



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:01 PM GMT

PDB ID : 4B8Z  
Title : Crystal structure of human GDP-L-fucose synthase with bound NADP and GDP, rhombohedral crystal form  
Authors : Vollmar, M.; Krojer, T.; Shafqat, N.; Rojkova, A.; Bunkoczi, G.; Yue, W.W.; Kavanagh, K.; von Delft, F.; Weigelt, J.; Arrowsmith, C.H.; Bountra, C.; Edwards, A.; Oppermann, U.  
Deposited on : 2012-08-31  
Resolution : 2.75 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

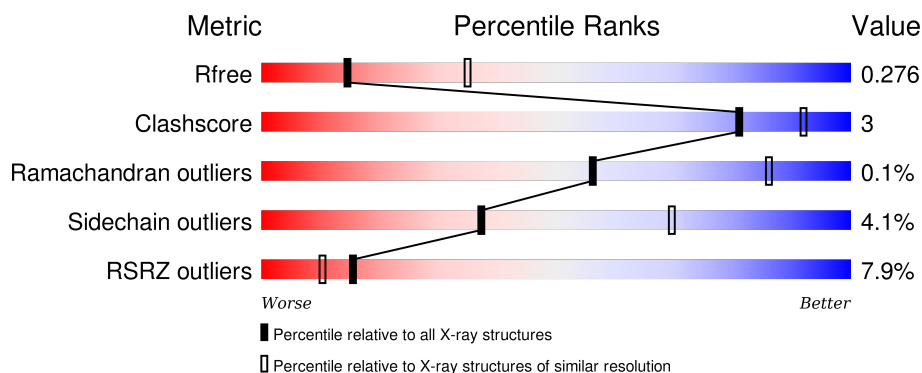
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.75 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	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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	320	<div> <div>11%</div> <div>89%</div> <div>7%</div> <div>.</div> </div>
1	B	320	<div> <div>4%</div> <div>88%</div> <div>11%</div> <div>.</div> </div>
1	C	320	<div> <div>8%</div> <div>88%</div> <div>10%</div> <div>.</div> </div>
1	D	320	<div> <div>8%</div> <div>88%</div> <div>9%</div> <div>.</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9492 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 GDP-L-FUCOSE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	308	Total	C	N	O	S	0	0	0
			2176	1379	379	409	9			
1	B	320	Total	C	N	O	S	0	0	0
			2427	1557	413	447	10			
1	C	313	Total	C	N	O	S	0	0	0
			2301	1472	392	428	9			
1	D	313	Total	C	N	O	S	0	0	0
			2266	1446	388	423	9			

There are 28 discrepancies between the modelled and reference sequences:

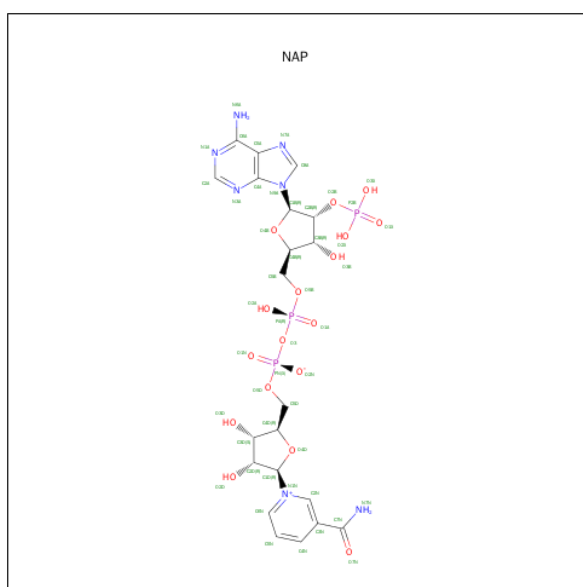
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLU	-	EXPRESSION TAG	UNP Q13630
A	-6	ASN	-	EXPRESSION TAG	UNP Q13630
A	-5	LEU	-	EXPRESSION TAG	UNP Q13630
A	-4	TYR	-	EXPRESSION TAG	UNP Q13630
A	-3	PHE	-	EXPRESSION TAG	UNP Q13630
A	-2	GLN	-	EXPRESSION TAG	UNP Q13630
A	-1	SER	-	EXPRESSION TAG	UNP Q13630
B	-7	GLU	-	EXPRESSION TAG	UNP Q13630
B	-6	ASN	-	EXPRESSION TAG	UNP Q13630
B	-5	LEU	-	EXPRESSION TAG	UNP Q13630
B	-4	TYR	-	EXPRESSION TAG	UNP Q13630
B	-3	PHE	-	EXPRESSION TAG	UNP Q13630
B	-2	GLN	-	EXPRESSION TAG	UNP Q13630
B	-1	SER	-	EXPRESSION TAG	UNP Q13630
C	-7	GLU	-	EXPRESSION TAG	UNP Q13630
C	-6	ASN	-	EXPRESSION TAG	UNP Q13630
C	-5	LEU	-	EXPRESSION TAG	UNP Q13630
C	-4	TYR	-	EXPRESSION TAG	UNP Q13630
C	-3	PHE	-	EXPRESSION TAG	UNP Q13630
C	-2	GLN	-	EXPRESSION TAG	UNP Q13630
C	-1	SER	-	EXPRESSION TAG	UNP Q13630

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-7	GLU	-	EXPRESSION TAG	UNP Q13630
D	-6	ASN	-	EXPRESSION TAG	UNP Q13630
D	-5	LEU	-	EXPRESSION TAG	UNP Q13630
D	-4	TYR	-	EXPRESSION TAG	UNP Q13630
D	-3	PHE	-	EXPRESSION TAG	UNP Q13630
D	-2	GLN	-	EXPRESSION TAG	UNP Q13630
D	-1	SER	-	EXPRESSION TAG	UNP Q13630

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula:  $C_{21}H_{28}N_7O_{17}P_3$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula:  $C_{10}H_{15}N_5O_{11}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 28	C 10	N 5	O 11	P 2	0	0
3	B	1	Total 28	C 10	N 5	O 11	P 2	0	0
3	C	1	Total 28	C 10	N 5	O 11	P 2	0	0
3	D	1	Total 28	C 10	N 5	O 11	P 2	0	0

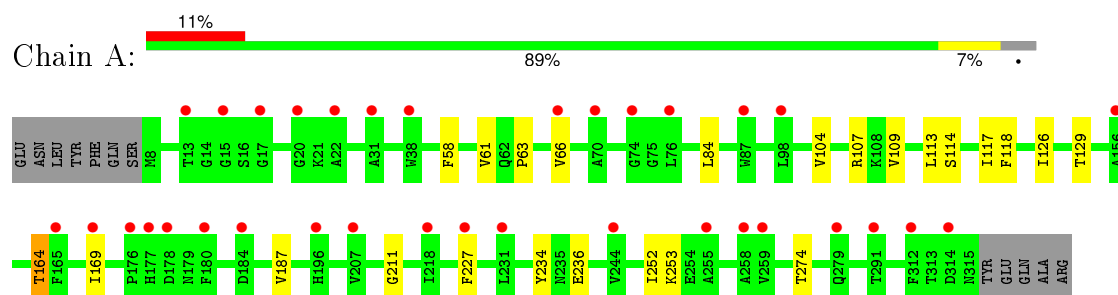
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	3	Total O 3 3	0	0
4	B	7	Total O 7 7	0	0
4	C	5	Total O 5 5	0	0
4	D	3	Total O 3 3	0	0

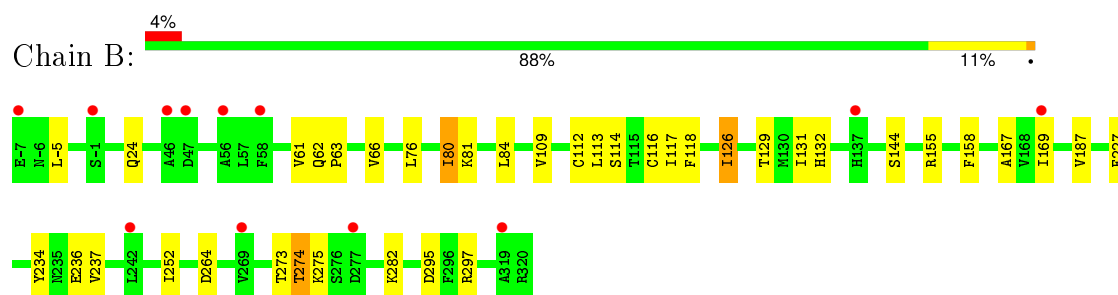
### 3 Residue-property plots [i](#)

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.

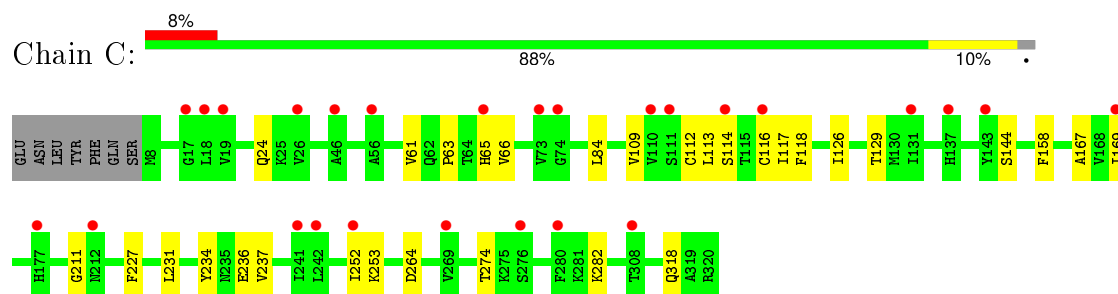
#### • Molecule 1: GDP-L-FUCOSE SYNTHASE



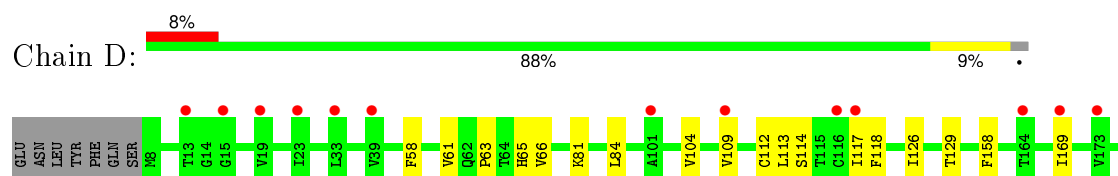
#### • Molecule 1: GDP-L-FUCOSE SYNTHASE

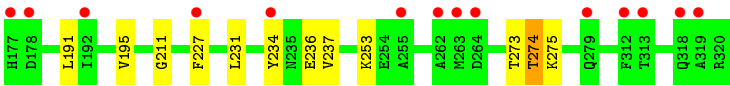


#### • Molecule 1: GDP-L-FUCOSE SYNTHASE



#### • Molecule 1: GDP-L-FUCOSE SYNTHASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.45Å 91.45Å 429.13Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	77.88 – 2.75 77.88 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.3 (77.88-2.75) 99.3 (77.88-2.75)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.52 (at 2.73Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, $R_{free}$	0.233 , 0.261 0.248 , 0.276	Depositor DCC
$R_{free}$ test set	1742 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	75.4	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 65.9	EDS
Estimated twinning fraction	0.064 for -h-k,k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 34584 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	9492	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.36 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.8892e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: GDP, NAP

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	0.40	0/2236	0.55	0/3062
1	B	0.41	0/2492	0.58	0/3401
1	C	0.40	0/2363	0.56	0/3233
1	D	0.41	0/2327	0.57	0/3184
All	All	0.40	0/9418	0.56	0/12880

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	2176	0	1813	11	0
1	B	2427	0	2218	23	0
1	C	2301	0	2052	13	0
1	D	2266	0	1983	20	0
2	A	48	0	25	1	0
2	B	48	0	25	1	0
2	C	48	0	25	2	0
2	D	48	0	25	1	0
3	A	28	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	28	0	12	0	0
3	C	28	0	12	0	0
3	D	28	0	12	0	0
4	A	3	0	0	0	0
4	B	7	0	0	0	0
4	C	5	0	0	0	0
4	D	3	0	0	0	0
All	All	9492	0	8214	59	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 3.

All (59) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:LYS:HG3	1:D:273:THR:HG21	1.78	0.65
1:A:107:ARG:O	1:A:164:THR:HG23	1.98	0.63
1:B:80:ILE:HD11	1:D:273:THR:O	2.01	0.60
1:B:273:THR:HG21	1:D:81:LYS:HG3	1.85	0.59
2:C:900:NAP:O1N	2:C:900:NAP:H2N	2.05	0.56
1:B:113:LEU:O	2:B:900:NAP:H6N	2.07	0.55
1:B:131:ILE:HG23	1:B:132:HIS:CD2	2.42	0.54
1:C:65:HIS:HB3	1:C:231:LEU:HD11	1.90	0.54
1:D:191:LEU:O	1:D:195:VAL:HG12	2.08	0.54
1:D:65:HIS:HB3	1:D:231:LEU:HD11	1.90	0.53
1:B:275:LYS:HG3	1:D:274:THR:HG22	1.91	0.52
1:C:61:VAL:O	1:C:63:PRO:HD3	2.11	0.51
1:C:114:SER:O	1:C:117:ILE:HG12	2.11	0.51
1:D:113:LEU:O	2:D:900:NAP:H6N	2.10	0.51
1:B:61:VAL:O	1:B:63:PRO:HD3	2.11	0.51
1:D:114:SER:O	1:D:117:ILE:HG12	2.11	0.50
1:B:114:SER:O	1:B:117:ILE:HG12	2.11	0.50
1:A:114:SER:O	1:A:117:ILE:HG12	2.11	0.50
1:B:76:LEU:O	1:B:80:ILE:HG23	2.12	0.49
1:C:113:LEU:HD13	1:C:167:ALA:HB1	1.94	0.49
1:D:61:VAL:O	1:D:63:PRO:HD3	2.13	0.49
1:B:113:LEU:HD13	1:B:167:ALA:HB1	1.94	0.49
1:A:61:VAL:O	1:A:63:PRO:HD3	2.13	0.49
1:B:274:THR:HG21	1:D:275:LYS:HE3	1.96	0.47
1:C:211:GLY:HA2	1:C:252:ILE:HB	1.97	0.47
1:B:274:THR:HG22	1:D:275:LYS:HG3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:TYR:CZ	1:A:236:GLU:HB2	2.50	0.47
1:C:234:TYR:CZ	1:C:236:GLU:HB2	2.51	0.46
1:D:234:TYR:CZ	1:D:236:GLU:HB2	2.50	0.46
1:D:118:PHE:HE2	1:D:169:ILE:HG21	1.82	0.45
1:B:234:TYR:CZ	1:B:236:GLU:HB2	2.51	0.45
1:B:80:ILE:HD11	1:D:273:THR:HG23	1.98	0.45
1:B:274:THR:CG2	1:D:275:LYS:HE3	2.47	0.45
1:C:118:PHE:HE2	1:C:169:ILE:HG21	1.82	0.44
1:A:113:LEU:O	2:A:900:NAP:H6N	2.17	0.44
1:A:118:PHE:HE2	1:A:169:ILE:HG21	1.83	0.44
1:B:80:ILE:CG1	1:D:273:THR:HG23	2.47	0.44
1:B:158:PHE:CG	1:B:237:VAL:HG21	2.53	0.44
1:B:118:PHE:HE2	1:B:169:ILE:HG21	1.82	0.44
1:C:158:PHE:CG	1:C:237:VAL:HG21	2.52	0.44
1:D:158:PHE:CG	1:D:237:VAL:HG21	2.54	0.43
1:B:66:VAL:HB	1:B:109:VAL:HG22	2.01	0.43
1:A:58:PHE:HB3	1:A:104:VAL:HG21	2.01	0.43
1:C:113:LEU:O	2:C:900:NAP:H6N	2.19	0.43
1:B:187:VAL:HG21	1:B:252:ILE:HD12	2.00	0.43
1:C:117:ILE:HG22	1:C:144:SER:HA	2.01	0.42
1:A:187:VAL:HG21	1:A:252:ILE:HD12	2.01	0.42
1:B:117:ILE:HG22	1:B:144:SER:HA	2.01	0.42
1:C:116:CYS:HB3	1:C:282:LYS:NZ	2.35	0.42
1:D:211:GLY:HA3	1:D:253:LYS:HG3	2.01	0.42
1:C:66:VAL:HB	1:C:109:VAL:HG22	2.02	0.42
1:D:58:PHE:HB3	1:D:104:VAL:HG21	2.00	0.42
1:D:66:VAL:HB	1:D:109:VAL:HG22	2.02	0.42
1:A:66:VAL:HB	1:A:109:VAL:HG22	2.02	0.42
1:C:211:GLY:HA3	1:C:253:LYS:HG3	2.02	0.41
1:A:211:GLY:HA3	1:A:253:LYS:HG3	2.02	0.41
1:A:211:GLY:HA2	1:A:252:ILE:HB	2.02	0.40
1:B:116:CYS:HB3	1:B:282:LYS:NZ	2.36	0.40
1:B:118:PHE:CD1	1:B:126:ILE:HG21	2.56	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	306/320 (96%)	291 (95%)	15 (5%)	0	100	100
1	B	318/320 (99%)	302 (95%)	15 (5%)	1 (0%)	46	77
1	C	311/320 (97%)	297 (96%)	14 (4%)	0	100	100
1	D	311/320 (97%)	297 (96%)	14 (4%)	0	100	100
All	All	1246/1280 (97%)	1187 (95%)	58 (5%)	1 (0%)	56	86

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	-5	LEU

### 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	183/274 (67%)	177 (97%)	6 (3%)	45	77
1	B	234/274 (85%)	221 (94%)	13 (6%)	26	56
1	C	215/274 (78%)	206 (96%)	9 (4%)	36	68
1	D	206/274 (75%)	200 (97%)	6 (3%)	50	81
All	All	838/1096 (76%)	804 (96%)	34 (4%)	37	69

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	LEU
1	A	126	ILE
1	A	129	THR
1	A	164	THR
1	A	227	PHE
1	A	274	THR
1	B	24	GLN
1	B	62	GLN
1	B	80	ILE
1	B	84	LEU
1	B	112	CYS
1	B	126	ILE
1	B	129	THR
1	B	155	ARG
1	B	227	PHE
1	B	264	ASP
1	B	274	THR
1	B	295	ASP
1	B	297	ARG
1	C	24	GLN
1	C	84	LEU
1	C	112	CYS
1	C	126	ILE
1	C	129	THR
1	C	227	PHE
1	C	264	ASP
1	C	274	THR
1	C	318	GLN
1	D	84	LEU
1	D	112	CYS
1	D	126	ILE
1	D	129	THR
1	D	227	PHE
1	D	274	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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$
2	NAP	A	900	-	42,52,52	0.58	0	54,80,80	1.01	3 (5%)
3	GDP	A	901	-	23,30,30	0.96	1 (4%)	30,47,47	2.17	5 (16%)
2	NAP	B	900	-	42,52,52	0.53	0	54,80,80	1.03	3 (5%)
3	GDP	B	901	-	23,30,30	0.92	1 (4%)	30,47,47	2.09	4 (13%)
2	NAP	C	900	-	42,52,52	0.52	0	54,80,80	1.07	3 (5%)
3	GDP	C	901	-	23,30,30	0.98	1 (4%)	30,47,47	2.12	4 (13%)
2	NAP	D	900	-	42,52,52	0.62	1 (2%)	54,80,80	0.99	2 (3%)
3	GDP	D	901	-	23,30,30	0.96	1 (4%)	30,47,47	2.15	5 (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
2	NAP	A	900	-	-	0/27/67/67	0/5/5/5
3	GDP	A	901	-	-	0/12/32/32	0/3/3/3
2	NAP	B	900	-	-	0/27/67/67	0/5/5/5
3	GDP	B	901	-	-	0/12/32/32	0/3/3/3
2	NAP	C	900	-	-	0/27/67/67	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GDP	C	901	-	-	0/12/32/32	0/3/3/3
2	NAP	D	900	-	-	0/27/67/67	0/5/5/5
3	GDP	D	901	-	-	0/12/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	900	NAP	O4D-C1D	-2.38	1.38	1.41
3	D	901	GDP	C6-N1	3.07	1.38	1.33
3	A	901	GDP	C6-N1	3.07	1.38	1.33
3	B	901	GDP	C6-N1	3.11	1.38	1.33
3	C	901	GDP	C6-N1	3.14	1.38	1.33

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	GDP	C5-C6-N1	-8.26	112.30	123.59
3	D	901	GDP	C5-C6-N1	-8.20	112.38	123.59
3	C	901	GDP	C5-C6-N1	-8.19	112.39	123.59
3	B	901	GDP	C5-C6-N1	-8.07	112.56	123.59
2	C	900	NAP	C4D-O4D-C1D	-3.16	106.25	109.72
3	B	901	GDP	N3-C2-N1	-2.67	123.38	127.44
3	C	901	GDP	N3-C2-N1	-2.65	123.41	127.44
3	D	901	GDP	N3-C2-N1	-2.60	123.48	127.44
3	A	901	GDP	N3-C2-N1	-2.52	123.61	127.44
3	A	901	GDP	C6-C5-C4	-2.36	118.08	120.90
3	D	901	GDP	C6-C5-C4	-2.18	118.29	120.90
3	B	901	GDP	C6-C5-C4	-2.11	118.37	120.90
3	C	901	GDP	C6-C5-C4	-2.09	118.41	120.90
2	B	900	NAP	C4D-O4D-C1D	2.08	112.01	109.72
2	A	900	NAP	O3-PA-O5B	2.09	108.47	102.94
3	A	901	GDP	PA-O3A-PB	2.16	139.90	132.67
3	D	901	GDP	PA-O3A-PB	2.28	140.30	132.67
2	D	900	NAP	O3-PA-O5B	2.34	109.16	102.94
2	A	900	NAP	O4D-C1D-N1N	2.46	110.83	108.13
2	C	900	NAP	O3-PA-O5B	2.48	109.50	102.94
2	B	900	NAP	O3-PA-O5B	2.66	109.99	102.94
2	B	900	NAP	O2B-P2B-O1X	4.97	119.53	107.11
2	C	900	NAP	O2B-P2B-O1X	5.08	119.80	107.11
2	A	900	NAP	O2B-P2B-O1X	5.15	119.97	107.11
2	D	900	NAP	O2B-P2B-O1X	5.29	120.31	107.11
3	B	901	GDP	C6-N1-C2	6.60	125.09	115.94

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Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	C	901	GDP	C6-N1-C2	6.64	125.16	115.94
3	D	901	GDP	C6-N1-C2	6.68	125.20	115.94
3	A	901	GDP	C6-N1-C2	6.71	125.25	115.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	900	NAP	1	0
2	B	900	NAP	1	0
2	C	900	NAP	2	0
2	D	900	NAP	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	308/320 (96%)	0.77	34 (11%) 7 4	62, 97, 126, 157	0
1	B	320/320 (100%)	0.44	12 (3%) 44 37	48, 68, 97, 140	0
1	C	313/320 (97%)	0.65	26 (8%) 14 9	59, 78, 104, 125	0
1	D	313/320 (97%)	0.63	27 (8%) 13 8	55, 87, 120, 148	0
All	All	1254/1280 (97%)	0.62	99 (7%) 15 10	48, 80, 119, 157	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	70	ALA	6.6
1	C	56	ALA	6.2
1	A	227	PHE	5.4
1	D	39	VAL	4.5
1	D	116	CYS	4.2
1	D	169	ILE	4.1
1	A	38	TRP	4.1
1	A	177	HIS	3.9
1	C	177	HIS	3.8
1	D	177	HIS	3.7
1	A	178	ASP	3.7
1	A	20	GLY	3.6
1	D	263	MET	3.6
1	A	17	GLY	3.6
1	D	319	ALA	3.5
1	D	109	VAL	3.4
1	C	46	ALA	3.4
1	C	241	ILE	3.3
1	B	58	PHE	3.3
1	A	156	ALA	3.2
1	D	312	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	13	THR	3.2
1	D	313	THR	3.2
1	A	98	LEU	3.1
1	A	13	THR	3.1
1	D	19	VAL	3.0
1	B	46	ALA	3.0
1	C	114	SER	3.0
1	D	33	LEU	2.9
1	A	196	HIS	2.9
1	A	218	ILE	2.8
1	A	244	VAL	2.8
1	A	255	ALA	2.8
1	D	101	ALA	2.8
1	D	262	ALA	2.8
1	C	169	ILE	2.7
1	C	242	LEU	2.7
1	D	164	THR	2.7
1	B	47	ASP	2.6
1	A	169	ILE	2.6
1	A	207	VAL	2.6
1	D	117	ILE	2.6
1	D	279	GLN	2.6
1	A	231	LEU	2.6
1	D	178	ASP	2.6
1	A	66	VAL	2.6
1	D	255	ALA	2.5
1	A	291	THR	2.5
1	C	111	SER	2.5
1	B	242	LEU	2.5
1	A	87	TRP	2.5
1	C	18	LEU	2.4
1	C	65	HIS	2.4
1	C	26	VAL	2.4
1	D	23	ILE	2.4
1	C	116	CYS	2.4
1	D	318	GLN	2.4
1	D	15	GLY	2.3
1	B	56	ALA	2.3
1	B	319	ALA	2.3
1	C	73	VAL	2.3
1	C	74	GLY	2.3
1	D	227	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	277	ASP	2.3
1	D	234	TYR	2.3
1	C	308	THR	2.3
1	A	279	GLN	2.3
1	C	17	GLY	2.2
1	A	76	LEU	2.2
1	A	31	ALA	2.2
1	A	314	ASP	2.2
1	A	15	GLY	2.2
1	D	264	ASP	2.2
1	C	143	TYR	2.2
1	B	-7	GLU	2.2
1	D	173	VAL	2.2
1	A	259	VAL	2.2
1	B	-1	SER	2.2
1	B	269	VAL	2.2
1	C	269	VAL	2.2
1	C	137	HIS	2.1
1	A	258	ALA	2.1
1	C	19	VAL	2.1
1	B	137	HIS	2.1
1	C	276	SER	2.1
1	C	252	ILE	2.1
1	D	192	ILE	2.1
1	A	312	PHE	2.1
1	C	131	ILE	2.1
1	A	74	GLY	2.1
1	A	176	PRO	2.0
1	A	180	PHE	2.0
1	C	212	ASN	2.0
1	A	165	PHE	2.0
1	B	169	ILE	2.0
1	A	22	ALA	2.0
1	C	110	VAL	2.0
1	A	184	ASP	2.0
1	C	280	PHE	2.0

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

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
2	NAP	A	900	48/48	0.92	0.24	-0.20	93,98,109,114	0
2	NAP	D	900	48/48	0.94	0.21	-0.52	90,98,106,110	0
2	NAP	B	900	48/48	0.94	0.20	-0.58	59,65,71,76	0
3	GDP	C	901	28/28	0.94	0.20	-0.59	65,69,73,75	0
2	NAP	C	900	48/48	0.92	0.22	-0.61	68,76,88,92	0
3	GDP	D	901	28/28	0.95	0.18	-0.63	60,69,78,80	0
3	GDP	B	901	28/28	0.96	0.19	-0.65	51,55,58,59	0
3	GDP	A	901	28/28	0.93	0.19	-0.81	80,85,88,89	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.