



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 04:51 PM EST

PDB ID : 5B3V
Title : Crystal structure of biliverdin reductase in complex with biliverdin and NADP+ from Synechocystis sp. PCC 6803
Authors : Takao, H.; Wada, K.
Deposited on : 2016-03-13
Resolution : 2.59 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
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) : rb-20028442

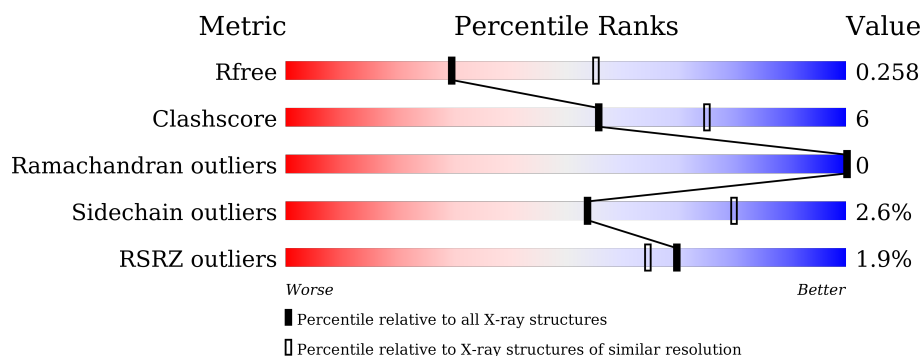
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	331	<div> <div>0%</div> <div>82%</div> <div>13%</div> <div>• •</div> </div>
1	B	331	<div> <div>2%</div> <div>82%</div> <div>13%</div> <div>• 5%</div> </div>
1	C	331	<div> <div>0%</div> <div>83%</div> <div>12%</div> <div>5%</div> </div>
1	D	331	<div> <div>2%</div> <div>82%</div> <div>13%</div> <div>• 5%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BLA	B	402	-	-	-	X
3	BLA	C	403	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10756 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.

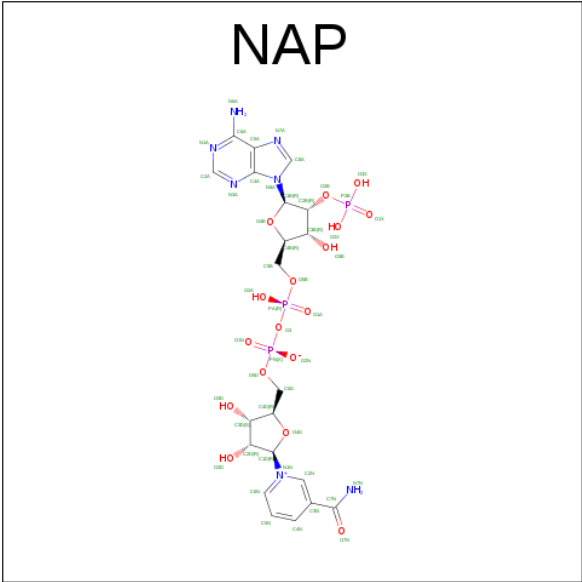
- Molecule 1 is a protein called Biliverdin reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	316	Total	C	N	O	S	0	0	0
			2497	1590	446	455	6			
1	A	317	Total	C	N	O	S	0	0	0
			2502	1593	447	456	6			
1	C	316	Total	C	N	O	S	0	0	0
			2497	1590	446	455	6			
1	D	316	Total	C	N	O	S	0	0	0
			2497	1590	446	455	6			

There are 12 discrepancies between the modelled and reference sequences:

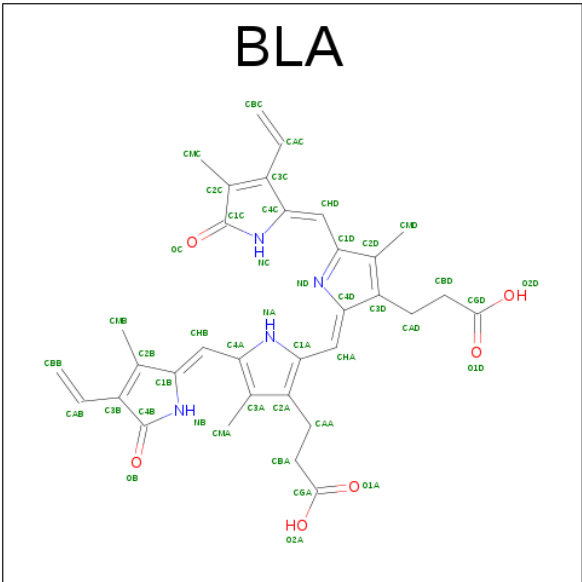
Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	expression tag	UNP P72782
B	-1	SER	-	expression tag	UNP P72782
B	0	HIS	-	expression tag	UNP P72782
A	-2	GLY	-	expression tag	UNP P72782
A	-1	SER	-	expression tag	UNP P72782
A	0	HIS	-	expression tag	UNP P72782
C	-2	GLY	-	expression tag	UNP P72782
C	-1	SER	-	expression tag	UNP P72782
C	0	HIS	-	expression tag	UNP P72782
D	-2	GLY	-	expression tag	UNP P72782
D	-1	SER	-	expression tag	UNP P72782
D	0	HIS	-	expression tag	UNP P72782

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	A	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 BILIVERDINE IX ALPHA (three-letter code: BLA) (formula: C₃₃H₃₄N₄O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			43	33	4	6		
3	B	1	Total	C	N	O	0	0
			43	33	4	6		
3	A	1	Total	C	N	O	0	0
			43	33	4	6		
3	A	1	Total	C	N	O	0	0
			43	33	4	6		
3	C	1	Total	C	N	O	0	0
			43	33	4	6		
3	C	1	Total	C	N	O	0	0
			43	33	4	6		
3	D	1	Total	C	N	O	0	0
			43	33	4	6		
3	D	1	Total	C	N	O	0	0
			43	33	4	6		

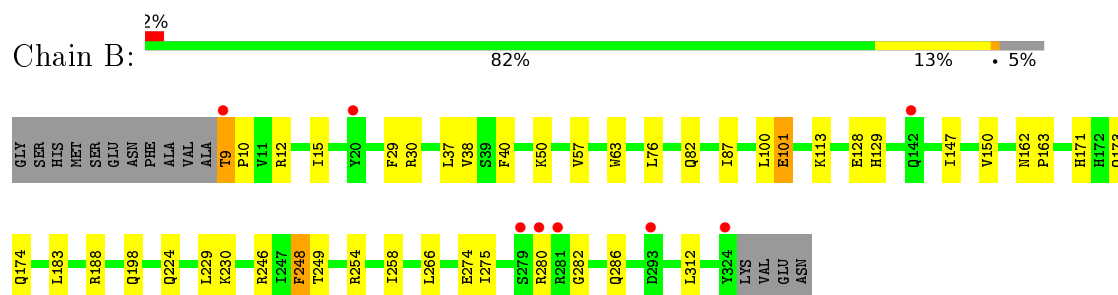
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	56	Total	O	0	0
			56	56		
4	A	71	Total	O	0	0
			71	71		
4	C	45	Total	O	0	0
			45	45		
4	D	55	Total	O	0	0
			55	55		

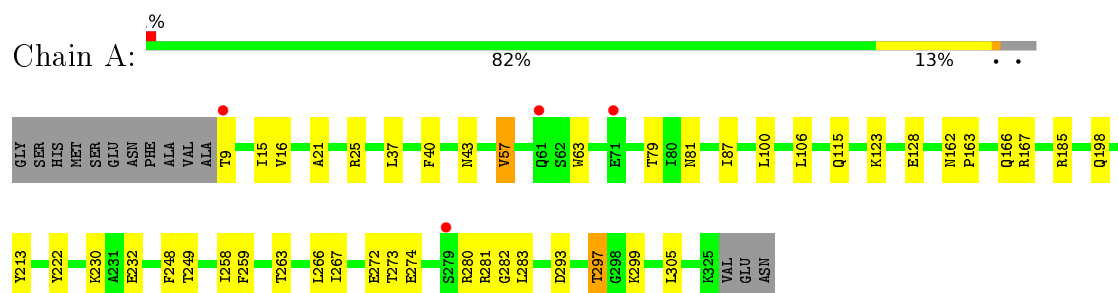
3 Residue-property plots

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.

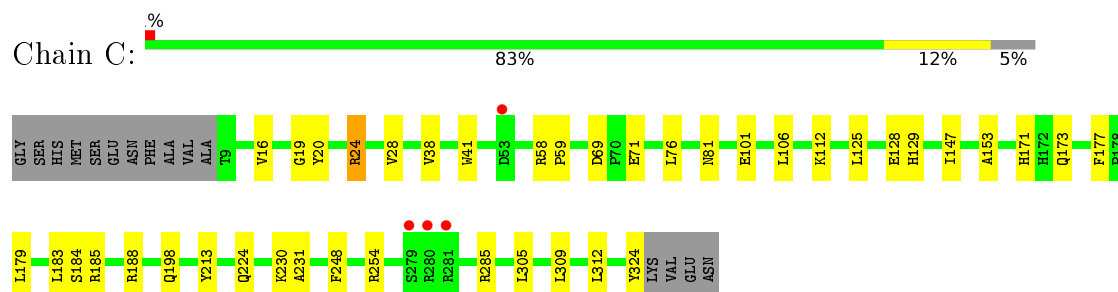
• Molecule 1: Biliverdin reductase



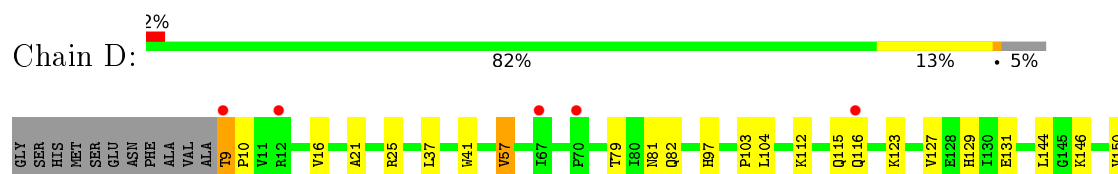
• Molecule 1: Biliverdin reductase



• Molecule 1: Biliverdin reductase



• Molecule 1: Biliverdin reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.50Å 100.41Å 138.03Å 90.00° 95.53° 90.00°	Depositor
Resolution (Å)	49.62 – 2.59 49.63 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.1 (49.62-2.59) 99.1 (49.63-2.59)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.58Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, R_{free}	0.208 , 0.258 0.206 , 0.258	Depositor DCC
R_{free} test set	2214 reflections (4.77%)	DCC
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10756	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.02% 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: NAP, BLA

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	0.28	0/2557	0.50	0/3467
1	B	0.37	0/2552	0.51	0/3460
1	C	0.37	0/2552	0.55	0/3460
1	D	0.29	0/2552	0.50	0/3460
All	All	0.33	0/10213	0.52	0/13847

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	2502	0	2467	24	0
1	B	2497	0	2465	27	0
1	C	2497	0	2465	23	0
1	D	2497	0	2465	24	0
2	A	48	0	24	3	0
2	B	48	0	25	1	0
2	C	48	0	24	3	0
2	D	48	0	25	1	0
3	A	86	0	64	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	86	0	64	11	0
3	C	86	0	64	9	0
3	D	86	0	64	7	0
4	A	71	0	0	0	0
4	B	56	0	0	0	0
4	C	45	0	0	0	0
4	D	55	0	0	0	0
All	All	10756	0	10216	128	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 6.

All (128) 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)
3:C:402:BLA:HBD2	3:C:403:BLA:HMA2	1.47	0.94
3:C:402:BLA:HMC1	3:C:402:BLA:HBC1	1.58	0.84
3:B:403:BLA:HBC1	3:B:403:BLA:HMC1	1.59	0.83
3:C:403:BLA:HBC1	3:C:403:BLA:HMC1	1.61	0.83
3:D:402:BLA:HMC1	3:D:402:BLA:HBC1	1.61	0.83
3:B:402:BLA:HBB1	3:B:402:BLA:HMB1	1.60	0.82
3:D:403:BLA:HMB1	3:D:403:BLA:HBB1	1.61	0.82
3:C:402:BLA:HMB1	3:C:402:BLA:HBB1	1.61	0.81
3:A:402:BLA:HBC1	3:A:402:BLA:HMC1	1.63	0.80
3:A:403:BLA:HBC1	3:A:403:BLA:HMC1	1.63	0.79
3:B:402:BLA:HAD2	3:B:403:BLA:HBD2	1.64	0.79
1:B:171:HIS:HE1	1:B:173:GLN:HB2	1.46	0.77
3:B:402:BLA:HHB	3:B:402:BLA:HBC1	1.67	0.75
1:B:280:ARG:HD2	3:B:402:BLA:HBC2	1.67	0.75
1:B:101:GLU:HB3	2:B:401:NAP:H1D	1.72	0.72
1:B:9:THR:HG22	1:B:10:PRO:HA	1.71	0.71
1:D:82:GLN:HB2	1:D:174:GLN:HA	1.72	0.71
1:B:171:HIS:CE1	1:B:173:GLN:HB2	2.24	0.71
1:D:37:LEU:HG	1:D:57:VAL:HG11	1.78	0.65
1:D:267:ILE:HG12	1:D:272:GLU:HG3	1.77	0.65
1:B:9:THR:CG2	1:B:10:PRO:HA	2.27	0.64
1:A:37:LEU:HG	1:A:57:VAL:HG11	1.80	0.62
1:C:185:ARG:NH2	3:C:403:BLA:HBA1	2.16	0.61
2:D:401:NAP:H72N	3:D:402:BLA:HB	1.49	0.61
1:A:115:GLN:HG3	1:A:305:LEU:HD23	1.83	0.60
1:D:129:HIS:HE1	1:D:183:LEU:HD22	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:ASN:HB3	2:A:401:NAP:O3X	2.03	0.59
1:C:58:ARG:HD3	1:C:59:PRO:HD2	1.85	0.59
3:D:403:BLA:HBC1	3:D:403:BLA:HHD	1.85	0.59
1:C:230:LYS:HD2	1:C:230:LYS:N	2.18	0.58
1:C:38:VAL:HG11	1:C:71:GLU:HG3	1.84	0.58
1:C:101:GLU:HB2	2:C:401:NAP:H1D	1.86	0.58
1:B:198:GLN:HB3	1:B:224:GLN:HB2	1.86	0.57
1:A:81:ASN:HB3	1:A:106:LEU:HD21	1.85	0.57
1:D:150:VAL:HG11	1:D:250:LEU:HD23	1.87	0.57
1:D:266:LEU:HB2	1:D:275:ILE:HD11	1.88	0.55
1:A:79:THR:HA	2:A:401:NAP:H4B	1.89	0.55
1:B:9:THR:HA	1:B:10:PRO:C	2.28	0.53
1:A:9:THR:O	1:A:9:THR:HG23	2.07	0.53
1:C:19:GLY:HA3	2:C:401:NAP:O1A	2.08	0.53
1:B:37:LEU:HG	1:B:57:VAL:HG21	1.90	0.53
1:C:112:LYS:HD3	1:C:309:LEU:HD21	1.89	0.52
1:D:281:ARG:HB3	1:D:282:GLY:HA3	1.90	0.52
1:B:249:THR:HG23	1:B:258:ILE:HG12	1.90	0.52
1:B:12:ARG:HB3	1:B:38:VAL:HG21	1.90	0.52
1:A:281:ARG:HB3	1:A:282:GLY:HA3	1.92	0.52
1:B:129:HIS:HE1	1:B:183:LEU:HD23	1.75	0.52
1:A:222:TYR:HB3	1:A:230:LYS:NZ	2.25	0.52
1:A:222:TYR:HD1	1:A:232:GLU:HG2	1.76	0.51
1:D:16:VAL:HG12	1:D:79:THR:HB	1.92	0.51
1:A:297:THR:HG22	1:A:299:LYS:H	1.76	0.50
1:A:249:THR:HG23	1:A:258:ILE:HG12	1.94	0.49
1:B:29:PHE:HE2	1:B:76:LEU:HD13	1.78	0.49
1:A:293:ASP:O	1:A:297:THR:HB	2.13	0.49
1:A:222:TYR:HB3	1:A:230:LYS:HZ2	1.79	0.48
3:D:402:BLA:C3A	3:D:403:BLA:HMA2	2.44	0.48
1:A:185:ARG:HH22	3:A:403:BLA:HBA1	1.79	0.48
1:C:184:SER:O	1:C:188:ARG:HG3	2.14	0.47
3:C:403:BLA:HB	3:C:403:BLA:HA	1.62	0.47
1:D:112:LYS:HE2	1:D:116:GLN:OE1	2.14	0.47
1:B:150:VAL:HG12	1:B:229:LEU:HD13	1.96	0.47
1:D:97:HIS:CE1	1:D:123:LYS:HE2	2.49	0.47
3:C:402:BLA:CBD	3:C:403:BLA:HMA2	2.34	0.47
1:D:202:GLN:HE22	1:D:222:TYR:HE2	1.58	0.47
2:A:401:NAP:H72N	3:A:402:BLA:HB	1.62	0.47
3:C:403:BLA:HBB1	3:C:403:BLA:OB	2.14	0.47
1:D:146:LYS:HA	1:D:254:ARG:NH2	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:129:HIS:HE1	1:C:183:LEU:HD23	1.79	0.47
3:D:403:BLA:HBC1	3:D:403:BLA:CHD	2.45	0.47
1:A:63:TRP:CD2	1:A:87:ILE:HG12	2.49	0.47
1:D:220:THR:HG22	1:D:234:ILE:HD12	1.96	0.47
1:A:16:VAL:HG11	1:A:87:ILE:HD13	1.97	0.46
1:B:82:GLN:HB2	1:B:174:GLN:HA	1.96	0.46
1:C:28:VAL:HG21	1:C:285:ARG:HA	1.96	0.46
1:C:125:LEU:HD23	1:C:305:LEU:HD21	1.98	0.46
1:C:81:ASN:HB3	1:C:106:LEU:HD21	1.97	0.46
1:A:123:LYS:HB3	1:A:123:LYS:HE3	1.82	0.45
1:A:166:GLN:HG2	1:A:213:TYR:CG	2.52	0.45
1:B:40:PHE:HD2	1:B:57:VAL:HG12	1.81	0.45
1:B:15:ILE:O	1:B:40:PHE:HA	2.16	0.45
2:C:401:NAP:H72N	3:C:402:BLA:HC	1.65	0.45
1:A:267:ILE:HG12	1:A:272:GLU:HG3	1.99	0.45
1:D:21:ALA:O	1:D:25:ARG:HG2	2.17	0.45
1:A:162:ASN:HA	1:A:163:PRO:HA	1.81	0.44
1:D:16:VAL:HG13	1:D:41:TRP:HE3	1.82	0.44
3:B:403:BLA:HBB1	3:B:403:BLA:OB	2.17	0.44
1:C:177:PHE:CE2	1:C:179:LEU:HB2	2.53	0.44
3:A:403:BLA:HBB1	3:A:403:BLA:OB	2.18	0.43
1:C:171:HIS:HD2	1:C:213:TYR:CZ	2.36	0.43
1:D:81:ASN:OD1	1:D:103:PRO:HD2	2.18	0.43
3:B:403:BLA:CBC	3:B:403:BLA:HMC1	2.41	0.43
1:C:198:GLN:HB3	1:C:224:GLN:HB2	2.01	0.43
3:B:402:BLA:HBC1	3:B:402:BLA:CHD	2.43	0.43
1:C:171:HIS:CE1	1:C:173:GLN:HB2	2.54	0.43
1:D:131:GLU:HB2	1:D:188:ARG:NH1	2.34	0.43
1:D:104:LEU:HB2	1:D:127:VAL:HG21	2.00	0.43
1:B:101:GLU:HG3	1:B:128:GLU:OE1	2.19	0.42
1:C:69:ASP:OD2	1:C:71:GLU:HG2	2.19	0.42
1:D:16:VAL:HG13	1:D:41:TRP:CE3	2.54	0.42
1:B:188:ARG:HD3	1:B:248:PHE:HE1	1.84	0.42
1:B:63:TRP:CE3	1:B:87:ILE:HG12	2.53	0.42
1:A:266:LEU:HB3	1:A:273:THR:HB	2.01	0.42
1:B:282:GLY:O	1:B:286:GLN:HG3	2.19	0.42
1:C:147:ILE:C	1:C:254:ARG:HB2	2.39	0.42
1:D:153:ALA:O	1:D:231:ALA:HA	2.19	0.42
1:B:183:LEU:HD11	1:B:312:LEU:HA	2.00	0.42
3:B:402:BLA:HMA1	3:B:402:BLA:HBB	1.80	0.42
3:B:402:BLA:HMC1	3:B:402:BLA:HAC	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:PHE:HA	1:A:263:THR:O	2.20	0.42
1:A:15:ILE:O	1:A:40:PHE:HA	2.20	0.42
1:B:147:ILE:C	1:B:254:ARG:HB2	2.40	0.41
1:B:266:LEU:HB2	1:B:275:ILE:HD11	2.02	0.41
1:B:162:ASN:HA	1:B:163:PRO:HA	1.94	0.41
1:C:20:TYR:O	1:C:24:ARG:HG2	2.19	0.41
1:D:9:THR:HA	1:D:10:PRO:HD3	1.92	0.41
1:C:183:LEU:HD11	1:C:312:LEU:HA	2.02	0.41
3:A:402:BLA:HHD	3:A:402:BLA:HAC	1.91	0.41
1:C:153:ALA:O	1:C:231:ALA:HA	2.21	0.41
3:D:402:BLA:HMC3	3:D:403:BLA:CHD	2.51	0.41
1:B:113:LYS:HE2	1:B:113:LYS:HB3	1.86	0.41
1:D:169:THR:HA	1:D:175:PHE:CD1	2.56	0.41
3:A:402:BLA:OB	3:A:402:BLA:HBB1	2.21	0.41
1:C:230:LYS:HD2	1:C:230:LYS:H	1.82	0.40
1:A:21:ALA:O	1:A:25:ARG:HG2	2.21	0.40
1:B:246:ARG:HD3	3:B:402:BLA:O1D	2.21	0.40
1:C:16:VAL:HG13	1:C:41:TRP:HE3	1.87	0.40
1:D:295:LEU:HD23	1:D:295:LEU:HA	1.94	0.40
1:D:115:GLN:HG3	1:D:305:LEU:HD23	2.04	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	315/331 (95%)	307 (98%)	8 (2%)	0	100	100
1	B	314/331 (95%)	302 (96%)	12 (4%)	0	100	100
1	C	314/331 (95%)	307 (98%)	7 (2%)	0	100	100
1	D	314/331 (95%)	304 (97%)	10 (3%)	0	100	100
All	All	1257/1324 (95%)	1220 (97%)	37 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	255/267 (96%)	245 (96%)	10 (4%)	39	68
1	B	255/267 (96%)	247 (97%)	8 (3%)	47	76
1	C	255/267 (96%)	250 (98%)	5 (2%)	63	85
1	D	255/267 (96%)	251 (98%)	4 (2%)	70	89
All	All	1020/1068 (96%)	993 (97%)	27 (3%)	54	80

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	9	THR
1	B	30	ARG
1	B	50	LYS
1	B	100	LEU
1	B	101	GLU
1	B	230	LYS
1	B	248	PHE
1	B	274	GLU
1	A	57	VAL
1	A	100	LEU
1	A	128	GLU
1	A	167	ARG
1	A	198	GLN
1	A	248	PHE
1	A	274	GLU
1	A	280	ARG
1	A	283	LEU
1	A	297	THR
1	C	24	ARG
1	C	76	LEU
1	C	128	GLU
1	C	248	PHE

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Mol	Chain	Res	Type
1	C	324	TYR
1	D	9	THR
1	D	57	VAL
1	D	144	LEU
1	D	248	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	43	ASN
1	B	174	GLN
1	A	173	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](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

12 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	401	-	45,52,52	2.19	10 (22%)	55,80,80	4.33	18 (32%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BLA	A	402	-	36,46,46	5.53	20 (55%)	45,67,67	1.59	10 (22%)
3	BLA	A	403	-	36,46,46	5.55	22 (61%)	45,67,67	1.64	11 (24%)
2	NAP	B	401	-	45,52,52	1.89	8 (17%)	55,80,80	4.09	17 (30%)
3	BLA	B	402	-	36,46,46	5.91	23 (63%)	45,67,67	2.69	16 (35%)
3	BLA	B	403	-	36,46,46	5.67	21 (58%)	45,67,67	1.66	7 (15%)
2	NAP	C	401	-	45,52,52	2.23	9 (20%)	55,80,80	3.86	14 (25%)
3	BLA	C	402	-	36,46,46	5.55	21 (58%)	45,67,67	1.62	8 (17%)
3	BLA	C	403	-	36,46,46	5.51	22 (61%)	45,67,67	1.66	9 (20%)
2	NAP	D	401	-	45,52,52	1.89	8 (17%)	55,80,80	4.08	16 (29%)
3	BLA	D	402	-	36,46,46	5.50	20 (55%)	45,67,67	1.58	10 (22%)
3	BLA	D	403	-	36,46,46	5.53	23 (63%)	45,67,67	1.78	7 (15%)

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	401	-	-	0/27/67/67	0/5/5/5
3	BLA	A	402	-	-	0/22/74/74	0/4/4/4
3	BLA	A	403	-	-	0/22/74/74	0/4/4/4
2	NAP	B	401	-	-	0/27/67/67	0/5/5/5
3	BLA	B	402	-	-	0/22/74/74	0/4/4/4
3	BLA	B	403	-	-	0/22/74/74	0/4/4/4
2	NAP	C	401	-	-	0/27/67/67	0/5/5/5
3	BLA	C	402	-	-	0/22/74/74	0/4/4/4
3	BLA	C	403	-	-	0/22/74/74	0/4/4/4
2	NAP	D	401	-	-	0/27/67/67	0/5/5/5
3	BLA	D	402	-	-	0/22/74/74	0/4/4/4
3	BLA	D	403	-	-	0/22/74/74	0/4/4/4

All (207) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	403	BLA	C3C-C4C	-8.14	1.32	1.45
3	D	402	BLA	C3C-C4C	-7.81	1.33	1.45
3	A	402	BLA	C3C-C4C	-7.68	1.33	1.45
3	C	403	BLA	C3C-C4C	-7.56	1.33	1.45
3	B	403	BLA	C3C-C4C	-7.46	1.33	1.45
3	C	402	BLA	C3C-C4C	-7.35	1.33	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	NAP	C3N-C7N	-7.28	1.39	1.50
2	A	401	NAP	C3N-C7N	-7.13	1.39	1.50
3	C	402	BLA	C4D-C3D	-7.09	1.33	1.45
2	D	401	NAP	C3N-C7N	-7.05	1.39	1.50
3	D	403	BLA	C4D-C3D	-7.01	1.33	1.45
2	C	401	NAP	C3N-C7N	-6.97	1.39	1.50
3	D	402	BLA	C4D-C3D	-6.88	1.34	1.45
3	A	403	BLA	C4D-C3D	-6.83	1.34	1.45
3	C	403	BLA	C1B-C2B	-6.78	1.32	1.45
3	C	403	BLA	C4D-C3D	-6.77	1.34	1.45
3	B	402	BLA	C1B-C2B	-6.76	1.32	1.45
3	A	402	BLA	C4D-C3D	-6.70	1.34	1.45
3	A	403	BLA	C1B-C2B	-6.63	1.33	1.45
3	B	403	BLA	C4D-C3D	-6.60	1.34	1.45
3	B	403	BLA	C1B-C2B	-6.54	1.33	1.45
3	D	402	BLA	C1B-C2B	-6.46	1.33	1.45
3	D	403	BLA	C3C-C4C	-6.46	1.35	1.45
3	C	402	BLA	C1B-C2B	-6.42	1.33	1.45
3	B	402	BLA	C1D-C2D	-6.37	1.31	1.45
3	D	403	BLA	C1B-C2B	-6.36	1.33	1.45
3	A	402	BLA	C1B-C2B	-6.33	1.33	1.45
3	B	402	BLA	C3C-C4C	-6.12	1.35	1.45
3	A	402	BLA	C1D-C2D	-6.00	1.32	1.45
3	B	403	BLA	C1D-C2D	-5.96	1.32	1.45
3	A	403	BLA	C1D-C2D	-5.95	1.32	1.45
3	D	402	BLA	C1D-C2D	-5.92	1.32	1.45
3	D	403	BLA	C1D-C2D	-5.90	1.32	1.45
3	C	403	BLA	C1D-C2D	-5.80	1.33	1.45
3	C	402	BLA	C1D-C2D	-5.80	1.33	1.45
3	B	402	BLA	C4D-C3D	-5.60	1.36	1.45
3	D	403	BLA	C1C-C2C	-5.09	1.32	1.47
3	B	402	BLA	C1C-C2C	-5.00	1.33	1.47
3	D	402	BLA	C1C-C2C	-4.93	1.33	1.47
3	C	402	BLA	C1C-C2C	-4.92	1.33	1.47
3	A	402	BLA	C1C-C2C	-4.88	1.33	1.47
3	B	403	BLA	C1C-C2C	-4.86	1.33	1.47
3	C	403	BLA	C1C-C2C	-4.80	1.33	1.47
3	A	403	BLA	C1C-C2C	-4.68	1.34	1.47
3	A	403	BLA	CAB-C3B	-4.56	1.34	1.47
3	C	403	BLA	CAB-C3B	-4.56	1.34	1.47
3	C	402	BLA	C3B-C4B	-4.51	1.33	1.47
3	B	403	BLA	CAB-C3B	-4.49	1.34	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	403	BLA	C3B-C4B	-4.43	1.33	1.47
3	B	402	BLA	C3B-C4B	-4.41	1.33	1.47
3	D	402	BLA	CAB-C3B	-4.39	1.34	1.47
3	A	403	BLA	CAC-C3C	-4.38	1.34	1.47
3	D	403	BLA	CAC-C3C	-4.37	1.34	1.47
3	A	402	BLA	CAB-C3B	-4.35	1.34	1.47
3	B	403	BLA	C3B-C4B	-4.32	1.34	1.47
3	B	402	BLA	CAB-C3B	-4.27	1.35	1.47
3	B	403	BLA	CAC-C3C	-4.26	1.35	1.47
3	A	403	BLA	C3B-C4B	-4.25	1.34	1.47
3	C	402	BLA	CAB-C3B	-4.21	1.35	1.47
3	D	402	BLA	C3B-C4B	-4.21	1.34	1.47
3	A	402	BLA	C3B-C4B	-4.21	1.34	1.47
3	C	403	BLA	CAC-C3C	-4.20	1.35	1.47
3	C	403	BLA	C3B-C4B	-4.20	1.34	1.47
3	B	402	BLA	CAC-C3C	-4.20	1.35	1.47
3	C	402	BLA	CAC-C3C	-4.18	1.35	1.47
3	D	402	BLA	CAC-C3C	-4.18	1.35	1.47
3	D	403	BLA	CAB-C3B	-4.13	1.35	1.47
3	A	402	BLA	CAC-C3C	-4.11	1.35	1.47
2	B	401	NAP	C5A-C4A	-2.92	1.33	1.40
2	D	401	NAP	C5A-C4A	-2.83	1.34	1.40
2	C	401	NAP	C5A-C4A	-2.69	1.34	1.40
3	B	402	BLA	C1B-NB	-2.63	1.33	1.37
3	C	402	BLA	C1B-NB	-2.60	1.33	1.37
2	B	401	NAP	C2D-C1D	-2.59	1.49	1.53
3	B	403	BLA	C4C-NC	-2.59	1.33	1.37
2	A	401	NAP	C5A-C4A	-2.57	1.34	1.40
3	D	403	BLA	C1B-NB	-2.55	1.33	1.37
3	C	402	BLA	C4C-NC	-2.52	1.33	1.37
3	C	403	BLA	C1B-NB	-2.52	1.33	1.37
3	D	402	BLA	C1B-NB	-2.52	1.33	1.37
3	A	403	BLA	C4C-NC	-2.52	1.33	1.37
3	C	403	BLA	C4C-NC	-2.51	1.33	1.37
3	A	402	BLA	C1B-NB	-2.50	1.33	1.37
3	D	402	BLA	C4C-NC	-2.40	1.33	1.37
3	D	403	BLA	C4C-NC	-2.37	1.33	1.37
3	A	402	BLA	C4C-NC	-2.37	1.33	1.37
3	A	403	BLA	C1B-NB	-2.37	1.33	1.37
3	B	403	BLA	C1B-NB	-2.36	1.33	1.37
3	B	402	BLA	C4C-NC	-2.33	1.33	1.37
3	A	403	BLA	C3B-C2B	-2.33	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	403	BLA	C3B-C2B	-2.33	1.32	1.37
2	D	401	NAP	C3B-C2B	-2.24	1.48	1.53
3	B	403	BLA	C3B-C2B	-2.23	1.32	1.37
3	B	402	BLA	C3C-C2C	-2.16	1.33	1.37
2	B	401	NAP	P2B-O2X	-2.15	1.47	1.54
2	C	401	NAP	P2B-O2X	-2.12	1.47	1.54
2	A	401	NAP	C2D-C3D	-2.12	1.47	1.53
2	D	401	NAP	P2B-O2X	-2.12	1.47	1.54
3	C	402	BLA	C1C-NC	-2.11	1.33	1.37
3	D	403	BLA	C3C-C2C	-2.09	1.33	1.37
2	A	401	NAP	P2B-O2X	-2.09	1.47	1.54
3	D	403	BLA	C4D-ND	-2.08	1.33	1.38
3	B	402	BLA	C3D-C2D	-2.06	1.32	1.36
3	D	403	BLA	C1C-NC	-2.05	1.33	1.37
3	A	403	BLA	C3D-C2D	-2.04	1.32	1.36
3	C	403	BLA	C3D-C2D	-2.02	1.32	1.36
2	D	401	NAP	C2D-C1D	-2.01	1.50	1.53
2	A	401	NAP	O4D-C4D	2.05	1.49	1.45
2	C	401	NAP	O4D-C4D	2.10	1.49	1.45
2	B	401	NAP	O4D-C4D	2.14	1.49	1.45
3	B	402	BLA	CAA-C2A	3.28	1.57	1.52
2	C	401	NAP	C2A-N1A	3.30	1.40	1.33
2	B	401	NAP	C2A-N1A	3.30	1.40	1.33
2	B	401	NAP	C2N-N1N	3.33	1.39	1.35
2	D	401	NAP	C2A-N1A	3.36	1.40	1.33
2	A	401	NAP	C2A-N1A	3.39	1.40	1.33
3	B	402	BLA	C4A-CHB	3.40	1.53	1.40
2	D	401	NAP	C2N-N1N	3.52	1.40	1.35
3	C	403	BLA	C4A-CHB	3.62	1.54	1.40
3	A	403	BLA	C4A-CHB	3.65	1.54	1.40
2	A	401	NAP	C2N-N1N	3.72	1.40	1.35
3	C	402	BLA	C1A-CHA	3.77	1.54	1.40
3	D	403	BLA	C4A-CHB	3.78	1.54	1.40
3	C	402	BLA	C4A-CHB	3.80	1.54	1.40
3	C	403	BLA	C1A-CHA	3.80	1.54	1.40
3	B	403	BLA	C4A-CHB	3.81	1.54	1.40
3	A	403	BLA	C1A-CHA	3.83	1.54	1.40
3	D	402	BLA	C4A-CHB	3.89	1.55	1.40
2	C	401	NAP	C2N-N1N	3.91	1.40	1.35
3	A	402	BLA	C4A-CHB	3.92	1.55	1.40
2	A	401	NAP	O4B-C1B	3.92	1.46	1.41
3	D	402	BLA	C1A-CHA	3.92	1.55	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	BLA	C1A-CHA	3.95	1.55	1.40
3	B	403	BLA	C1A-CHA	4.03	1.55	1.40
3	D	403	BLA	C1A-CHA	4.07	1.55	1.40
3	B	402	BLA	C1A-CHA	4.20	1.56	1.40
3	A	403	BLA	CBC-CAC	4.55	1.52	1.30
3	D	403	BLA	CBC-CAC	4.73	1.53	1.30
3	A	403	BLA	CBB-CAB	4.74	1.53	1.30
3	C	403	BLA	CBB-CAB	4.75	1.53	1.30
3	B	402	BLA	CBC-CAC	4.77	1.53	1.30
3	B	403	BLA	CBB-CAB	4.78	1.53	1.30
3	B	403	BLA	CBC-CAC	4.80	1.53	1.30
3	D	402	BLA	CBB-CAB	4.81	1.53	1.30
3	D	402	BLA	CBC-CAC	4.82	1.53	1.30
3	C	403	BLA	CBC-CAC	4.82	1.53	1.30
3	C	402	BLA	CBB-CAB	4.84	1.53	1.30
3	A	402	BLA	CBC-CAC	4.84	1.54	1.30
3	B	402	BLA	CBB-CAB	4.85	1.54	1.30
3	A	402	BLA	CBB-CAB	4.89	1.54	1.30
3	C	402	BLA	CBC-CAC	4.89	1.54	1.30
2	B	401	NAP	C2A-N3A	4.90	1.40	1.32
3	D	403	BLA	CBB-CAB	4.90	1.54	1.30
2	C	401	NAP	C2A-N3A	5.01	1.41	1.32
2	D	401	NAP	C2A-N3A	5.03	1.41	1.32
3	A	403	BLA	CHD-C1D	5.14	1.53	1.40
2	A	401	NAP	C2A-N3A	5.22	1.41	1.32
3	D	402	BLA	CHD-C1D	5.23	1.53	1.40
2	C	401	NAP	O4B-C1B	5.32	1.48	1.41
3	A	402	BLA	CHD-C1D	5.43	1.53	1.40
3	D	403	BLA	CHD-C1D	5.55	1.54	1.40
3	C	403	BLA	CHD-C1D	5.74	1.54	1.40
3	B	403	BLA	CHD-C1D	5.79	1.54	1.40
3	C	402	BLA	CHD-C1D	5.81	1.54	1.40
2	A	401	NAP	C3B-C2B	5.92	1.66	1.53
3	B	402	BLA	CHD-C1D	5.93	1.55	1.40
2	C	401	NAP	C3B-C2B	5.94	1.66	1.53
3	A	402	BLA	CHD-C4C	6.46	1.53	1.37
3	D	402	BLA	CHD-C4C	6.57	1.53	1.37
3	C	402	BLA	OC-C1C	6.65	1.36	1.23
3	A	403	BLA	CHD-C4C	6.70	1.53	1.37
3	D	402	BLA	OC-C1C	6.71	1.36	1.23
3	D	403	BLA	OC-C1C	6.72	1.36	1.23
3	B	402	BLA	OC-C1C	6.73	1.36	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	BLA	OB-C4B	6.73	1.36	1.23
3	B	403	BLA	OC-C1C	6.75	1.36	1.23
3	C	402	BLA	OB-C4B	6.78	1.36	1.23
3	A	402	BLA	OC-C1C	6.79	1.36	1.23
3	C	403	BLA	OC-C1C	6.80	1.36	1.23
3	A	402	BLA	OB-C4B	6.81	1.36	1.23
3	B	403	BLA	OB-C4B	6.81	1.36	1.23
3	D	403	BLA	OB-C4B	6.83	1.36	1.23
3	D	402	BLA	OB-C4B	6.83	1.36	1.23
3	A	403	BLA	OC-C1C	6.84	1.36	1.23
3	C	403	BLA	OB-C4B	6.88	1.36	1.23
3	C	403	BLA	CHD-C4C	6.90	1.54	1.37
3	A	403	BLA	OB-C4B	6.91	1.36	1.23
3	B	403	BLA	CHD-C4C	6.93	1.54	1.37
3	C	402	BLA	CHD-C4C	6.94	1.54	1.37
3	B	402	BLA	CHD-C4C	7.00	1.54	1.37
3	D	403	BLA	CHD-C4C	7.06	1.54	1.37
3	B	402	BLA	CHB-C1B	8.78	1.52	1.34
3	C	403	BLA	CHB-C1B	9.09	1.52	1.34
3	A	403	BLA	CHB-C1B	9.66	1.53	1.34
3	D	403	BLA	CHB-C1B	9.83	1.54	1.34
3	C	402	BLA	CHB-C1B	9.88	1.54	1.34
3	D	402	BLA	CHB-C1B	9.88	1.54	1.34
3	B	403	BLA	CHB-C1B	9.91	1.54	1.34
3	A	402	BLA	CHB-C1B	10.08	1.54	1.34
3	D	402	BLA	CHA-C4D	21.11	1.54	1.35
3	A	402	BLA	CHA-C4D	21.36	1.54	1.35
3	C	403	BLA	CHA-C4D	21.38	1.54	1.35
3	A	403	BLA	CHA-C4D	21.47	1.54	1.35
3	C	402	BLA	CHA-C4D	21.48	1.54	1.35
3	D	403	BLA	CHA-C4D	21.55	1.54	1.35
3	B	403	BLA	CHA-C4D	22.51	1.55	1.35
3	B	402	BLA	CHA-C4D	25.29	1.58	1.35

All (143) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	NAP	O4B-C1B-C2B	-17.71	74.76	106.60
2	D	401	NAP	O4B-C1B-C2B	-17.26	75.58	106.60
2	D	401	NAP	N3A-C2A-N1A	-13.38	118.36	128.87
2	B	401	NAP	N3A-C2A-N1A	-13.33	118.40	128.87
2	C	401	NAP	N3A-C2A-N1A	-13.16	118.53	128.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NAP	N3A-C2A-N1A	-12.56	119.00	128.87
2	B	401	NAP	O4B-C4B-C5B	-8.98	77.17	109.29
2	C	401	NAP	C3B-C2B-C1B	-8.91	85.57	102.63
2	D	401	NAP	O4B-C4B-C5B	-8.89	77.49	109.29
2	C	401	NAP	C5B-C4B-C3B	-8.88	80.84	115.20
2	A	401	NAP	C3B-C2B-C1B	-8.80	85.77	102.63
2	A	401	NAP	O4D-C4D-C5D	-8.61	78.46	109.29
2	B	401	NAP	O4D-C4D-C5D	-8.57	78.63	109.29
2	D	401	NAP	O4D-C4D-C5D	-8.54	78.73	109.29
2	C	401	NAP	O4D-C4D-C5D	-8.52	78.80	109.29
2	C	401	NAP	O4D-C4D-C3D	-8.22	88.48	105.16
2	D	401	NAP	O4D-C4D-C3D	-7.80	89.34	105.16
2	A	401	NAP	C5B-C4B-C3B	-7.75	85.21	115.20
2	B	401	NAP	O4D-C4D-C3D	-7.72	89.51	105.16
2	A	401	NAP	O4D-C4D-C3D	-7.71	89.51	105.16
2	D	401	NAP	O4B-C4B-C3B	-7.27	90.42	105.16
3	D	403	BLA	C4C-CHD-C1D	-6.84	110.79	128.03
2	B	401	NAP	O4B-C4B-C3B	-6.73	91.50	105.16
3	B	402	BLA	CAD-C3D-C2D	-6.37	116.32	127.96
3	C	402	BLA	C4C-CHD-C1D	-6.11	112.62	128.03
2	A	401	NAP	O4B-C4B-C3B	-5.79	93.41	105.16
2	C	401	NAP	O4B-C4B-C3B	-5.66	93.68	105.16
2	D	401	NAP	C3B-C2B-C1B	-5.56	91.98	102.63
2	B	401	NAP	C3B-C2B-C1B	-5.10	92.86	102.63
3	B	403	BLA	C4C-CHD-C1D	-5.01	115.39	128.03
3	A	402	BLA	C4C-CHD-C1D	-4.63	116.36	128.03
3	D	402	BLA	C4C-CHD-C1D	-4.55	116.56	128.03
3	C	403	BLA	C4C-CHD-C1D	-4.11	117.66	128.03
3	A	403	BLA	C4C-CHD-C1D	-3.86	118.31	128.03
3	B	402	BLA	C1B-NB-C4B	-3.78	105.25	110.73
3	D	402	BLA	C4C-NC-C1C	-3.54	105.61	110.73
3	A	402	BLA	C4C-NC-C1C	-3.53	105.62	110.73
3	A	403	BLA	C4C-NC-C1C	-3.49	105.67	110.73
2	A	401	NAP	O4B-C1B-N9A	-3.39	101.69	108.11
3	B	403	BLA	CAD-C3D-C2D	-3.29	121.95	127.96
2	B	401	NAP	C4B-O4B-C1B	-3.26	106.19	109.64
3	B	402	BLA	CAC-C3C-C2C	-3.26	119.61	128.90
3	C	403	BLA	C4C-NC-C1C	-3.25	106.03	110.73
3	A	403	BLA	CHD-C4C-C3C	-3.20	119.61	128.02
3	B	403	BLA	C4C-NC-C1C	-3.11	106.23	110.73
3	D	403	BLA	CAC-C3C-C2C	-3.08	120.12	128.90
2	D	401	NAP	C4B-O4B-C1B	-3.06	106.40	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	BLA	CHD-C1D-C2D	-3.03	117.45	124.71
3	B	402	BLA	C4C-CHD-C1D	-3.03	120.38	128.03
3	C	402	BLA	C1B-NB-C4B	-3.02	106.36	110.73
3	C	403	BLA	CAD-C3D-C2D	-2.93	122.61	127.96
3	B	402	BLA	CHB-C1B-C2B	-2.93	121.28	126.96
3	D	403	BLA	C1B-NB-C4B	-2.91	106.52	110.73
3	C	402	BLA	C4C-NC-C1C	-2.88	106.56	110.73
3	C	403	BLA	C1B-NB-C4B	-2.81	106.66	110.73
2	A	401	NAP	C3N-C7N-N7N	-2.74	114.73	117.82
2	B	401	NAP	C3N-C7N-N7N	-2.72	114.75	117.82
2	D	401	NAP	C3N-C7N-N7N	-2.65	114.82	117.82
2	C	401	NAP	O2B-P2B-O1X	-2.65	101.16	107.48
3	D	402	BLA	CHD-C4C-C3C	-2.63	121.10	128.02
2	B	401	NAP	O2B-P2B-O1X	-2.62	101.22	107.48
2	D	401	NAP	C2D-C1D-N1N	-2.61	108.41	113.53
3	B	402	BLA	C4C-NC-C1C	-2.60	106.97	110.73
2	C	401	NAP	C2D-C1D-N1N	-2.56	108.52	113.53
3	B	403	BLA	C1B-NB-C4B	-2.53	107.06	110.73
2	D	401	NAP	C4D-O4D-C1D	-2.51	106.98	109.64
2	B	401	NAP	C2B-C3B-C4B	-2.49	95.95	101.85
3	A	403	BLA	C1B-NB-C4B	-2.47	107.15	110.73
2	A	401	NAP	C4D-O4D-C1D	-2.43	107.06	109.64
3	A	402	BLA	CHD-C4C-C3C	-2.43	121.64	128.02
3	D	403	BLA	CAD-C3D-C2D	-2.40	123.59	127.96
2	C	401	NAP	C3N-C7N-N7N	-2.38	115.14	117.82
3	D	402	BLA	C1B-NB-C4B	-2.37	107.30	110.73
2	B	401	NAP	C2D-C1D-N1N	-2.37	108.89	113.53
3	D	403	BLA	CHD-C4C-NC	-2.31	121.18	126.15
2	A	401	NAP	C2D-C1D-N1N	-2.29	109.04	113.53
3	C	402	BLA	CAD-C3D-C2D	-2.28	123.79	127.96
3	A	403	BLA	CAD-C3D-C2D	-2.26	123.83	127.96
3	B	402	BLA	CMB-C2B-C1B	-2.26	121.25	124.20
3	A	402	BLA	CAD-C3D-C2D	-2.22	123.92	127.96
3	C	402	BLA	CMB-C2B-C1B	-2.21	121.32	124.20
3	A	402	BLA	C1B-NB-C4B	-2.20	107.54	110.73
3	C	402	BLA	CHB-C1B-C2B	-2.19	122.71	126.96
3	A	403	BLA	CAB-C3B-C2B	-2.17	122.72	128.90
2	B	401	NAP	C4D-O4D-C1D	-2.15	107.36	109.64
3	C	403	BLA	CHD-C4C-C3C	-2.14	122.39	128.02
3	A	403	BLA	CAC-C3C-C4C	-2.12	118.42	123.67
3	D	402	BLA	OB-C4B-NB	-2.12	120.09	125.21
3	D	402	BLA	CAD-C3D-C2D	-2.11	124.10	127.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	NAP	O4B-C1B-N9A	-2.10	104.15	108.11
3	A	402	BLA	OB-C4B-NB	-2.09	120.17	125.21
3	A	403	BLA	CAB-C3B-C4B	2.01	129.35	123.58
3	D	403	BLA	CBA-CAA-C2A	2.01	116.02	112.49
3	C	402	BLA	C3C-C4C-NC	2.04	109.97	106.85
3	D	402	BLA	CAB-C3B-C4B	2.04	129.45	123.58
2	C	401	NAP	N6A-C6A-N1A	2.05	121.96	118.52
3	B	403	BLA	C3C-C4C-NC	2.06	110.02	106.85
3	A	403	BLA	CAD-C3D-C4D	2.08	128.64	125.00
2	B	401	NAP	C2N-C3N-C4N	2.10	120.65	118.27
3	C	403	BLA	C1D-C2D-C3D	2.10	108.86	106.43
3	A	402	BLA	CAD-C3D-C4D	2.13	128.73	125.00
3	D	402	BLA	CBA-CAA-C2A	2.15	116.27	112.49
3	B	403	BLA	CHD-C1D-ND	2.19	128.98	124.89
2	D	401	NAP	N6A-C6A-N1A	2.19	122.19	118.52
3	C	403	BLA	C3C-C4C-NC	2.21	110.23	106.85
2	A	401	NAP	O2B-C2B-C1B	2.24	118.89	110.12
3	C	402	BLA	CBA-CAA-C2A	2.26	116.47	112.49
3	A	403	BLA	C3C-C4C-NC	2.28	110.35	106.85
3	D	402	BLA	CHD-C1D-ND	2.29	129.17	124.89
3	A	402	BLA	CBA-CAA-C2A	2.37	116.66	112.49
2	A	401	NAP	N6A-C6A-N1A	2.37	122.49	118.52
2	A	401	NAP	C4B-O4B-C1B	2.41	112.20	109.64
2	A	401	NAP	O5B-C5B-C4B	2.43	117.87	109.09
3	D	402	BLA	C3C-C4C-NC	2.45	110.61	106.85
3	A	402	BLA	C3C-C4C-NC	2.49	110.66	106.85
2	D	401	NAP	P2B-O2B-C2B	2.52	128.01	121.56
2	D	401	NAP	C1B-N9A-C4A	2.56	129.66	126.81
3	B	402	BLA	C2B-C1B-NB	2.71	110.77	106.94
3	A	402	BLA	CHD-C1D-ND	2.72	129.97	124.89
3	C	403	BLA	CAD-C3D-C4D	2.80	129.92	125.00
2	B	401	NAP	P2B-O2B-C2B	2.82	128.78	121.56
3	B	402	BLA	C1D-C2D-C3D	2.82	109.69	106.43
3	C	403	BLA	CHD-C1D-ND	2.90	130.30	124.89
3	A	403	BLA	CHD-C1D-ND	2.90	130.31	124.89
3	B	403	BLA	CAD-C3D-C4D	2.98	130.22	125.00
3	B	402	BLA	CMA-C3A-C2A	3.24	132.01	125.24
3	D	403	BLA	CAC-C3C-C4C	3.87	133.25	123.67
3	B	402	BLA	CAC-C3C-C4C	4.07	133.75	123.67
2	A	401	NAP	P2B-O2B-C2B	4.62	133.40	121.56
3	B	402	BLA	CHD-C1D-ND	4.90	134.03	124.89
2	B	401	NAP	C5B-C4B-C3B	4.91	134.20	115.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	NAP	C5D-C4D-C3D	5.07	134.81	115.20
2	B	401	NAP	C5D-C4D-C3D	5.17	135.22	115.20
2	D	401	NAP	C5B-C4B-C3B	5.21	135.37	115.20
2	D	401	NAP	C5D-C4D-C3D	5.26	135.56	115.20
2	A	401	NAP	C5D-C4D-C3D	5.32	135.80	115.20
3	B	402	BLA	CAA-CBA-CGA	5.85	124.17	112.78
3	B	402	BLA	CAD-C3D-C4D	6.18	135.84	125.00
3	B	402	BLA	CBA-CAA-C2A	7.24	125.22	112.49
2	C	401	NAP	C2B-C1B-N9A	8.43	137.57	113.48
2	A	401	NAP	C2B-C1B-N9A	9.48	140.56	113.48
2	C	401	NAP	C1B-N9A-C4A	12.31	140.54	126.81
2	A	401	NAP	C1B-N9A-C4A	18.56	147.51	126.81

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 39 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAP	3	0
3	A	402	BLA	4	0
3	A	403	BLA	3	0
2	B	401	NAP	1	0
3	B	402	BLA	8	0
3	B	403	BLA	4	0
2	C	401	NAP	3	0
3	C	402	BLA	5	0
3	C	403	BLA	6	0
2	D	401	NAP	1	0
3	D	402	BLA	4	0
3	D	403	BLA	5	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, 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	317/331 (95%)	0.14	4 (1%) 79 75	15, 22, 36, 53	0
1	B	316/331 (95%)	0.24	8 (2%) 61 54	17, 24, 38, 79	0
1	C	316/331 (95%)	0.28	4 (1%) 79 75	20, 29, 41, 80	0
1	D	316/331 (95%)	0.37	8 (2%) 61 54	20, 27, 43, 59	0
All	All	1265/1324 (95%)	0.26	24 (1%) 70 64	15, 26, 41, 80	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	280	ARG	7.3
1	C	279	SER	4.3
1	B	280	ARG	4.2
1	D	271	THR	3.5
1	B	279	SER	3.3
1	D	9	THR	3.0
1	D	270	GLN	3.0
1	B	281	ARG	2.8
1	B	9	THR	2.8
1	D	246	ARG	2.8
1	A	9	THR	2.7
1	C	281	ARG	2.6
1	C	53	ASP	2.5
1	B	20	TYR	2.4
1	A	71	GLU	2.4
1	B	293	ASP	2.3
1	A	279	SER	2.2
1	B	142	GLN	2.2
1	B	324	TYR	2.1
1	D	67	ILE	2.1
1	D	116	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	12	ARG	2.0
1	A	61	GLN	2.0
1	D	70	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	BLA	B	402	43/43	0.59	0.54	3.86	35,55,75,81	0
3	BLA	A	403	43/43	0.88	0.21	1.79	19,25,31,34	0
3	BLA	C	403	43/43	0.62	0.49	1.76	33,63,80,83	0
3	BLA	C	402	43/43	0.83	0.23	1.30	31,47,61,65	0
2	NAP	A	401	48/48	0.91	0.18	1.18	18,21,27,29	0
3	BLA	B	403	43/43	0.87	0.21	0.80	25,40,53,59	0
2	NAP	C	401	48/48	0.90	0.20	0.72	29,31,36,40	0
3	BLA	D	403	43/43	0.88	0.21	0.61	23,29,32,38	0
3	BLA	A	402	43/43	0.92	0.17	0.37	17,21,24,26	0
3	BLA	D	402	43/43	0.92	0.17	0.30	22,25,27,29	0
2	NAP	B	401	48/48	0.94	0.16	-0.21	22,25,30,34	0
2	NAP	D	401	48/48	0.95	0.15	-0.26	22,23,29,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.