
4	SO4	I	404	5/5	0.79	0.38	-	122,122,123,123	0
4	SO4	D	402	5/5	0.90	0.18	-	68,69,70,71	0
4	SO4	I	403	5/5	0.91	0.20	-	93,93,94,94	0
4	SO4	A	402	5/5	0.95	0.12	-	71,71,72,72	0
4	SO4	D	403	5/5	0.95	0.17	-	87,87,88,89	0
4	SO4	H	101	5/5	0.88	0.20	-	104,104,105,106	0
4	SO4	I	402	5/5	0.91	0.19	-	70,72,74,74	0
4	SO4	B	403	5/5	0.93	0.13	-	72,73,75,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.