



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 21, 2018 – 11:57 AM EST

PDB ID : 6BWC  
Title : X-ray structure of Pen from Bacillus thuringiensis  
Authors : Delvaux, N.A.; Thoden, J.B.; Holden, H.M.  
Deposited on : 2017-12-14  
Resolution : 2.70 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030736  
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-20030736

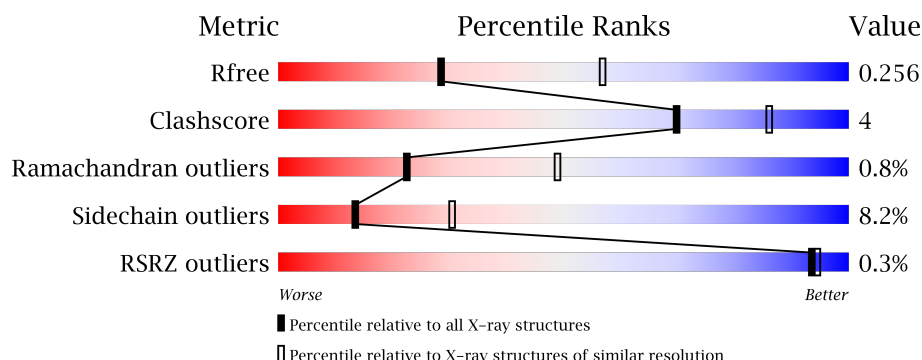
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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	343	
1	B	343	
1	C	343	
1	D	343	
1	E	343	

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Mol	Chain	Length	Quality of chain
1	F	343	<div> <div></div> <div>79%</div> <div>13%</div> <div>6%</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16002 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 Polysaccharide biosynthesis protein CapD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2559	1631	427	486	15			
1	B	316	Total	C	N	O	S	0	0	0
			2504	1595	420	474	15			
1	C	327	Total	C	N	O	S	0	0	0
			2590	1650	436	489	15			
1	D	324	Total	C	N	O	S	0	0	0
			2566	1636	431	484	15			
1	E	322	Total	C	N	O	S	0	0	0
			2553	1625	429	484	15			
1	F	323	Total	C	N	O	S	0	0	0
			2549	1624	425	485	15			

There are 24 discrepancies between the modelled and reference sequences:

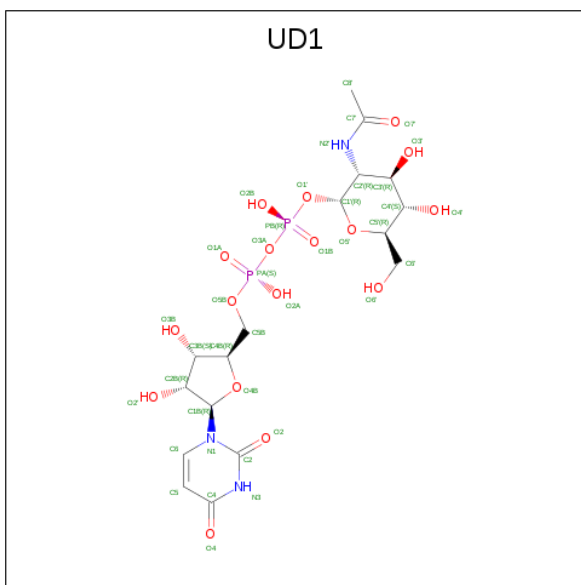
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP J3UJH9
A	0	HIS	-	expression tag	UNP J3UJH9
A	128	ASN	ASP	engineered mutation	UNP J3UJH9
A	129	ALA	LYS	engineered mutation	UNP J3UJH9
B	-1	GLY	-	expression tag	UNP J3UJH9
B	0	HIS	-	expression tag	UNP J3UJH9
B	128	ASN	ASP	engineered mutation	UNP J3UJH9
B	129	ALA	LYS	engineered mutation	UNP J3UJH9
C	-1	GLY	-	expression tag	UNP J3UJH9
C	0	HIS	-	expression tag	UNP J3UJH9
C	128	ASN	ASP	engineered mutation	UNP J3UJH9
C	129	ALA	LYS	engineered mutation	UNP J3UJH9
D	-1	GLY	-	expression tag	UNP J3UJH9
D	0	HIS	-	expression tag	UNP J3UJH9
D	128	ASN	ASP	engineered mutation	UNP J3UJH9
D	129	ALA	LYS	engineered mutation	UNP J3UJH9
E	-1	GLY	-	expression tag	UNP J3UJH9

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Chain	Residue	Modelled	Actual	Comment	Reference
E	0	HIS	-	expression tag	UNP J3UJH9
E	128	ASN	ASP	engineered mutation	UNP J3UJH9
E	129	ALA	LYS	engineered mutation	UNP J3UJH9
F	-1	GLY	-	expression tag	UNP J3UJH9
F	0	HIS	-	expression tag	UNP J3UJH9
F	128	ASN	ASP	engineered mutation	UNP J3UJH9
F	129	ALA	LYS	engineered mutation	UNP J3UJH9

- # NAP
- 
- The chemical structure of Naproxen (NAP) is shown, illustrating its enantiomers and stereochemistry. The molecule consists of a naphthalene ring system substituted with a carboxylic acid group and a chiral center. The stereochemistry is indicated by wedged and dashed bonds, showing the spatial arrangement of the atoms. The structure is labeled with various atoms and groups, including NH<sub>2</sub>, NH, COOH, and CO<sub>2</sub>H, and is associated with the chemical formula C<sub>15</sub>H<sub>11</sub>NO<sub>2</sub>.

- Molecule 3 is URIDINE-DIPHOSPHATE-N-ACETYLGUCOSAMINE (three-letter code: UD1) (formula:  $C_{17}H_{27}N_3O_{17}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 39	C 17	N 3	O 17	P 2	0	0
3	B	1	Total 39	C 17	N 3	O 17	P 2	0	0
3	C	1	Total 39	C 17	N 3	O 17	P 2	0	0
3	D	1	Total 39	C 17	N 3	O 17	P 2	0	0
3	E	1	Total 39	C 17	N 3	O 17	P 2	0	0
3	F	1	Total 39	C 17	N 3	O 17	P 2	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		

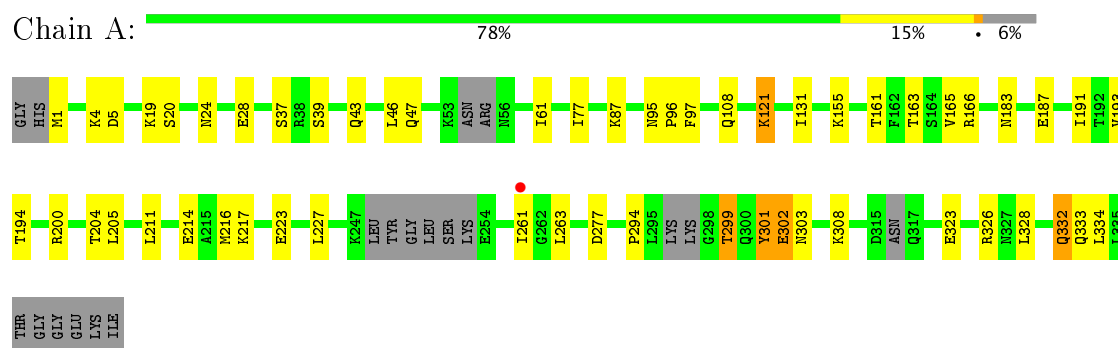
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	34	Total	O	0	0
			34	34		
5	B	21	Total	O	0	0
			21	21		
5	C	25	Total	O	0	0
			25	25		
5	D	21	Total	O	0	0
			21	21		
5	E	18	Total	O	0	0
			18	18		
5	F	15	Total	O	0	0
			15	15		

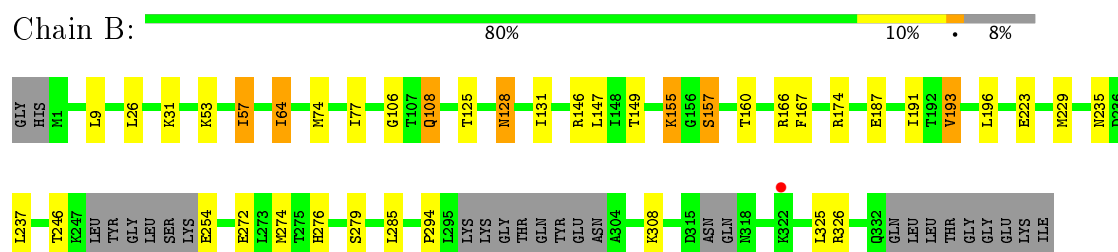
### 3 Residue-property plots [i](#)

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.

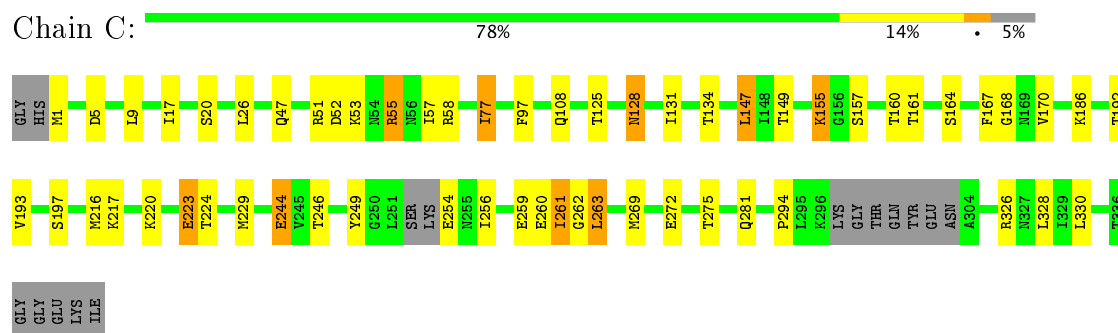
#### • Molecule 1: Polysaccharide biosynthesis protein CapD



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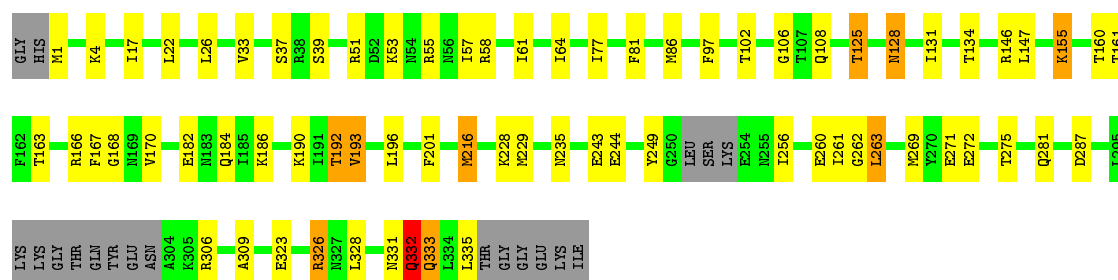
#### • Molecule 1: Polysaccharide biosynthesis protein CapD



#### • Molecule 1: Polysaccharide biosynthesis protein CapD

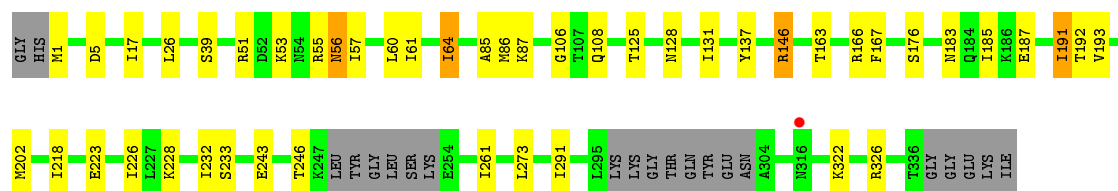






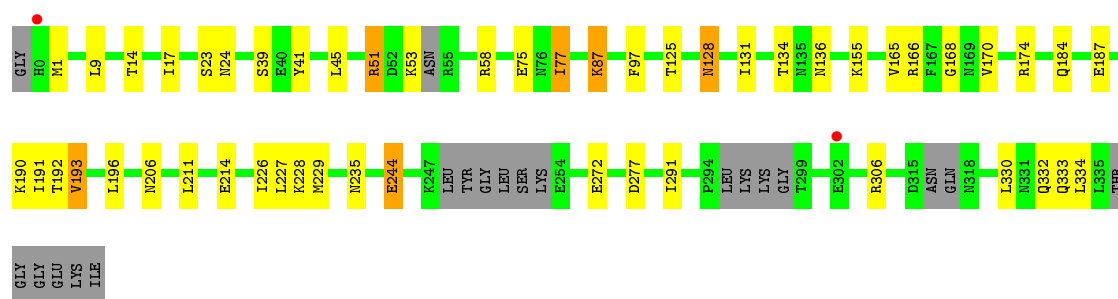
- Molecule 1: Polysaccharide biosynthesis protein CapD

Chain E: 80% 13% 6%



- Molecule 1: Polysaccharide biosynthesis protein CapD

Chain F: 79% 13% 6%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.52Å 141.88Å 93.24Å 90.00° 111.66° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 29.82 – 2.70	Depositor EDS
% Data completeness (in resolution range)	90.7 (30.00-2.70) 88.7 (29.82-2.70)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.95 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.8.0124	Depositor
R, $R_{free}$	0.201 , 0.251 0.206 , 0.256	Depositor DCC
$R_{free}$ test set	2898 reflections (5.49%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.3	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , -2.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.399 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	16002	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAP, UD1, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.64	1/2595 (0.0%)	0.82	1/3486 (0.0%)
1	B	0.57	0/2540	0.75	0/3412
1	C	0.59	0/2628	0.80	1/3533 (0.0%)
1	D	0.61	0/2604	0.83	4/3501 (0.1%)
1	E	0.59	0/2590	0.78	0/3482
1	F	0.59	0/2585	0.79	1/3475 (0.0%)
All	All	0.60	1/15542 (0.0%)	0.80	7/20889 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	301	TYR	CE2-CZ	-5.25	1.31	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	333	GLN	N-CA-C	5.87	126.86	111.00
1	D	58	ARG	NE-CZ-NH1	5.55	123.07	120.30
1	D	326	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	F	174	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	A	200	ARG	NE-CZ-NH1	5.23	122.91	120.30
1	D	326	ARG	NE-CZ-NH2	-5.19	117.70	120.30
1	C	326	ARG	NE-CZ-NH1	5.09	122.85	120.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2559	0	2598	20	0
1	B	2504	0	2552	18	0
1	C	2590	0	2642	28	0
1	D	2566	0	2616	38	0
1	E	2553	0	2604	20	0
1	F	2549	0	2576	21	0
2	A	48	0	25	1	0
2	B	48	0	25	0	0
2	C	48	0	25	1	0
2	D	48	0	25	2	0
2	E	48	0	25	0	0
2	F	48	0	25	2	0
3	A	39	0	25	1	0
3	B	39	0	25	0	0
3	C	39	0	25	0	0
3	D	39	0	25	0	0
3	E	39	0	25	0	0
3	F	39	0	25	1	0
4	A	10	0	0	0	0
4	B	5	0	0	0	0
4	F	10	0	0	1	0
5	A	34	0	0	0	0
5	B	21	0	0	1	0
5	C	25	0	0	0	0
5	D	21	0	0	0	0
5	E	18	0	0	0	0
5	F	15	0	0	0	0
All	All	16002	0	15888	133	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 4.

All (133) 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:D:182:GLU:OE1	1:D:335:LEU:HD12	1.82	0.79
1:D:17:ILE:HD11	1:D:170:VAL:HG21	1.72	0.72
1:F:17:ILE:HD11	1:F:170:VAL:HG21	1.72	0.71
1:B:31:LYS:NZ	5:B:501:HOH:O	2.18	0.70
1:B:64:ILE:HD13	1:B:106:GLY:HA3	1.74	0.70
1:A:131:ILE:HD11	1:A:166:ARG:CZ	2.23	0.69
1:C:244:GLU:HG2	1:C:330:LEU:HD21	1.76	0.65
1:C:26:LEU:HD11	1:C:57:ILE:HD11	1.79	0.64
1:C:97:PHE:CD1	1:D:108:GLN:HG3	2.32	0.64
1:E:26:LEU:HD11	1:E:57:ILE:HD11	1.80	0.63
1:D:243:GLU:OE1	1:D:326:ARG:NH2	2.32	0.63
1:C:17:ILE:HD11	1:C:170:VAL:HG21	1.81	0.63
1:E:64:ILE:HD13	1:E:106:GLY:HA3	1.80	0.62
1:E:56:ASN:HD22	1:E:56:ASN:H	1.48	0.61
1:A:299:THR:HG22	1:A:301:TYR:CE2	2.36	0.61
1:B:155:LYS:HE2	1:B:160:THR:O	2.01	0.61
1:C:246:THR:HG21	1:C:254:GLU:N	2.16	0.60
1:D:261:ILE:HG22	1:D:262:GLY:O	2.01	0.60
1:C:128:ASN:HB2	1:C:272:GLU:O	2.03	0.58
1:C:108:GLN:HG3	1:D:97:PHE:CD1	2.38	0.58
1:C:108:GLN:HE21	1:C:147:LEU:HG	1.69	0.58
1:C:168:GLY:HA3	1:C:229:MET:HE1	1.85	0.58
1:A:121:LYS:HG3	1:A:161:THR:HB	1.87	0.57
1:F:14:THR:HG1	2:F:401:NAP:HO3A	1.53	0.57
1:E:108:GLN:HG3	1:F:97:PHE:CD1	2.40	0.57
1:F:87:LYS:HD3	3:F:402:UD1:H8'3	1.87	0.56
1:D:64:ILE:HD13	1:D:106:GLY:HA3	1.86	0.56
1:A:4:LYS:HE2	1:A:5:ASP:HB2	1.88	0.55
1:F:58:ARG:NH1	4:F:404:SO4:O2	2.37	0.55
1:D:192:THR:HG23	1:D:261:ILE:HD12	1.88	0.54
1:B:157:SER:HB2	1:F:75:GLU:OE1	2.08	0.54
1:C:261:ILE:HG22	1:C:262:GLY:O	2.07	0.54
1:A:97:PHE:CD1	1:B:108:GLN:HG3	2.43	0.54
1:D:22:LEU:HD13	1:D:33:VAL:HG11	1.90	0.53
1:E:243:GLU:OE1	1:E:326:ARG:NH2	2.37	0.53
1:A:223:GLU:HG2	1:A:301:TYR:OH	2.09	0.53
1:C:55:ARG:O	1:C:58:ARG:NH2	2.42	0.52
1:F:9:LEU:HB2	1:F:77:ILE:HD13	1.91	0.52
1:E:202:MET:HB2	1:E:232:ILE:HD12	1.90	0.52
1:D:193:VAL:HG21	1:D:196:LEU:HD21	1.92	0.52
1:F:41:TYR:CZ	1:F:45:LEU:HD11	2.44	0.52
1:C:125:THR:HG23	1:C:167:PHE:CE1	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:64:ILE:HD11	1:D:102:THR:HG23	1.91	0.51
1:A:302:GLU:O	1:A:303:ASN:HB2	2.11	0.51
1:F:184:GLN:NE2	1:F:192:THR:OG1	2.40	0.51
1:B:125:THR:HG23	1:B:167:PHE:CE1	2.47	0.50
1:D:201:PHE:HB3	1:D:229:MET:HE2	1.91	0.50
1:E:191:ILE:HG23	1:E:193:VAL:HG23	1.93	0.50
1:D:331:ASN:O	1:D:332:GLN:HG3	2.11	0.50
1:D:168:GLY:HA3	1:D:229:MET:HE1	1.94	0.50
1:C:263:LEU:HD21	1:C:269:MET:HE2	1.93	0.49
1:E:192:THR:HG23	1:E:261:ILE:HG21	1.94	0.49
1:A:131:ILE:HD12	1:A:131:ILE:N	2.27	0.49
1:D:201:PHE:HB3	1:D:229:MET:CE	2.43	0.49
1:A:211:LEU:HD22	1:A:227:LEU:HB2	1.95	0.48
1:A:19:LYS:HD3	1:A:46:LEU:HD13	1.95	0.48
1:C:131:ILE:N	1:C:131:ILE:HD12	2.28	0.48
1:D:128:ASN:HB2	1:D:272:GLU:O	2.13	0.48
1:A:332:GLN:O	1:A:334:LEU:N	2.47	0.48
1:B:128:ASN:HB2	1:B:272:GLU:O	2.14	0.48
1:D:61:ILE:HG21	1:E:86:MET:HE1	1.96	0.48
1:D:128:ASN:ND2	1:D:271:GLU:HB3	2.29	0.47
1:F:131:ILE:HD11	1:F:166:ARG:CZ	2.43	0.47
1:E:17:ILE:HD13	1:E:167:PHE:HE1	1.78	0.47
1:C:108:GLN:NE2	1:C:147:LEU:HG	2.27	0.47
1:D:81:PHE:CZ	1:D:216:MET:HG2	2.49	0.47
1:F:168:GLY:HA3	1:F:229:MET:HE1	1.97	0.47
1:F:128:ASN:HB2	1:F:272:GLU:O	2.14	0.47
1:B:26:LEU:HD11	1:B:57:ILE:HD11	1.97	0.46
1:C:9:LEU:HB2	1:C:77:ILE:HD13	1.97	0.46
1:C:155:LYS:HE3	1:C:160:THR:O	2.16	0.46
1:C:192:THR:HG21	1:C:261:ILE:HD12	1.97	0.46
1:F:332:GLN:HB3	1:F:334:LEU:HG	1.98	0.46
1:D:131:ILE:HD11	1:D:166:ARG:NH1	2.31	0.46
1:F:14:THR:OG1	2:F:401:NAP:O3B	2.28	0.46
1:B:237:LEU:HA	1:B:325:LEU:HD21	1.98	0.46
1:F:211:LEU:HD22	1:F:227:LEU:HB2	1.97	0.45
1:B:246:THR:HG21	1:B:254:GLU:N	2.31	0.45
1:C:275:THR:HG21	1:D:134:THR:HG23	1.96	0.45
1:A:37:SER:O	1:A:61:ILE:HA	2.16	0.45
1:C:167:PHE:HB2	2:C:400:NAP:C5N	2.46	0.45
1:D:155:LYS:CE	1:D:160:THR:O	2.65	0.45
1:E:87:LYS:HA	1:E:137:TYR:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:155:LYS:HE2	1:D:160:THR:O	2.17	0.44
1:A:87:LYS:HE3	2:A:401:NAP:O2N	2.17	0.44
1:D:186:LYS:HE2	1:D:249:TYR:CZ	2.52	0.44
1:D:61:ILE:HG21	1:E:86:MET:CE	2.48	0.44
1:D:86:MET:CE	1:E:61:ILE:HG21	2.48	0.44
1:F:184:GLN:HB3	1:F:190:LYS:O	2.17	0.44
1:F:193:VAL:HG21	1:F:196:LEU:HD21	2.00	0.44
1:B:149:THR:HG21	1:B:294:PRO:HD2	2.00	0.44
1:D:167:PHE:HB2	2:D:400:NAP:C5N	2.47	0.44
1:B:26:LEU:HD11	1:B:57:ILE:CD1	2.47	0.44
1:C:220:LYS:N	1:C:223:GLU:OE1	2.51	0.44
1:D:131:ILE:HG23	1:D:275:THR:HG23	1.99	0.44
1:D:39:SER:HB3	2:D:400:NAP:O1X	2.18	0.43
1:D:131:ILE:HD11	1:D:166:ARG:CZ	2.49	0.43
1:B:276:HIS:O	1:B:279:SER:OG	2.32	0.43
1:D:26:LEU:HD21	1:D:57:ILE:CD1	2.48	0.43
1:D:184:GLN:HB3	1:D:190:LYS:O	2.19	0.43
1:E:125:THR:HG23	1:E:167:PHE:CE1	2.54	0.43
1:A:194:THR:HG23	3:A:402:UD1:C2	2.48	0.43
1:C:155:LYS:CE	1:C:160:THR:O	2.66	0.43
1:A:223:GLU:CG	1:A:301:TYR:OH	2.67	0.43
1:C:168:GLY:HA3	1:C:229:MET:CE	2.47	0.43
1:B:9:LEU:HB2	1:B:77:ILE:HD13	2.01	0.42
1:D:22:LEU:CD1	1:D:33:VAL:HG11	2.49	0.42
1:C:164:SER:O	1:C:224:THR:HA	2.19	0.42
1:A:96:PRO:HB2	1:B:147:LEU:HD12	2.02	0.42
1:A:28:GLU:OE2	1:A:217:LYS:NZ	2.52	0.42
1:B:193:VAL:HG11	1:B:196:LEU:HG	2.01	0.42
1:A:43:GLN:O	1:A:47:GLN:HB2	2.20	0.41
1:E:17:ILE:HD13	1:E:167:PHE:CE1	2.55	0.41
1:C:186:LYS:CE	1:C:249:TYR:CZ	3.04	0.41
1:E:146:ARG:HG2	1:F:136:ASN:HB2	2.02	0.41
1:F:191:ILE:HG23	1:F:193:VAL:HG23	2.02	0.41
1:C:149:THR:HG21	1:C:294:PRO:HD2	2.02	0.41
1:E:226:ILE:HG21	1:E:273:LEU:HB3	2.03	0.41
1:E:131:ILE:HD11	1:E:166:ARG:NH1	2.35	0.41
1:A:20:SER:HB2	1:A:205:LEU:HD11	2.01	0.41
1:B:74:MET:HA	1:B:77:ILE:HD12	2.02	0.41
1:C:125:THR:HG22	1:C:125:THR:O	2.21	0.41
1:F:244:GLU:HG2	1:F:330:LEU:HD21	2.01	0.41
1:D:86:MET:HE2	1:E:61:ILE:HG21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:263:LEU:HD21	1:D:269:MET:CE	2.51	0.41
1:A:204:THR:HG22	1:A:332:GLN:OE1	2.21	0.41
1:D:125:THR:HG22	1:D:125:THR:O	2.20	0.41
1:D:184:GLN:HE22	1:D:192:THR:HG22	1.85	0.41
1:F:226:ILE:HD11	1:F:291:ILE:HD12	2.03	0.41
1:E:226:ILE:HD11	1:E:291:ILE:HD12	2.03	0.41
1:C:263:LEU:HD11	1:C:269:MET:CE	2.51	0.40
1:B:131:ILE:HD11	1:B:166:ARG:CZ	2.50	0.40
1:D:128:ASN:HD22	1:D:128:ASN:C	2.25	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	314/343 (92%)	294 (94%)	17 (5%)	3 (1%)	18	43
1	B	308/343 (90%)	297 (96%)	10 (3%)	1 (0%)	44	73
1	C	321/343 (94%)	300 (94%)	19 (6%)	2 (1%)	28	56
1	D	318/343 (93%)	300 (94%)	15 (5%)	3 (1%)	20	46
1	E	316/343 (92%)	296 (94%)	16 (5%)	4 (1%)	14	35
1	F	313/343 (91%)	292 (93%)	19 (6%)	2 (1%)	28	56
All	All	1890/2058 (92%)	1779 (94%)	96 (5%)	15 (1%)	22	49

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	155	LYS
1	E	322	LYS
1	F	51	ARG

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Mol	Chain	Res	Type
1	F	155	LYS
1	A	155	LYS
1	E	51	ARG
1	C	55	ARG
1	D	155	LYS
1	B	155	LYS
1	E	185	ILE
1	A	333	GLN
1	D	309	ALA
1	A	294	PRO
1	D	332	GLN
1	E	85	ALA

### 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	284/302 (94%)	259 (91%)	25 (9%)	12	27
1	B	279/302 (92%)	261 (94%)	18 (6%)	20	44
1	C	288/302 (95%)	262 (91%)	26 (9%)	11	25
1	D	285/302 (94%)	256 (90%)	29 (10%)	8	20
1	E	285/302 (94%)	265 (93%)	20 (7%)	18	40
1	F	282/302 (93%)	260 (92%)	22 (8%)	15	33
All	All	1703/1812 (94%)	1563 (92%)	140 (8%)	13	30

All (140) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	24	ASN
1	A	39	SER
1	A	77	ILE
1	A	95	ASN
1	A	108	GLN

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Mol	Chain	Res	Type
1	A	121	LYS
1	A	163	THR
1	A	165	VAL
1	A	183	ASN
1	A	187	GLU
1	A	191	ILE
1	A	193	VAL
1	A	214	GLU
1	A	216	MET
1	A	261	ILE
1	A	263	LEU
1	A	277	ASP
1	A	299	THR
1	A	302	GLU
1	A	308	LYS
1	A	323	GLU
1	A	326	ARG
1	A	328	LEU
1	A	332	GLN
1	B	53	LYS
1	B	57	ILE
1	B	64	ILE
1	B	108	GLN
1	B	128	ASN
1	B	146	ARG
1	B	157	SER
1	B	174	ARG
1	B	187	GLU
1	B	191	ILE
1	B	193	VAL
1	B	223	GLU
1	B	229	MET
1	B	235	ASN
1	B	274	MET
1	B	285	LEU
1	B	308	LYS
1	B	326	ARG
1	C	1	MET
1	C	5	ASP
1	C	20	SER
1	C	47	GLN
1	C	51	ARG

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Mol	Chain	Res	Type
1	C	52	ASP
1	C	53	LYS
1	C	77	ILE
1	C	128	ASN
1	C	134	THR
1	C	147	LEU
1	C	157	SER
1	C	161	THR
1	C	193	VAL
1	C	197	SER
1	C	216	MET
1	C	217	LYS
1	C	223	GLU
1	C	244	GLU
1	C	256	ILE
1	C	259	GLU
1	C	260	GLU
1	C	261	ILE
1	C	263	LEU
1	C	281	GLN
1	C	328	LEU
1	D	1	MET
1	D	4	LYS
1	D	37	SER
1	D	51	ARG
1	D	53	LYS
1	D	55	ARG
1	D	77	ILE
1	D	125	THR
1	D	128	ASN
1	D	146	ARG
1	D	147	LEU
1	D	161	THR
1	D	163	THR
1	D	192	THR
1	D	193	VAL
1	D	216	MET
1	D	228	LYS
1	D	235	ASN
1	D	244	GLU
1	D	256	ILE
1	D	260	GLU

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Mol	Chain	Res	Type
1	D	263	LEU
1	D	281	GLN
1	D	287	ASP
1	D	306	ARG
1	D	323	GLU
1	D	328	LEU
1	D	332	GLN
1	D	333	GLN
1	E	1	MET
1	E	5	ASP
1	E	39	SER
1	E	53	LYS
1	E	55	ARG
1	E	56	ASN
1	E	60	LEU
1	E	64	ILE
1	E	128	ASN
1	E	146	ARG
1	E	163	THR
1	E	176	SER
1	E	183	ASN
1	E	187	GLU
1	E	191	ILE
1	E	218	ILE
1	E	223	GLU
1	E	228	LYS
1	E	233	SER
1	E	246	THR
1	F	1	MET
1	F	23	SER
1	F	24	ASN
1	F	39	SER
1	F	51	ARG
1	F	53	LYS
1	F	77	ILE
1	F	87	LYS
1	F	125	THR
1	F	128	ASN
1	F	134	THR
1	F	165	VAL
1	F	187	GLU
1	F	193	VAL

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Mol	Chain	Res	Type
1	F	206	ASN
1	F	214	GLU
1	F	228	LYS
1	F	235	ASN
1	F	244	GLU
1	F	277	ASP
1	F	306	ARG
1	F	333	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	GLN
1	A	128	ASN
1	A	183	ASN
1	A	300	GLN
1	B	108	GLN
1	B	128	ASN
1	B	184	GLN
1	C	108	GLN
1	C	128	ASN
1	C	189	GLN
1	D	47	GLN
1	D	108	GLN
1	D	128	ASN
1	D	327	ASN
1	E	56	ASN
1	E	188	ASN
1	F	24	ASN
1	F	47	GLN
1	F	183	ASN
1	F	235	ASN

### 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 ⓘ

17 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	-	44,52,52	1.68	4 (9%)	51,80,80	1.70	7 (13%)
3	UD1	A	402	-	34,41,41	0.63	1 (2%)	43,62,62	1.58	3 (6%)
4	SO4	A	403	-	4,4,4	0.44	0	6,6,6	0.38	0
4	SO4	A	404	-	4,4,4	0.52	0	6,6,6	0.39	0
2	NAP	B	401	-	44,52,52	1.51	3 (6%)	51,80,80	2.15	13 (25%)
3	UD1	B	402	-	34,41,41	0.73	0	43,62,62	1.98	7 (16%)
4	SO4	B	403	-	4,4,4	0.46	0	6,6,6	0.63	0
2	NAP	C	400	-	44,52,52	1.45	3 (6%)	51,80,80	1.86	10 (19%)
3	UD1	C	401	-	34,41,41	0.60	0	43,62,62	1.72	4 (9%)
2	NAP	D	400	-	44,52,52	1.62	5 (11%)	51,80,80	1.82	9 (17%)
3	UD1	D	401	-	34,41,41	0.67	0	43,62,62	1.55	5 (11%)
2	NAP	E	400	-	44,52,52	1.60	5 (11%)	51,80,80	1.96	10 (19%)
3	UD1	E	401	-	34,41,41	0.75	1 (2%)	43,62,62	1.70	4 (9%)
2	NAP	F	401	-	44,52,52	1.53	4 (9%)	51,80,80	1.82	8 (15%)
3	UD1	F	402	-	34,41,41	0.76	1 (2%)	43,62,62	1.86	9 (20%)
4	SO4	F	403	-	4,4,4	0.42	0	6,6,6	0.26	0
4	SO4	F	404	-	4,4,4	0.42	0	6,6,6	0.57	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAP	A	401	-	-	0/27/67/67	0/5/5/5
3	UD1	A	402	-	-	0/22/63/63	0/3/3/3
4	SO4	A	403	-	-	0/0/0/0	0/0/0/0
4	SO4	A	404	-	-	0/0/0/0	0/0/0/0
2	NAP	B	401	-	-	0/27/67/67	0/5/5/5
3	UD1	B	402	-	-	0/22/63/63	0/3/3/3
4	SO4	B	403	-	-	0/0/0/0	0/0/0/0
2	NAP	C	400	-	-	0/27/67/67	0/5/5/5
3	UD1	C	401	-	-	0/22/63/63	0/3/3/3
2	NAP	D	400	-	-	0/27/67/67	0/5/5/5
3	UD1	D	401	-	-	0/22/63/63	0/3/3/3
2	NAP	E	400	-	-	0/27/67/67	0/5/5/5
3	UD1	E	401	-	-	0/22/63/63	0/3/3/3
2	NAP	F	401	-	-	0/27/67/67	0/5/5/5
3	UD1	F	402	-	-	0/22/63/63	0/3/3/3
4	SO4	F	403	-	-	0/0/0/0	0/0/0/0
4	SO4	F	404	-	-	0/0/0/0	0/0/0/0

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	402	UD1	C2-N3	-2.47	1.33	1.38
2	E	400	NAP	C5A-N7A	-2.10	1.32	1.39
3	A	402	UD1	C2-N3	-2.03	1.34	1.38
3	E	401	UD1	C4-N3	2.11	1.36	1.33
2	F	401	NAP	O4D-C1D	2.16	1.44	1.41
2	D	400	NAP	O4D-C1D	2.17	1.44	1.41
2	E	400	NAP	O4D-C1D	2.18	1.44	1.41
2	D	400	NAP	C2A-N3A	2.35	1.36	1.32
2	B	401	NAP	C5A-C4A	2.73	1.46	1.40
2	A	401	NAP	C5A-C4A	2.85	1.46	1.40
2	A	401	NAP	O4B-C1B	2.89	1.45	1.41
2	E	400	NAP	C5A-C4A	2.94	1.47	1.40
2	F	401	NAP	C5A-C4A	3.06	1.47	1.40
2	B	401	NAP	C5N-C4N	3.15	1.44	1.38
2	C	400	NAP	C5A-C4A	3.55	1.48	1.40
2	D	400	NAP	C5A-C4A	3.58	1.48	1.40
2	E	400	NAP	C5N-C4N	3.81	1.46	1.38
2	F	401	NAP	C5N-C4N	3.96	1.46	1.38
2	C	400	NAP	C5N-C4N	4.32	1.47	1.38
2	A	401	NAP	C5N-C4N	4.75	1.48	1.38
2	D	400	NAP	C5N-C4N	4.75	1.48	1.38
2	C	400	NAP	C4N-C3N	5.84	1.49	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	400	NAP	C4N-C3N	6.44	1.49	1.39
2	F	401	NAP	C4N-C3N	7.08	1.51	1.39
2	A	401	NAP	C4N-C3N	7.64	1.51	1.39
2	E	400	NAP	C4N-C3N	7.71	1.52	1.39
2	B	401	NAP	C4N-C3N	7.79	1.52	1.39

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	NAP	C5N-C4N-C3N	-7.26	111.81	120.35
2	C	400	NAP	N3A-C2A-N1A	-7.17	122.62	128.86
2	B	401	NAP	N3A-C2A-N1A	-6.96	122.80	128.86
2	E	400	NAP	N3A-C2A-N1A	-6.87	122.87	128.86
2	D	400	NAP	C5N-C4N-C3N	-6.75	112.41	120.35
2	B	401	NAP	C5N-C4N-C3N	-6.72	112.45	120.35
2	E	400	NAP	C5N-C4N-C3N	-6.26	112.99	120.35
2	D	400	NAP	N3A-C2A-N1A	-6.22	123.44	128.86
2	C	400	NAP	C5N-C4N-C3N	-5.95	113.35	120.35
2	A	401	NAP	N3A-C2A-N1A	-5.74	123.86	128.86
2	F	401	NAP	N3A-C2A-N1A	-5.67	123.92	128.86
2	A	401	NAP	C5N-C4N-C3N	-5.37	114.03	120.35
2	B	401	NAP	C4A-C5A-N7A	-3.73	105.81	109.41
3	F	402	UD1	O5'-C1'-C2'	-3.56	103.35	110.65
2	E	400	NAP	O4B-C1B-C2B	-3.53	100.41	106.59
2	A	401	NAP	C4A-C5A-N7A	-3.48	106.05	109.41
3	C	401	UD1	O5'-C1'-O1'	-3.28	107.07	111.36
2	B	401	NAP	O7N-C7N-N7N	-3.02	118.29	122.58
2	E	400	NAP	O7N-C7N-N7N	-2.98	118.35	122.58
2	D	400	NAP	O2B-P2B-O1X	-2.93	97.77	109.26
3	F	402	UD1	C4B-O4B-C1B	-2.89	106.69	109.77
2	F	401	NAP	O7N-C7N-C3N	-2.87	116.27	119.62
2	D	400	NAP	C4A-C5A-N7A	-2.80	106.70	109.41
2	F	401	NAP	C1B-N9A-C4A	-2.80	121.80	126.64
2	E	400	NAP	C4A-C5A-N7A	-2.67	106.83	109.41
2	A	401	NAP	O7N-C7N-C3N	-2.60	116.58	119.62
2	C	400	NAP	C4A-C5A-N7A	-2.60	106.90	109.41
2	B	401	NAP	C1B-N9A-C4A	-2.59	122.17	126.64
2	C	400	NAP	O2B-P2B-O1X	-2.58	99.15	109.26
3	F	402	UD1	C1'-C2'-N2'	-2.55	106.71	111.03
2	E	400	NAP	O2B-P2B-O1X	-2.50	99.44	109.26
2	B	401	NAP	O2B-P2B-O1X	-2.43	99.73	109.26
3	A	402	UD1	C1'-C2'-N2'	-2.41	106.96	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	400	NAP	C1B-N9A-C4A	-2.36	122.55	126.64
2	F	401	NAP	C4A-C5A-N7A	-2.35	107.14	109.41
3	B	402	UD1	O4'-C4'-C5'	-2.25	103.63	109.28
2	C	400	NAP	O3B-C3B-C4B	-2.21	104.62	111.09
3	E	401	UD1	C1'-C2'-N2'	-2.20	107.31	111.03
3	B	402	UD1	C6-N1-C2	-2.20	117.72	121.28
2	A	401	NAP	C4D-O4D-C1D	-2.18	107.44	109.77
2	B	401	NAP	C4D-O4D-C1D	-2.12	107.51	109.77
2	B	401	NAP	O7N-C7N-C3N	-2.12	117.14	119.62
3	C	401	UD1	C6-N1-C2	-2.12	117.85	121.28
3	D	401	UD1	O7'-C7'-C8'	-2.07	118.28	122.06
2	D	400	NAP	C1B-N9A-C4A	-2.04	123.11	126.64
2	B	401	NAP	O2X-P2B-O2B	-2.01	96.87	106.00
3	F	402	UD1	O4B-C4B-C3B	2.00	109.15	105.17
2	E	400	NAP	C2A-N1A-C6A	2.02	122.30	118.77
3	E	401	UD1	C3'-C4'-C5'	2.03	113.80	110.22
3	D	401	UD1	C3'-C4'-C5'	2.05	113.83	110.22
2	C	400	NAP	C2A-N1A-C6A	2.06	122.38	118.77
2	C	400	NAP	C6N-C5N-C4N	2.08	122.58	119.44
2	A	401	NAP	O3X-P2B-O2X	2.09	116.03	107.61
3	D	401	UD1	O7'-C7'-N2'	2.10	125.97	121.92
3	A	402	UD1	C3'-C4'-C5'	2.15	114.01	110.22
2	D	400	NAP	O2N-PN-O1N	2.18	123.57	112.28
3	B	402	UD1	O2A-PA-O1A	2.20	123.65	112.28
2	E	400	NAP	N6A-C6A-N1A	2.22	123.16	118.77
3	C	401	UD1	PB-O1'-C1'	2.23	128.60	119.74
2	F	401	NAP	C6N-C5N-C4N	2.27	122.86	119.44
2	E	400	NAP	C4B-O4B-C1B	2.27	112.19	109.77
3	D	401	UD1	PB-O1'-C1'	2.29	128.82	119.74
2	B	401	NAP	C2A-N1A-C6A	2.31	122.82	118.77
3	E	401	UD1	O4B-C1B-N1	2.35	112.80	108.08
2	F	401	NAP	C3N-C2N-N1N	2.36	122.81	120.43
3	B	402	UD1	O5'-C5'-C6'	2.39	112.14	106.41
3	F	402	UD1	C3'-C2'-N2'	2.47	115.35	110.61
2	D	400	NAP	C6N-C5N-C4N	2.52	123.24	119.44
2	B	401	NAP	O2X-P2B-O1X	2.60	120.67	110.50
3	F	402	UD1	C3'-C4'-C5'	2.66	114.90	110.22
2	B	401	NAP	C2N-C3N-C4N	2.72	121.37	118.26
3	B	402	UD1	O4B-C1B-N1	2.79	113.67	108.08
2	C	400	NAP	C3N-C2N-N1N	2.91	123.36	120.43
2	D	400	NAP	C4B-O4B-C1B	2.94	112.90	109.77
2	D	400	NAP	C3N-C2N-N1N	2.97	123.42	120.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	NAP	C3N-C7N-N7N	3.24	121.47	117.77
3	F	402	UD1	O1'-C1'-C2'	3.45	114.64	108.40
3	B	402	UD1	C3'-C4'-C5'	3.45	116.30	110.22
2	C	400	NAP	C3N-C7N-N7N	3.48	121.75	117.77
3	F	402	UD1	C4'-C3'-C2'	3.55	115.59	110.33
2	E	400	NAP	C3N-C7N-N7N	4.07	122.42	117.77
2	A	401	NAP	C3N-C7N-N7N	4.17	122.54	117.77
2	B	401	NAP	C3N-C7N-N7N	5.90	124.51	117.77
3	F	402	UD1	C4-N3-C2	7.22	120.33	114.13
3	D	401	UD1	C4-N3-C2	7.37	120.46	114.13
3	A	402	UD1	C4-N3-C2	7.91	120.92	114.13
3	E	401	UD1	C4-N3-C2	8.43	121.37	114.13
3	C	401	UD1	C4-N3-C2	8.78	121.68	114.13
3	B	402	UD1	C4-N3-C2	9.59	122.37	114.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAP	1	0
3	A	402	UD1	1	0
2	C	400	NAP	1	0
2	D	400	NAP	2	0
2	F	401	NAP	2	0
3	F	402	UD1	1	0
4	F	404	SO4	1	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	324/343 (94%)	-0.42	1 (0%) 93 94	10, 23, 45, 61	0
1	B	316/343 (92%)	-0.34	1 (0%) 93 94	9, 26, 53, 74	0
1	C	327/343 (95%)	-0.43	0 100 100	10, 22, 40, 68	0
1	D	324/343 (94%)	-0.43	0 100 100	9, 21, 43, 60	0
1	E	322/343 (93%)	-0.31	1 (0%) 93 94	10, 28, 57, 75	0
1	F	323/343 (94%)	-0.39	2 (0%) 89 90	10, 24, 49, 77	0
All	All	1936/2058 (94%)	-0.39	5 (0%) 93 94	9, 24, 49, 77	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	316	ASN	2.8
1	A	261	ILE	2.7
1	F	0	HIS	2.3
1	B	322	LYS	2.3
1	F	302	GLU	2.2

### 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 ⓘ

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(Å <sup>2</sup> )	Q<0.9
4	SO4	A	404	5/5	0.97	0.13	0.20	57,58,61,63	0
2	NAP	B	401	48/48	0.97	0.12	-0.13	14,19,30,34	0
2	NAP	E	400	48/48	0.97	0.12	-0.34	14,17,27,29	0
3	UD1	B	402	39/39	0.97	0.14	-0.35	19,20,31,32	0
2	NAP	D	400	48/48	0.97	0.12	-0.50	11,14,24,27	0
3	UD1	E	401	39/39	0.97	0.13	-0.51	23,25,30,31	0
3	UD1	F	402	39/39	0.97	0.13	-0.55	16,19,26,27	0
2	NAP	A	401	48/48	0.97	0.12	-0.67	10,12,23,25	0
2	NAP	F	401	48/48	0.97	0.11	-0.80	13,14,24,26	0
3	UD1	A	402	39/39	0.97	0.12	-0.85	17,20,25,25	0
3	UD1	C	401	39/39	0.97	0.13	-0.85	12,15,19,20	0
3	UD1	D	401	39/39	0.97	0.12	-0.91	13,16,18,18	0
2	NAP	C	400	48/48	0.98	0.10	-1.19	11,14,19,20	0
4	SO4	F	403	5/5	0.95	0.10	-	57,58,62,66	0
4	SO4	B	403	5/5	0.96	0.14	-	49,50,52,53	0
4	SO4	F	404	5/5	0.89	0.21	-	66,68,72,73	0
4	SO4	A	403	5/5	0.93	0.16	-	60,62,64,65	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.