



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 11, 2017 – 03:34 PM EDT

PDB ID : 5BTF
Title : Crystal structure of a topoisomerase II complex
Authors : Blower, T.R.; Williamson, B.H.; Kerns, R.J.; Berger, J.M.
Deposited on : unknown
Resolution : 2.61 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20029824

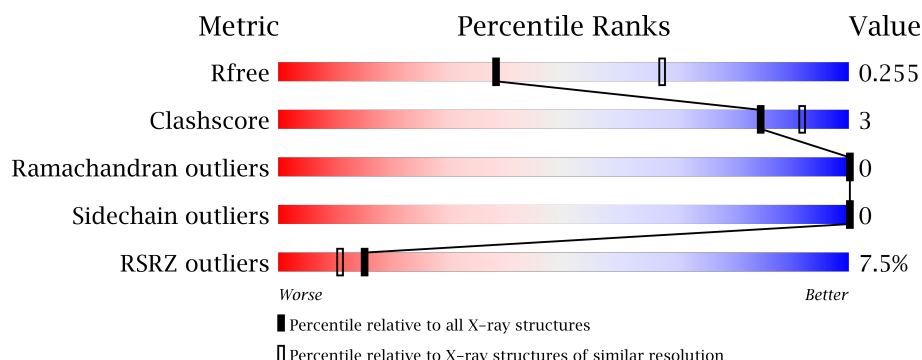
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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.



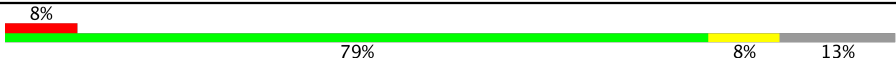


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	2983 (2.64-2.60)
Clashscore	112137	3351 (2.64-2.60)
Ramachandran outliers	110173	3298 (2.64-2.60)
Sidechain outliers	110143	3298 (2.64-2.60)
RSRZ outliers	101464	2992 (2.64-2.60)

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	503	<div> <div>6%</div> <div>93%</div> <div>• •</div> </div>
1	C	503	<div> <div>10%</div> <div>91%</div> <div>5% •</div> </div>
2	B	253	<div> <div>5%</div> <div>92%</div> <div>5% •</div> </div>
2	D	253	<div> <div>3%</div> <div>91%</div> <div>7% •</div> </div>
3	E	24	<div> <div>21%</div> <div>75%</div> <div>13% 13%</div> </div>

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Mol	Chain	Length	Quality of chain
3	H	24	
4	F	24	
4	G	24	

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
1	PTR	A	129	-	-	X	-
1	PTR	C	129	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 25958 atoms, of which 12513 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 DNA gyrase subunit A.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	487	Total	C	H	N	O	P	S	0	1	0
			7692	2388	3853	704	733	1	13			
1	C	487	Total	C	H	N	O	P	S	0	1	0
			7691	2388	3852	704	733	1	13			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	90	SER	ALA	engineered mutation	UNP P9WG47
A	501	ILE	-	expression tag	UNP P9WG47
A	502	GLY	-	expression tag	UNP P9WG47
A	503	SER	-	expression tag	UNP P9WG47
A	504	GLY	-	expression tag	UNP P9WG47
C	90	SER	ALA	engineered mutation	UNP P9WG47
C	501	ILE	-	expression tag	UNP P9WG47
C	502	GLY	-	expression tag	UNP P9WG47
C	503	SER	-	expression tag	UNP P9WG47
C	504	GLY	-	expression tag	UNP P9WG47

- Molecule 2 is a protein called DNA gyrase subunit B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	245	Total	C	H	N	O	S	0	0	0
			3906	1217	1974	348	360	7			
2	D	247	Total	C	H	N	O	S	0	0	0
			3930	1224	1984	351	364	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	423	SER	-	expression tag	UNP P9WG45
B	424	ASN	-	expression tag	UNP P9WG45

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Chain	Residue	Modelled	Actual	Comment	Reference
B	425	ALA	-	expression tag	UNP P9WG45
D	423	SER	-	expression tag	UNP P9WG45
D	424	ASN	-	expression tag	UNP P9WG45
D	425	ALA	-	expression tag	UNP P9WG45

- Molecule 3 is a DNA chain called DNA substrate 24-mer GGTCATGAATGACTATGCAC GTAA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	E	21	Total	C	H	N	O	P	0	21	0
			618	196	208	74	120	20			
3	H	21	Total	C	H	N	O	P	0	21	0
			612	196	202	74	120	20			

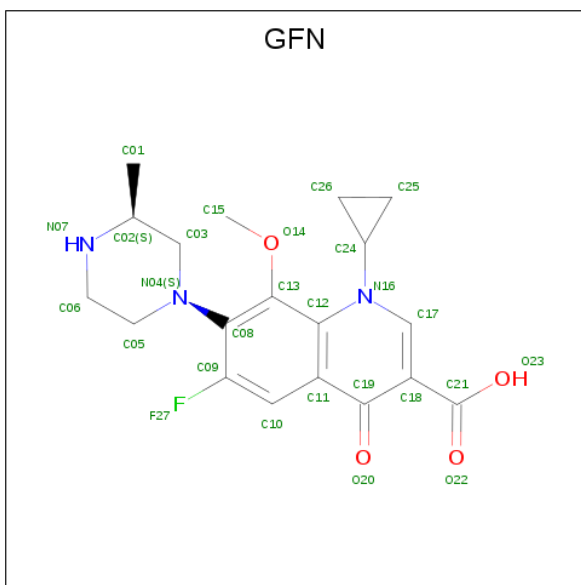
- Molecule 4 is a DNA chain called DNA substrate 24-mer TTACGTGCATAGTCATTTCAT GACC.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	F	21	Total	C	H	N	O	P	0	21	0
			613	196	203	74	120	20			
4	G	21	Total	C	H	N	O	P	0	21	0
			605	196	195	74	120	20			

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		
5	D	1	Total	Mg	0	0
			1	1		
5	C	1	Total	Mg	0	0
			1	1		
5	E	1	Total	Mg	0	0
			1	1		

- Molecule 6 is 1-cyclopropyl-6-fluoro-8-methoxy-7-[(3S)-3-methylpiperazin-1-yl]-4-oxo-1,4-dihydroquinoline-3-carboxylic acid (three-letter code: GFN) (formula: C₁₉H₂₂FN₃O₄).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	F	1	Total	C	F	H	N	O	0	0
			48	19	1	21	3	4		
6	G	1	Total	C	F	H	N	O	0	0
			48	19	1	21	3	4		

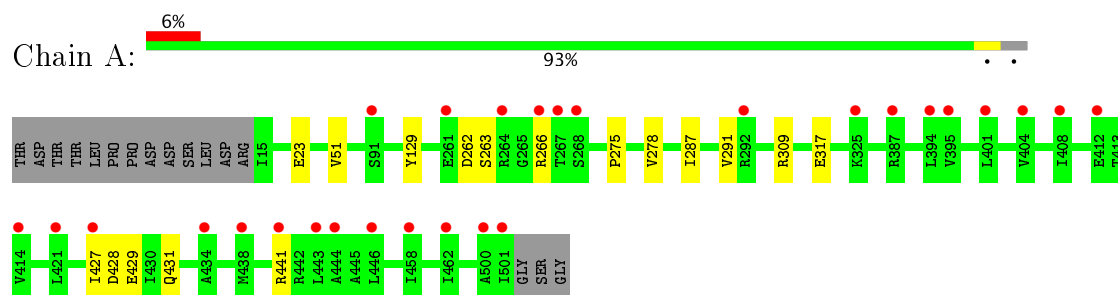
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	85	Total	O	0	0
			85	85		
7	B	32	Total	O	0	0
			32	32		
7	C	35	Total	O	0	0
			35	35		
7	D	17	Total	O	0	0
			17	17		
7	E	6	Total	O	0	0
			6	6		
7	F	6	Total	O	0	0
			6	6		
7	G	5	Total	O	0	0
			5	5		
7	H	5	Total	O	0	0
			5	5		

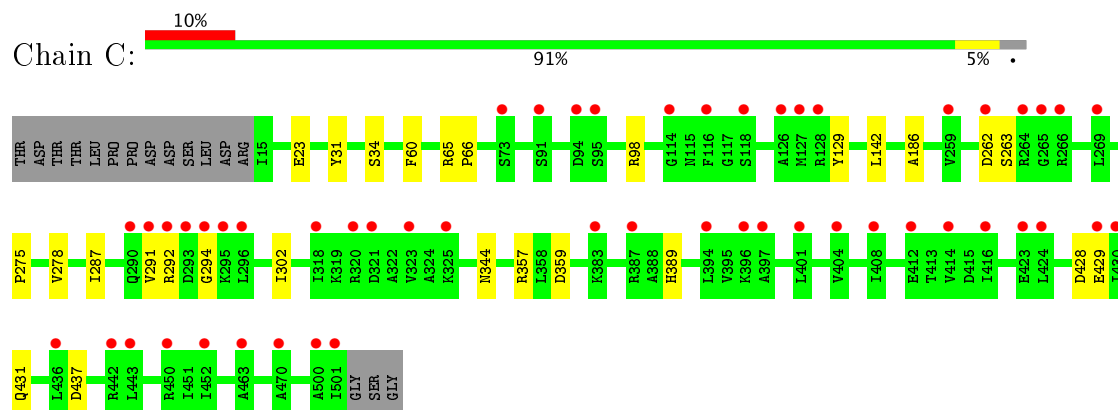
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

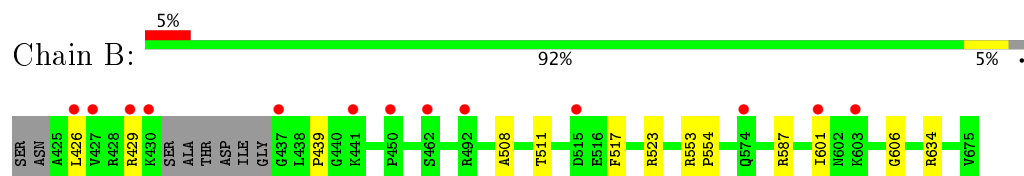
- Molecule 1: DNA gyrase subunit A



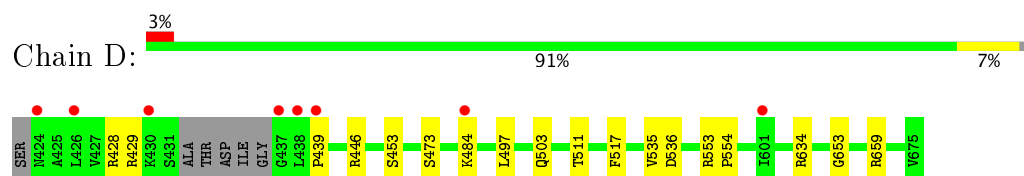
- Molecule 1: DNA gyrase subunit A




- Molecule 2: DNA gyrase subunit B

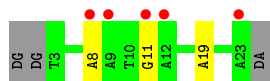


- Molecule 2: DNA gyrase subunit B




- Molecule 3: DNA substrate 24-mer GGTCATGAATGACTATGCACGTAA

Chain E: 




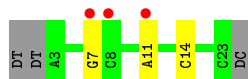
- Molecule 3: DNA substrate 24-mer GGTCATGAATGACTATGCACGTAA

Chain H: 




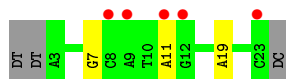
- Molecule 4: DNA substrate 24-mer TTACGTGCATAGTCATTCATGACC

Chain F: 



- Molecule 4: DNA substrate 24-mer TTACGTGCATAGTCATTCATGACC

Chain G: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	108.95Å 84.02Å 130.59Å 90.00° 108.87° 90.00°	Depositor
Resolution (Å)	49.78 – 2.61 49.78 – 2.61	Depositor EDS
% Data completeness (in resolution range)	93.9 (49.78-2.61) 93.7 (49.78-2.61)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 2.61Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.230 , 0.261 0.226 , 0.255	Depositor DCC
R_{free} test set	3211 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	58.9	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 44.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	25958	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.34% 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, 0.333 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: MG, GFN, PTR

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	0.22	0/3882	0.40	0/5254
1	C	0.22	0/3882	0.40	0/5254
2	B	0.23	0/1960	0.39	0/2634
2	D	0.23	0/1974	0.39	0/2653
3	E	0.52	0/458	0.95	0/704
3	H	0.58	0/458	0.97	0/704
4	F	0.53	0/458	0.94	0/704
4	G	0.55	0/458	0.92	0/704
All	All	0.29	0/13530	0.52	0/18611

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	3839	3853	3853	16	0
1	C	3839	3852	3853	24	0
2	B	1932	1974	1973	8	0
2	D	1946	1984	1984	15	0
3	E	410	208	208	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	410	202	202	4	0
4	F	410	203	203	5	0
4	G	410	195	195	5	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
6	F	27	21	0	0	0
6	G	27	21	0	0	0
7	A	85	0	0	1	0
7	B	32	0	0	1	0
7	C	35	0	0	0	0
7	D	17	0	0	0	0
7	E	6	0	0	0	0
7	F	6	0	0	0	0
7	G	5	0	0	0	0
7	H	5	0	0	0	0
All	All	13445	12513	12471	58	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 (58) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:129:PTR:O3P	4:G:11[B]:DA:C5'	2.04	1.06
1:C:287:ILE:O	1:C:291:VAL:HG23	1.57	1.04
1:A:129:PTR:O3P	3:H:11[B]:DG:C5'	2.10	0.98
1:C:129:PTR:O3P	3:E:11[A]:DG:H4'	1.66	0.95
1:A:287:ILE:O	1:A:291:VAL:HG23	1.66	0.94
1:A:129:PTR:O3P	4:F:11[A]:DA:C5'	2.19	0.90
1:C:98:ARG:NH1	4:F:7[A]:DG:OP1	2.12	0.82
2:D:653:GLY:O	2:D:659:ARG:NH2	2.13	0.81
1:A:266:ARG:HD3	1:A:317:GLU:OE2	1.83	0.78
1:A:129:PTR:O3P	3:H:11[B]:DG:C4'	2.31	0.77
1:A:129:PTR:O3P	4:F:11[A]:DA:H4'	1.84	0.77
1:C:129:PTR:O3P	3:E:11[A]:DG:C4'	2.33	0.76
1:A:129:PTR:O3P	3:H:11[B]:DG:H4'	1.86	0.76
1:A:129:PTR:O3P	4:F:11[A]:DA:C4'	2.37	0.72
1:C:31:TYR:HA	2:D:535:VAL:CG1	2.22	0.70
2:D:429:ARG:NH1	2:D:439:PRO:O	2.26	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:129:PTR:O3P	4:G:11[B]:DA:C4'	2.39	0.68
1:C:357:ARG:NH1	1:C:359:ASP:OD1	2.27	0.67
1:C:129:PTR:O3P	3:E:11[A]:DG:C5'	2.43	0.67
4:G:7[B]:DG:N2	3:H:18[B]:DC:O2	2.28	0.65
1:C:31:TYR:HA	2:D:535:VAL:HG13	1.81	0.62
2:B:508:ALA:O	2:B:523:ARG:NE	2.32	0.62
1:C:129:PTR:O3P	4:G:11[B]:DA:H4'	1.99	0.61
2:D:484:LYS:NZ	2:D:536:ASP:O	2.36	0.57
2:B:429:ARG:NH1	2:B:439:PRO:O	2.37	0.57
1:C:262:ASP:OD1	1:C:263:SER:N	2.36	0.55
1:C:34:SER:OG	2:D:535:VAL:HG11	2.08	0.54
2:D:659:ARG:NH1	3:E:19[A]:DA:OP1	2.42	0.53
2:B:426:LEU:HD21	2:B:523:ARG:HD2	1.91	0.52
1:C:291:VAL:HG21	1:C:302:ILE:HD11	1.92	0.51
2:D:428:ARG:NH2	2:D:453:SER:OG	2.43	0.51
1:C:428:ASP:OD1	1:C:429:GLU:N	2.43	0.50
2:D:659:ARG:NH1	4:G:19[B]:DA:OP1	2.45	0.50
3:E:11[A]:DG:N2	4:F:14[A]:DC:O2	2.45	0.49
2:D:497:LEU:O	2:D:503:GLN:NE2	2.44	0.48
2:D:446:ARG:NH2	2:D:473:SER:OG	2.48	0.46
2:B:511:THR:HG21	2:B:517:PHE:CE2	2.51	0.46
1:C:389:HIS:O	1:C:431:GLN:NE2	2.46	0.46
1:A:51:VAL:HG11	3:E:8[A]:DA:H5'	1.98	0.45
1:C:186:ALA:O	1:C:344:ASN:ND2	2.50	0.45
1:A:309:ARG:NH1	7:A:605:HOH:O	2.46	0.44
1:C:31:TYR:HD1	2:D:535:VAL:HG12	1.82	0.44
1:C:292:ARG:C	1:C:294:GLY:H	2.21	0.44
1:C:23:GLU:OE1	2:D:634:ARG:NH1	2.50	0.44
2:B:587:ARG:NH2	7:B:804:HOH:O	2.50	0.43
1:A:441:ARG:NE	1:C:437:ASP:OD1	2.48	0.43
1:A:275:PRO:HG2	1:A:278:VAL:HG21	2.00	0.43
1:A:427:ILE:CG2	1:A:431:GLN:HB2	2.49	0.43
1:A:262:ASP:OD2	1:A:263:SER:N	2.51	0.42
2:D:511:THR:HG21	2:D:517:PHE:CE2	2.54	0.42
2:B:553:ARG:N	2:B:554:PRO:CD	2.82	0.42
1:A:23:GLU:OE1	2:B:634:ARG:NH1	2.49	0.42
2:D:553:ARG:N	2:D:554:PRO:CD	2.83	0.42
1:C:60:PHE:HB2	1:C:142:LEU:HD13	2.03	0.41
1:C:275:PRO:HG2	1:C:278:VAL:HG21	2.03	0.41
1:A:428:ASP:OD1	1:A:429:GLU:N	2.44	0.40
2:B:601:ILE:HG22	2:B:606:GLY:HA2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:65:ARG:HB3	1:C:66:PRO:HD2	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	485/503 (96%)	469 (97%)	16 (3%)	0	100	100
1	C	485/503 (96%)	469 (97%)	16 (3%)	0	100	100
2	B	241/253 (95%)	233 (97%)	8 (3%)	0	100	100
2	D	243/253 (96%)	235 (97%)	8 (3%)	0	100	100
All	All	1454/1512 (96%)	1406 (97%)	48 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	408/421 (97%)	408 (100%)	0	100	100
1	C	408/421 (97%)	408 (100%)	0	100	100
2	B	205/211 (97%)	205 (100%)	0	100	100
2	D	207/211 (98%)	207 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1228/1264 (97%)	1228 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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 ⓘ

2 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$
1	PTR	A	129	1	15,16,17	0.65	0	19,22,24	0.97	1 (5%)
1	PTR	C	129	1	15,16,17	0.64	0	19,22,24	1.03	1 (5%)

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
1	PTR	A	129	1	-	0/9/11/13	0/1/1/1
1	PTR	C	129	1	-	0/9/11/13	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	129	PTR	CB-CA-C	-2.72	106.17	111.41
1	A	129	PTR	CB-CA-C	-2.46	106.67	111.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	129	PTR	6	0
1	C	129	PTR	6	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GFN	F	101	5	21,30,30	1.57	6 (28%)	26,45,45	1.50	4 (15%)
6	GFN	G	101	5	21,30,30	1.54	6 (28%)	26,45,45	1.98	6 (23%)

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
6	GFN	F	101	5	-	0/10/26/26	0/3/4/4
6	GFN	G	101	5	-	0/10/26/26	0/3/4/4

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	G	101	GFN	C11-C12	-3.25	1.39	1.41
6	F	101	GFN	C11-C12	-3.21	1.39	1.41
6	F	101	GFN	C02-N07	-3.16	1.43	1.48
6	G	101	GFN	C02-N07	-2.71	1.44	1.48
6	G	101	GFN	C05-N04	-2.64	1.42	1.46
6	F	101	GFN	C05-N04	-2.33	1.43	1.46
6	F	101	GFN	C08-N04	2.35	1.47	1.40
6	G	101	GFN	C08-N04	2.38	1.47	1.40
6	G	101	GFN	O14-C13	2.47	1.41	1.38
6	F	101	GFN	O14-C13	2.49	1.42	1.38
6	G	101	GFN	C10-C09	2.60	1.39	1.35
6	F	101	GFN	C10-C09	2.73	1.39	1.35

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	G	101	GFN	C10-C11-C19	-2.60	117.45	121.69
6	G	101	GFN	C17-C18-C19	-2.57	118.79	119.99
6	F	101	GFN	C10-C11-C19	-2.56	117.52	121.69
6	F	101	GFN	C17-C18-C19	-2.44	118.85	119.99
6	G	101	GFN	C24-N16-C12	2.06	123.50	120.87
6	F	101	GFN	C19-C11-C12	2.72	121.92	119.90
6	G	101	GFN	C19-C11-C12	2.95	122.09	119.90
6	F	101	GFN	C13-C08-C09	3.32	119.24	116.01
6	G	101	GFN	C13-C08-C09	3.35	119.28	116.01
6	G	101	GFN	C06-N07-C02	6.14	117.04	111.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	486/503 (96%)	0.62	28 (5%) 24 18	34, 60, 124, 162	0
1	C	486/503 (96%)	0.79	52 (10%) 7 4	37, 62, 118, 167	0
2	B	245/253 (96%)	0.66	13 (5%) 27 21	43, 69, 106, 134	0
2	D	247/253 (97%)	0.50	8 (3%) 48 41	38, 65, 108, 139	0
3	E	21/24 (87%)	1.27	5 (23%) 1 0	65, 99, 164, 178	21 (100%)
3	H	21/24 (87%)	1.12	2 (9%) 9 6	27, 58, 128, 169	21 (100%)
4	F	21/24 (87%)	1.15	3 (14%) 3 1	55, 102, 174, 201	21 (100%)
4	G	21/24 (87%)	1.27	5 (23%) 1 0	32, 57, 147, 169	21 (100%)
All	All	1548/1608 (96%)	0.69	116 (7%) 15 11	27, 64, 124, 201	84 (5%)

All (116) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	501	ILE	9.8
1	C	404	VAL	6.8
1	A	501	ILE	6.8
2	B	426	LEU	5.2
1	A	266	ARG	5.2
2	D	437	GLY	5.1
1	A	401	LEU	4.9
1	C	408	ILE	4.7
1	A	292	ARG	4.7
1	A	412	GLU	4.7
1	A	414	VAL	4.6
1	A	387	ARG	4.5
1	C	294	GLY	4.4
1	A	421	LEU	4.3
1	C	414	VAL	4.2
3	E	23[A]	DA	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	292	ARG	4.1
1	C	265	GLY	4.0
4	G	23[B]	DC	4.0
1	C	442	ARG	3.9
2	D	426	LEU	3.8
2	B	603	LYS	3.5
2	B	437	GLY	3.5
1	C	264	ARG	3.4
1	C	500	ALA	3.4
1	A	395	VAL	3.3
1	A	427	ILE	3.3
2	B	601	ILE	3.3
1	C	394	LEU	3.3
1	C	325	LYS	3.3
1	A	408	ILE	3.2
2	D	424	ASN	3.2
1	C	266	ARG	3.1
1	A	394	LEU	3.0
1	A	443	LEU	3.0
1	C	412	GLU	3.0
4	G	11[B]	DA	3.0
3	E	11[A]	DG	3.0
1	C	463	ALA	3.0
1	C	430	ILE	2.9
1	C	259	VAL	2.9
1	C	114	GLY	2.9
3	H	11[B]	DG	2.9
1	C	450	ARG	2.8
1	C	295	LYS	2.8
1	A	404	VAL	2.8
1	C	323	VAL	2.8
1	C	383	LYS	2.8
1	C	396	LYS	2.8
3	H	8[B]	DA	2.7
1	A	261	GLU	2.7
1	C	94	ASP	2.7
3	E	8[A]	DA	2.7
1	C	128[A]	ARG	2.7
1	A	268	SER	2.7
2	D	430	LYS	2.7
1	C	118	SER	2.6
1	C	452	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	127	MET	2.6
2	B	427	VAL	2.6
2	B	429	ARG	2.5
1	C	423	GLU	2.5
1	C	318	ILE	2.5
1	A	444	ALA	2.5
1	A	500	ALA	2.5
1	C	387	ARG	2.5
4	G	8[B]	DC	2.5
4	F	8[A]	DC	2.5
1	A	441	ARG	2.4
1	C	397	ALA	2.4
1	C	296	LEU	2.4
4	G	12[B]	DG	2.4
2	D	439	PRO	2.4
2	B	430	LYS	2.4
2	D	484	LYS	2.4
1	C	320	ARG	2.4
1	C	126	ALA	2.4
1	C	401	LEU	2.3
1	C	470	ALA	2.3
1	A	264	ARG	2.3
1	A	438	MET	2.3
1	C	321	ASP	2.3
4	G	9[B]	DA	2.3
2	D	438	LEU	2.3
1	C	95	SER	2.3
1	A	462	ILE	2.3
2	B	441	LYS	2.3
4	F	11[A]	DA	2.3
1	C	269	LEU	2.3
2	B	492	ARG	2.3
3	E	9[A]	DA	2.2
3	E	12[A]	DA	2.2
2	B	462	SER	2.2
1	A	446	LEU	2.2
1	C	91	SER	2.2
1	A	434	ALA	2.2
1	C	262	ASP	2.2
1	C	290	GLN	2.2
1	A	91	SER	2.1
4	F	7[A]	DG	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	436	LEU	2.1
1	C	291	VAL	2.1
1	C	416	ILE	2.1
2	B	450	PRO	2.1
1	A	458	ILE	2.1
1	C	443	LEU	2.1
1	A	325	LYS	2.1
1	C	73	SER	2.1
1	C	424	LEU	2.1
1	A	267	THR	2.1
1	C	293	ASP	2.1
1	C	429	GLU	2.0
2	B	515	ASP	2.0
2	D	601	ILE	2.0
2	B	574	GLN	2.0
1	C	116	PHE	2.0

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. 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
1	PTR	C	129	16/17	0.97	0.32	-	52,55,65,66	0
1	PTR	A	129	16/17	0.97	0.24	-	50,54,64,65	0

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, 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(\AA^2)	Q<0.9
5	MG	E	101	1/1	0.85	0.32	0.66	68,68,68,68	0
5	MG	C	601	1/1	0.75	0.30	0.49	83,83,83,83	0
6	GFN	G	101	27/27	0.92	0.28	0.22	48,58,70,70	0
6	GFN	F	101	27/27	0.92	0.28	0.05	46,55,68,70	0
5	MG	B	701	1/1	0.96	0.28	-0.06	41,41,41,41	0
5	MG	D	701	1/1	0.93	0.21	-0.46	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.