



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:55 PM GMT

PDB ID : 4QAA
Title : X-RAY STRUCTURE OF ACETYLCHOLINE BINDING PROTEIN (ACHBP) IN COMPLEX WITH 6-(4-Methoxyphenyl)-N4-octylpyrimidine-2,4-diamine
Authors : Kaczanowska, K.; Harel, M.; Radic', Z.; Changeux, J.-P.; Finn, M.G.; Taylor, P.
Deposited on : 2014-05-03
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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

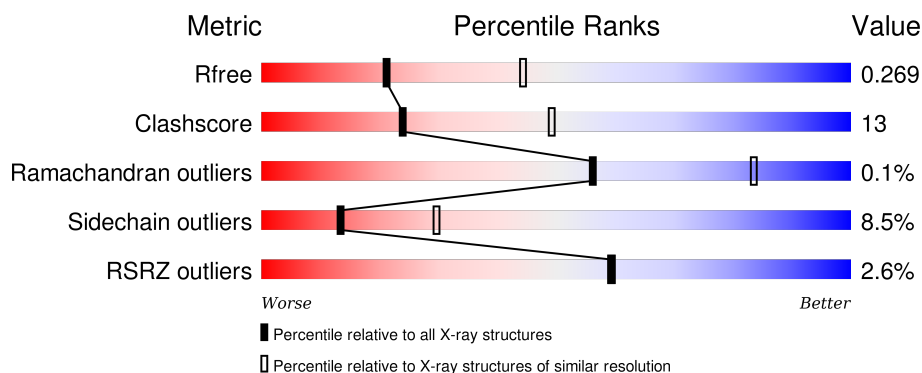
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (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 ≥ 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	217	<div> <div>4%</div> <div>74%</div> <div>18%</div> <div>• 5%</div> </div>
1	B	217	<div> <div>72%</div> <div>21%</div> <div>• 5%</div> </div>
1	C	217	<div> <div>%</div> <div>74%</div> <div>18%</div> <div>• •</div> </div>
1	D	217	<div> <div>4%</div> <div>70%</div> <div>25%</div> <div>• •</div> </div>
1	E	217	<div> <div>%</div> <div>72%</div> <div>20%</div> <div>• 6%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	217	
1	G	217	
1	H	217	
1	I	217	
1	J	217	

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
2	KK1	B	301	-	-	-	X
2	KK1	C	301	-	-	-	X
2	KK1	E	301	-	-	-	X
2	KK1	F	301	-	-	-	X
2	KK1	G	301	-	-	-	X
3	NAG	I	302	-	-	-	X
3	NAG	J	302	-	-	-	X
4	PO4	B	303	-	-	-	X
4	PO4	C	305	-	-	-	X
4	PO4	E	305	-	-	-	X
4	PO4	F	304	-	-	-	X
4	PO4	G	304	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17293 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 Acetylcholine-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1662	1037	284	338	3			
1	B	207	Total	C	N	O	S	0	0	0
			1656	1033	284	336	3			
1	C	209	Total	C	N	O	S	0	1	0
			1682	1050	286	343	3			
1	D	213	Total	C	N	O	S	0	0	0
			1706	1064	290	347	5			
1	E	205	Total	C	N	O	S	0	0	0
			1645	1028	281	333	3			
1	F	209	Total	C	N	O	S	0	0	0
			1675	1045	286	340	4			
1	G	208	Total	C	N	O	S	0	0	0
			1672	1047	285	337	3			
1	H	207	Total	C	N	O	S	0	0	0
			1664	1042	283	336	3			
1	I	207	Total	C	N	O	S	0	0	0
			1664	1042	283	336	3			
1	J	206	Total	C	N	O	S	0	0	0
			1655	1032	283	337	3			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	ASP	-	EXPRESSION TAG	UNP P58154
A	-6	TYR	-	EXPRESSION TAG	UNP P58154
A	-5	LYS	-	EXPRESSION TAG	UNP P58154
A	-4	ASP	-	EXPRESSION TAG	UNP P58154
A	-3	ASP	-	EXPRESSION TAG	UNP P58154
A	-2	ASP	-	EXPRESSION TAG	UNP P58154
A	-1	ASP	-	EXPRESSION TAG	UNP P58154
A	0	LYS	-	EXPRESSION TAG	UNP P58154
B	-7	ASP	-	EXPRESSION TAG	UNP P58154

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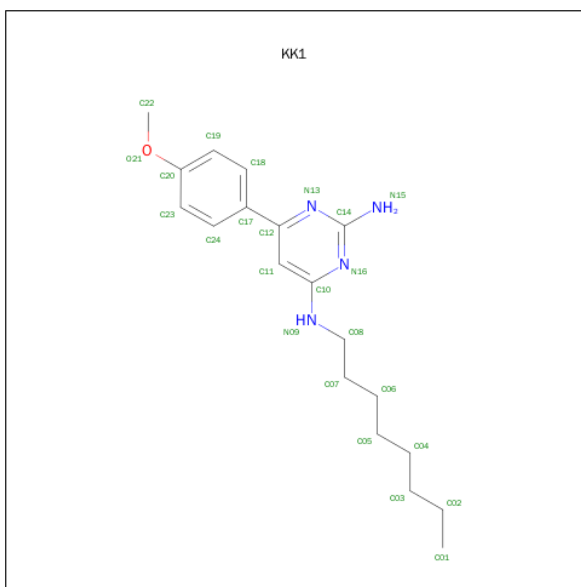
Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	TYR	-	EXPRESSION TAG	UNP P58154
B	-5	LYS	-	EXPRESSION TAG	UNP P58154
B	-4	ASP	-	EXPRESSION TAG	UNP P58154
B	-3	ASP	-	EXPRESSION TAG	UNP P58154
B	-2	ASP	-	EXPRESSION TAG	UNP P58154
B	-1	ASP	-	EXPRESSION TAG	UNP P58154
B	0	LYS	-	EXPRESSION TAG	UNP P58154
C	-7	ASP	-	EXPRESSION TAG	UNP P58154
C	-6	TYR	-	EXPRESSION TAG	UNP P58154
C	-5	LYS	-	EXPRESSION TAG	UNP P58154
C	-4	ASP	-	EXPRESSION TAG	UNP P58154
C	-3	ASP	-	EXPRESSION TAG	UNP P58154
C	-2	ASP	-	EXPRESSION TAG	UNP P58154
C	-1	ASP	-	EXPRESSION TAG	UNP P58154
C	0	LYS	-	EXPRESSION TAG	UNP P58154
D	-7	ASP	-	EXPRESSION TAG	UNP P58154
D	-6	TYR	-	EXPRESSION TAG	UNP P58154
D	-5	LYS	-	EXPRESSION TAG	UNP P58154
D	-4	ASP	-	EXPRESSION TAG	UNP P58154
D	-3	ASP	-	EXPRESSION TAG	UNP P58154
D	-2	ASP	-	EXPRESSION TAG	UNP P58154
D	-1	ASP	-	EXPRESSION TAG	UNP P58154
D	0	LYS	-	EXPRESSION TAG	UNP P58154
E	-7	ASP	-	EXPRESSION TAG	UNP P58154
E	-6	TYR	-	EXPRESSION TAG	UNP P58154
E	-5	LYS	-	EXPRESSION TAG	UNP P58154
E	-4	ASP	-	EXPRESSION TAG	UNP P58154
E	-3	ASP	-	EXPRESSION TAG	UNP P58154
E	-2	ASP	-	EXPRESSION TAG	UNP P58154
E	-1	ASP	-	EXPRESSION TAG	UNP P58154
E	0	LYS	-	EXPRESSION TAG	UNP P58154
F	-7	ASP	-	EXPRESSION TAG	UNP P58154
F	-6	TYR	-	EXPRESSION TAG	UNP P58154
F	-5	LYS	-	EXPRESSION TAG	UNP P58154
F	-4	ASP	-	EXPRESSION TAG	UNP P58154
F	-3	ASP	-	EXPRESSION TAG	UNP P58154
F	-2	ASP	-	EXPRESSION TAG	UNP P58154
F	-1	ASP	-	EXPRESSION TAG	UNP P58154
F	0	LYS	-	EXPRESSION TAG	UNP P58154
G	-7	ASP	-	EXPRESSION TAG	UNP P58154
G	-6	TYR	-	EXPRESSION TAG	UNP P58154
G	-5	LYS	-	EXPRESSION TAG	UNP P58154

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-4	ASP	-	EXPRESSION TAG	UNP P58154
G	-3	ASP	-	EXPRESSION TAG	UNP P58154
G	-2	ASP	-	EXPRESSION TAG	UNP P58154
G	-1	ASP	-	EXPRESSION TAG	UNP P58154
G	0	LYS	-	EXPRESSION TAG	UNP P58154
H	-7	ASP	-	EXPRESSION TAG	UNP P58154
H	-6	TYR	-	EXPRESSION TAG	UNP P58154
H	-5	LYS	-	EXPRESSION TAG	UNP P58154
H	-4	ASP	-	EXPRESSION TAG	UNP P58154
H	-3	ASP	-	EXPRESSION TAG	UNP P58154
H	-2	ASP	-	EXPRESSION TAG	UNP P58154
H	-1	ASP	-	EXPRESSION TAG	UNP P58154
H	0	LYS	-	EXPRESSION TAG	UNP P58154
I	-7	ASP	-	EXPRESSION TAG	UNP P58154
I	-6	TYR	-	EXPRESSION TAG	UNP P58154
I	-5	LYS	-	EXPRESSION TAG	UNP P58154
I	-4	ASP	-	EXPRESSION TAG	UNP P58154
I	-3	ASP	-	EXPRESSION TAG	UNP P58154
I	-2	ASP	-	EXPRESSION TAG	UNP P58154
I	-1	ASP	-	EXPRESSION TAG	UNP P58154
I	0	LYS	-	EXPRESSION TAG	UNP P58154
J	-7	ASP	-	EXPRESSION TAG	UNP P58154
J	-6	TYR	-	EXPRESSION TAG	UNP P58154
J	-5	LYS	-	EXPRESSION TAG	UNP P58154
J	-4	ASP	-	EXPRESSION TAG	UNP P58154
J	-3	ASP	-	EXPRESSION TAG	UNP P58154
J	-2	ASP	-	EXPRESSION TAG	UNP P58154
J	-1	ASP	-	EXPRESSION TAG	UNP P58154
J	0	LYS	-	EXPRESSION TAG	UNP P58154

- Molecule 2 is 6-(4-METHOXYPHENYL)-N 4 -OCTYLPYRIMIDINE-2,4-DIAMINE (three-letter code: KK1) (formula: C₁₉H₂₈N₄O).



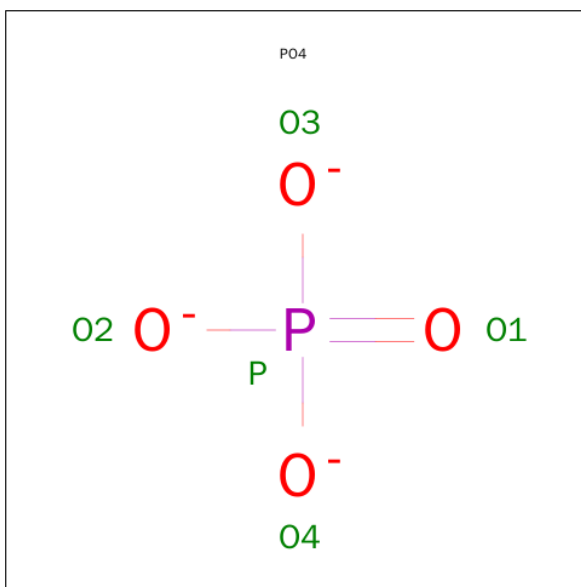
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			24	19	4	1		
2	B	1	Total	C	N	O	0	0
			24	19	4	1		
2	C	1	Total	C	N	O	0	0
			24	19	4	1		
2	D	1	Total	C	N	O	0	0
			24	19	4	1		
2	E	1	Total	C	N	O	0	0
			24	19	4	1		
2	F	1	Total	C	N	O	0	0
			24	19	4	1		
2	G	1	Total	C	N	O	0	0
			16	11	4	1		
2	H	1	Total	C	N	O	0	0
			16	11	4	1		
2	I	1	Total	C	N	O	0	0
			24	19	4	1		
2	J	1	Total	C	N	O	0	0
			16	11	4	1		

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	F	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		
3	H	1	Total	C	N	O	0	0
			14	8	1	5		
3	I	1	Total	C	N	O	0	0
			14	8	1	5		
3	J	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		
4	D	1	Total	O	P	0	0
			5	4	1		
4	D	1	Total	O	P	0	0
			5	4	1		
4	E	1	Total	O	P	0	0
			5	4	1		
4	E	1	Total	O	P	0	0
			5	4	1		
4	E	1	Total	O	P	0	0
			5	4	1		
4	F	1	Total	O	P	0	0
			5	4	1		
4	F	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	G	1	Total	O	P	0	0
			5	4	1		
4	G	1	Total	O	P	0	0
			5	4	1		
4	H	1	Total	O	P	0	0
			5	4	1		
4	H	1	Total	O	P	0	0
			5	4	1		
4	J	1	Total	O	P	0	0
			5	4	1		

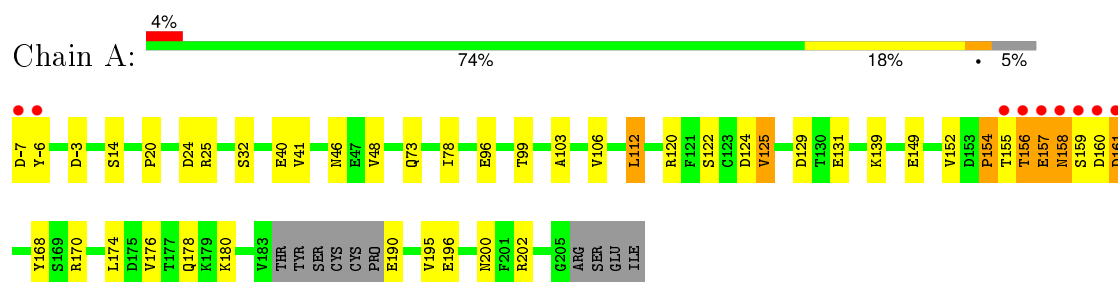
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	18	Total	O	0	0
			18	18		
5	B	17	Total	O	0	0
			17	17		
5	C	26	Total	O	0	0
			26	26		
5	D	22	Total	O	0	0
			22	22		
5	E	18	Total	O	0	0
			18	18		
5	F	9	Total	O	0	0
			9	9		
5	G	9	Total	O	0	0
			9	9		
5	H	14	Total	O	0	0
			14	14		
5	I	11	Total	O	0	0
			11	11		
5	J	17	Total	O	0	0
			17	17		

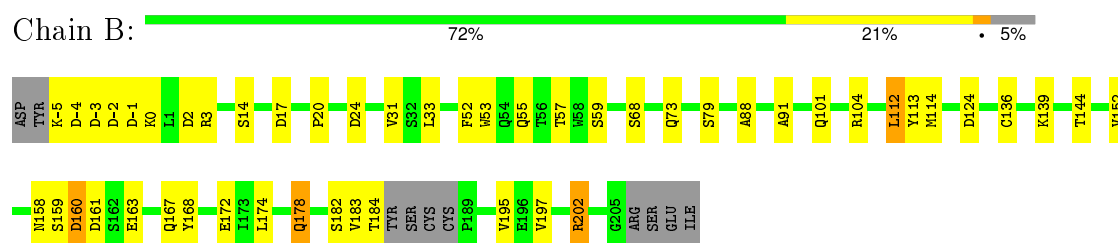
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

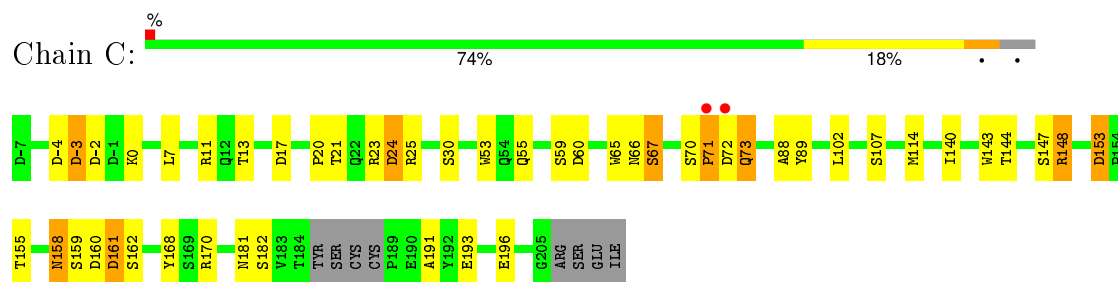
• Molecule 1: Acetylcholine-binding protein



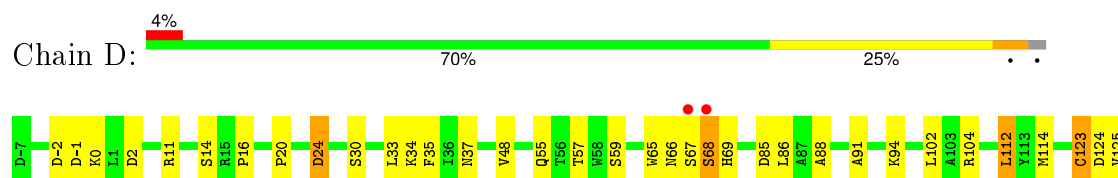
• Molecule 1: Acetylcholine-binding protein

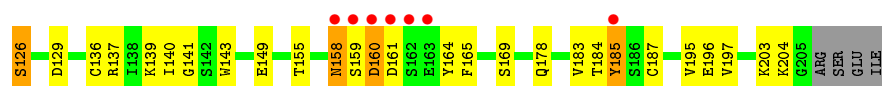


• Molecule 1: Acetylcholine-binding protein

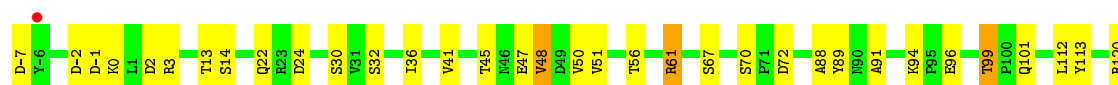


• Molecule 1: Acetylcholine-binding protein

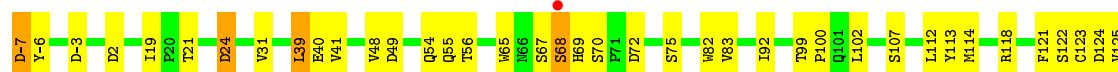




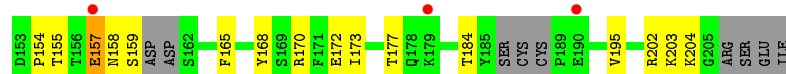
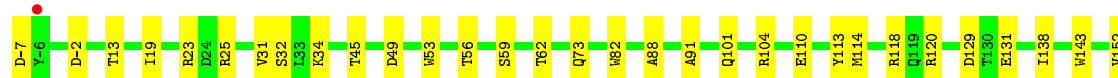
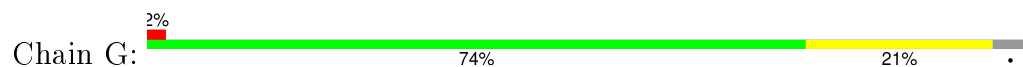
- Molecule 1: Acetylcholine-binding protein



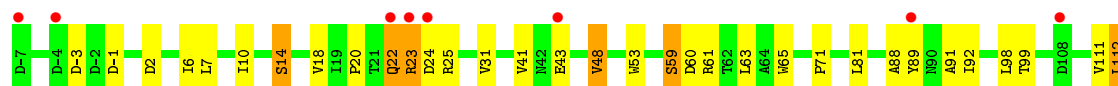
- Molecule 1: Acetylcholine-binding protein



- Molecule 1: Acetylcholine-binding protein

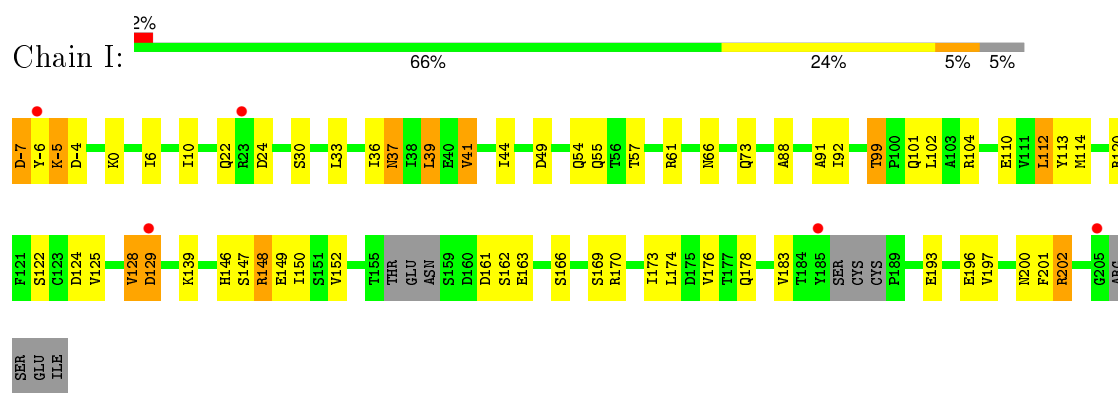


- Molecule 1: Acetylcholine-binding protein

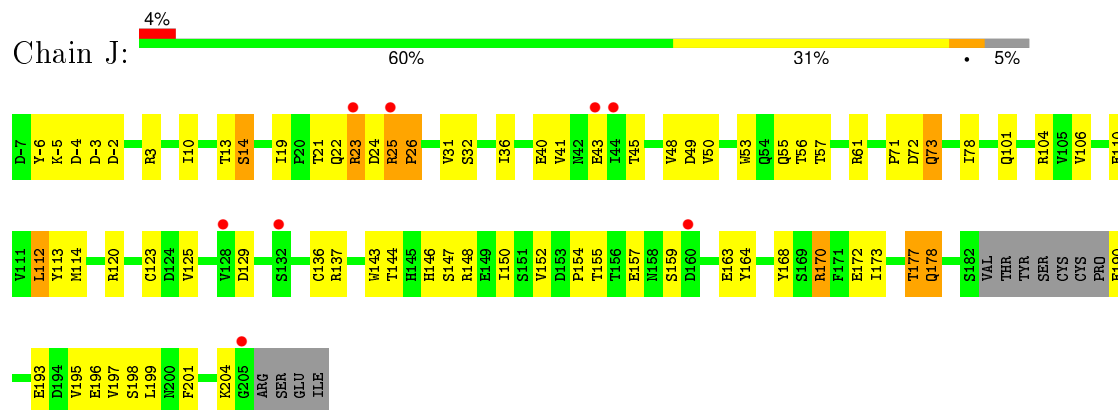


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- Molecule 1: Acetylcholine-binding protein



• Molecule 1: Acetylcholine-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.41Å 129.93Å 122.75Å 90.00° 106.21° 90.00°	Depositor
Resolution (Å)	49.65 – 2.70 49.65 – 2.18	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.65-2.70) 84.0 (49.65-2.18)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 2.18Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8_1069)	Depositor
R, R_{free}	0.193 , 0.262 0.216 , 0.269	Depositor DCC
R_{free} test set	3492 reflections (5.63%)	DCC
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 38.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	3 of 130627 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	17293	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.00% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, KK1, NAG

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	0.45	0/1696	0.59	0/2310
1	B	0.45	0/1690	0.59	0/2302
1	C	0.60	1/1720 (0.1%)	0.66	3/2343 (0.1%)
1	D	0.44	0/1743	0.61	0/2377
1	E	0.41	0/1678	0.57	0/2284
1	F	0.40	0/1709	0.59	0/2327
1	G	0.41	0/1707	0.57	0/2324
1	H	0.40	0/1699	0.59	0/2313
1	I	0.41	0/1699	0.58	0/2313
1	J	0.48	1/1689 (0.1%)	0.61	0/2300
All	All	0.45	2/17030 (0.0%)	0.59	3/23193 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	26	PRO	N-CD	5.55	1.55	1.47
1	C	71	PRO	N-CD	5.05	1.54	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	70	SER	C-N-CD	5.74	140.46	128.40
1	C	60	ASP	CB-CG-OD1	5.32	123.08	118.30
1	C	161	ASP	CB-CG-OD1	-5.14	113.67	118.30

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	1662	0	1600	52	0
1	B	1656	0	1600	41	0
1	C	1682	0	1619	47	0
1	D	1706	0	1637	69	0
1	E	1645	0	1584	27	0
1	F	1675	0	1610	64	0
1	G	1672	0	1615	27	0
1	H	1664	0	1604	38	0
1	I	1664	0	1604	55	0
1	J	1655	0	1590	67	0
2	A	24	0	28	4	0
2	B	24	0	28	6	0
2	C	24	0	28	2	0
2	D	24	0	28	1	0
2	E	24	0	28	1	0
2	F	24	0	28	8	0
2	G	16	0	10	0	0
2	H	16	0	10	6	0
2	I	24	0	28	4	0
2	J	16	0	10	8	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
3	C	14	0	13	6	0
3	D	14	0	13	0	0
3	E	14	0	13	0	0
3	F	14	0	13	1	0
3	G	14	0	13	0	0
3	H	14	0	13	0	0
3	I	14	0	13	1	0
3	J	14	0	13	0	0
4	A	15	0	0	1	0
4	B	5	0	0	0	0
4	C	15	0	0	1	0
4	D	10	0	0	0	0
4	E	15	0	0	0	0
4	F	10	0	0	0	0
4	G	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	10	0	0	0	0
4	J	5	0	0	1	0
5	A	18	0	0	0	0
5	B	17	0	0	0	0
5	C	26	0	0	0	0
5	D	22	0	0	0	0
5	E	18	0	0	0	0
5	F	9	0	0	3	0
5	G	9	0	0	0	0
5	H	14	0	0	1	0
5	I	11	0	0	0	0
5	J	17	0	0	0	0
All	All	17293	0	16419	433	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 13.

All (433) 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:-3:ASP:HA	1:C:71:PRO:CG	1.75	1.16
1:C:-3:ASP:OD2	1:C:71:PRO:HB3	1.48	1.12
1:A:158:ASN:HB3	1:A:176:VAL:O	1.46	1.10
1:C:-3:ASP:CA	1:C:71:PRO:HG3	1.79	1.10
1:F:183:VAL:HG11	2:F:301:KK1:H4	1.33	1.10
1:C:66:ASN:HD22	3:C:302:NAG:H82	1.15	1.07
1:D:34:LYS:HZ3	1:D:160:ASP:HB2	1.15	1.06
1:C:66:ASN:ND2	3:C:302:NAG:H82	1.70	1.04
1:G:131:GLU:HG2	1:G:202:ARG:NH1	1.71	1.03
1:F:183:VAL:CG1	2:F:301:KK1:H4	1.88	1.02
1:D:34:LYS:HB3	1:D:161:ASP:HB2	1.43	0.99
1:J:-4:ASP:O	1:J:71:PRO:HG3	1.60	0.99
1:J:22:GLN:HE22	1:J:61:ARG:HD2	1.31	0.96
1:C:-3:ASP:HA	1:C:71:PRO:HG3	0.97	0.95
1:D:34:LYS:NZ	1:D:160:ASP:HB2	1.82	0.93
1:D:34:LYS:CB	1:D:161:ASP:HB2	1.99	0.91
1:A:-6:TYR:HB3	1:E:148:ARG:HH12	1.36	0.90
1:C:66:ASN:HD22	3:C:302:NAG:C8	1.85	0.90
1:D:34:LYS:HA	1:D:161:ASP:CB	2.02	0.89
1:D:37:ASN:HA	1:D:165:PHE:CE1	2.07	0.89
1:F:114:MET:H	2:J:301:KK1:H26	1.36	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:102:LEU:HD12	1:J:144:THR:CG2	2.04	0.88
1:B:160:ASP:HB2	1:B:163:GLU:OE1	1.74	0.88
1:D:35:PHE:N	1:D:161:ASP:HB3	1.90	0.87
1:D:68:SER:O	1:D:69:HIS:CG	2.28	0.87
1:F:123:CYS:CB	1:F:136:CYS:SG	2.63	0.86
1:F:99:THR:OG1	1:F:100:PRO:HD2	1.75	0.86
1:J:22:GLN:O	1:J:23:ARG:HD3	1.75	0.85
1:F:102:LEU:CB	1:J:144:THR:HG21	2.06	0.85
1:D:34:LYS:CA	1:D:161:ASP:HB2	2.07	0.85
1:C:55:GLN:HG3	1:C:114:MET:SD	2.17	0.84
1:D:35:PHE:H	1:D:161:ASP:HB3	1.42	0.84
1:H:22:GLN:HE22	1:H:60:ASP:HA	1.40	0.84
1:A:157:GLU:HB2	1:A:178:GLN:H	1.42	0.83
1:A:161:ASP:OD1	1:A:176:VAL:HG21	1.79	0.81
1:J:-3:ASP:HA	1:J:71:PRO:HB3	1.59	0.81
1:H:22:GLN:NE2	1:H:60:ASP:HA	1.94	0.81
1:D:34:LYS:HA	1:D:161:ASP:HB2	1.62	0.80
1:D:34:LYS:HD2	1:D:160:ASP:HB3	1.64	0.80
1:D:34:LYS:CA	1:D:161:ASP:CB	2.59	0.80
1:B:158:ASN:O	1:B:161:ASP:CG	2.21	0.79
1:F:159:SER:O	1:F:176:VAL:HG23	1.81	0.79
1:A:160:ASP:O	1:A:176:VAL:HB	1.82	0.79
1:F:102:LEU:HB2	1:J:144:THR:HG21	1.65	0.78
1:D:178:GLN:HG2	1:D:197:VAL:HG22	1.64	0.78
1:I:146:HIS:CD2	1:I:149:GLU:HG3	2.19	0.78
1:F:114:MET:N	2:J:301:KK1:H26	1.98	0.77
1:D:69:HIS:CE1	1:F:69:HIS:NE2	2.53	0.77
1:I:146:HIS:NE2	1:I:149:GLU:HG3	2.00	0.77
1:H:178:GLN:HG2	1:H:197:VAL:HG22	1.67	0.77
1:C:66:ASN:ND2	3:C:302:NAG:C8	2.47	0.77
1:I:37:ASN:ND2	1:I:166:SER:HB3	2.01	0.76
1:J:22:GLN:HE22	1:J:61:ARG:CD	2.00	0.75
1:D:158:ASN:HD21	1:D:178:GLN:HB2	1.49	0.74
1:I:178:GLN:HG2	1:I:197:VAL:HG22	1.69	0.74
1:C:158:ASN:O	1:C:159:SER:OG	2.06	0.73
2:B:301:KK1:H16	1:C:53:TRP:HZ3	1.53	0.72
1:A:157:GLU:HG3	1:A:178:GLN:O	1.90	0.72
1:D:34:LYS:HA	1:D:161:ASP:HB3	1.72	0.72
1:F:181:ASN:HD21	2:F:301:KK1:H1	1.55	0.72
1:I:183:VAL:HG21	2:I:301:KK1:H14	1.72	0.71
1:F:99:THR:HG23	5:F:403:HOH:O	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:-3:ASP:HA	1:H:71:PRO:HG3	1.72	0.69
1:D:20:PRO:HA	1:E:0:LYS:HE2	1.73	0.69
1:A:41:VAL:HG22	1:A:48:VAL:HG23	1.75	0.69
1:F:102:LEU:HD12	1:J:144:THR:HG21	1.74	0.69
1:I:148:ARG:HD2	1:J:-6:TYR:HE2	1.58	0.69
1:G:170:ARG:HH21	1:G:203:LYS:HE3	1.58	0.69
1:A:41:VAL:HG13	1:A:125:VAL:HG11	1.75	0.68
1:J:22:GLN:NE2	1:J:61:ARG:HD2	2.09	0.68
1:G:73:GLN:HG3	1:G:104:ARG:HH21	1.59	0.68
1:F:137:ARG:NH1	1:F:198:SER:OG	2.27	0.68
1:D:37:ASN:CA	1:D:165:PHE:CE1	2.77	0.67
1:J:123:CYS:CB	1:J:136:CYS:HG	2.07	0.67
2:A:301:KK1:H16	1:B:53:TRP:HH2	1.59	0.67
1:H:172:GLU:HB3	1:H:204:LYS:HE3	1.77	0.67
1:H:53:TRP:CD1	1:H:116:SER:HB2	2.29	0.67
1:A:160:ASP:HB3	1:A:161:ASP:CG	2.15	0.67
1:G:131:GLU:CG	1:G:202:ARG:NH1	2.54	0.66
1:D:34:LYS:CA	1:D:161:ASP:HB3	2.25	0.66
1:F:102:LEU:CD1	1:J:144:THR:HG21	2.25	0.65
1:D:57:THR:HG22	1:D:112:LEU:HB3	1.78	0.65
1:F:39:LEU:O	1:F:49:ASP:HB3	1.97	0.65
1:C:30:SER:HB3	1:C:155:THR:CG2	2.27	0.64
1:A:124:ASP:HB2	1:B:168:TYR:CE1	2.33	0.64
1:F:155:THR:HB	1:F:157:GLU:HG2	1.80	0.64
1:I:174:LEU:HD21	1:I:202:ARG:HG2	1.80	0.64
1:A:96:GLU:HG3	1:E:94:LYS:HE3	1.79	0.64
1:C:67:SER:HB2	1:C:107:SER:CB	2.28	0.64
1:I:147:SER:HB3	1:I:193:GLU:HG3	1.80	0.63
1:G:49:ASP:OD1	1:G:120:ARG:HG2	1.99	0.63
1:F:19:ILE:HG13	1:F:21:THR:HG23	1.81	0.63
1:C:-3:ASP:HA	1:C:71:PRO:CB	2.27	0.63
1:H:23:ARG:O	1:H:25:ARG:HD2	1.99	0.63
1:J:26:PRO:HB3	1:J:148:ARG:O	1.97	0.63
1:C:-3:ASP:CG	1:C:71:PRO:HB3	2.20	0.62
1:F:102:LEU:HD12	1:J:144:THR:HG23	1.81	0.62
1:B:88:ALA:HB3	1:B:91:ALA:HB2	1.80	0.62
1:A:158:ASN:CB	1:A:176:VAL:O	2.36	0.62
1:H:92:ILE:HD12	1:I:39:LEU:HD11	1.80	0.62
1:F:123:CYS:HG	1:F:136:CYS:HG	0.64	0.61
1:E:41:VAL:HG13	1:E:125:VAL:HG11	1.82	0.61
1:C:24:ASP:HA	1:D:-2:ASP:HB2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:34:LYS:HB3	1:D:161:ASP:CB	2.24	0.61
1:B:160:ASP:HA	1:B:163:GLU:HB2	1.82	0.61
1:D:68:SER:O	1:D:69:HIS:ND1	2.32	0.61
1:G:172:GLU:HB3	1:G:204:LYS:HE3	1.82	0.61
1:H:120:ARG:NH2	5:H:410:HOH:O	2.33	0.61
1:H:88:ALA:HB3	1:H:91:ALA:HB2	1.83	0.61
1:A:157:GLU:HB3	1:A:178:GLN:HB2	1.81	0.60
1:F:41:VAL:HG22	1:F:48:VAL:HG23	1.82	0.60
1:D:161:ASP:CG	1:D:164:TYR:HD2	2.04	0.60
1:A:157:GLU:CB	1:A:178:GLN:H	2.13	0.60
1:A:-6:TYR:HB3	1:E:148:ARG:NH1	2.13	0.60
1:B:139:LYS:NZ	2:B:301:KK1:H5	2.16	0.60
2:B:301:KK1:H13	1:C:53:TRP:HH2	1.65	0.60
1:B:152:VAL:HG12	1:B:195:VAL:HG23	1.84	0.60
1:D:34:LYS:HD2	1:D:160:ASP:CB	2.30	0.60
1:B:159:SER:O	1:B:160:ASP:OD1	2.19	0.60
1:F:-7:ASP:OD2	1:F:68:SER:O	2.20	0.59
1:J:123:CYS:SG	1:J:136:CYS:CB	2.91	0.59
1:C:170:ARG:NH2	4:C:305:PO4:O3	2.35	0.59
1:I:146:HIS:NE2	1:I:149:GLU:CG	2.66	0.59
1:G:34:LYS:HB2	1:G:53:TRP:HB2	1.85	0.59
1:F:178:GLN:HB3	1:F:197:VAL:HG22	1.85	0.59
1:F:124:ASP:HB2	1:G:168:TYR:CE2	2.37	0.58
1:I:-7:ASP:H3	1:I:-4:ASP:HB2	1.67	0.58
1:C:181:ASN:OD1	1:C:182:SER:N	2.35	0.58
1:B:24:ASP:OD1	1:C:-2:ASP:HB2	2.03	0.58
1:G:152:VAL:HG12	1:G:195:VAL:HG13	1.85	0.58
1:I:129:ASP:N	1:I:129:ASP:OD2	2.37	0.58
1:J:123:CYS:SG	1:J:136:CYS:HB2	2.44	0.57
1:B:183:VAL:HG11	2:B:301:KK1:H12	1.86	0.57
1:J:147:SER:HB3	1:J:193:GLU:HG3	1.86	0.57
1:H:123:CYS:SG	1:H:136:CYS:CB	2.91	0.57
1:E:152:VAL:HG23	1:E:195:VAL:HG23	1.85	0.57
1:B:-1:ASP:HB3	1:B:2:ASP:HB2	1.86	0.57
1:J:172:GLU:CG	1:J:204:LYS:HG2	2.35	0.57
1:D:34:LYS:C	1:D:161:ASP:HB3	2.25	0.56
1:A:46:ASN:OD1	1:A:125:VAL:HG23	2.05	0.56
2:I:301:KK1:H16	1:J:53:TRP:CH2	2.41	0.56
1:C:65:TRP:CZ2	1:C:72:ASP:O	2.58	0.56
1:H:22:GLN:NE2	1:H:59:SER:O	2.39	0.56
1:B:20:PRO:HA	1:C:0:LYS:HE3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:73:GLN:HG3	1:G:104:ARG:NH2	2.20	0.56
1:I:146:HIS:NE2	1:I:149:GLU:CD	2.59	0.56
1:B:139:LYS:HZ2	2:B:301:KK1:H5	1.71	0.56
1:D:34:LYS:CD	1:D:160:ASP:CB	2.84	0.55
1:I:146:HIS:NE2	1:I:149:GLU:OE2	2.39	0.55
1:G:165:PHE:CE1	1:G:173:ILE:HD11	2.41	0.55
1:B:172:GLU:OE2	1:B:202:ARG:NH1	2.39	0.55
1:A:40:GLU:OE1	1:A:120:ARG:NH2	2.39	0.55
2:A:301:KK1:H16	1:B:53:TRP:CH2	2.41	0.55
1:J:146:HIS:CD2	1:J:148:ARG:HB2	2.42	0.55
2:F:301:KK1:H26	1:G:114:MET:H	1.71	0.55
1:C:72:ASP:CG	1:C:73:GLN:HG2	2.27	0.55
1:J:49:ASP:OD1	1:J:120:ARG:HG3	2.06	0.55
1:F:102:LEU:O	1:F:113:TYR:HD1	1.89	0.55
1:B:178:GLN:HG3	1:B:197:VAL:HG22	1.88	0.55
1:H:123:CYS:SG	1:H:136:CYS:HB2	2.48	0.54
1:J:22:GLN:C	1:J:23:ARG:HD3	2.27	0.54
1:A:170:ARG:HD2	1:E:45:THR:HG22	1.88	0.54
1:D:143:TRP:CE2	1:E:99:THR:HG21	2.42	0.54
1:F:155:THR:HG1	1:F:157:GLU:N	2.05	0.54
2:H:301:KK1:H25	1:I:104:ARG:CG	2.38	0.54
1:I:41:VAL:HG21	1:I:201:PHE:HE2	1.72	0.54
1:A:168:TYR:CE1	1:E:124:ASP:HB2	2.43	0.54
3:C:302:NAG:C1	3:C:302:NAG:C8	2.85	0.54
1:J:3:ARG:NH1	4:J:303:PO4:O3	2.41	0.54
1:F:123:CYS:HB2	1:F:136:CYS:SG	2.48	0.54
1:A:139:LYS:NZ	2:A:301:KK1:H4	2.23	0.54
1:G:88:ALA:HB3	1:G:91:ALA:HB2	1.90	0.53
1:J:101:GLN:OE1	1:J:113:TYR:OH	2.26	0.53
1:F:174:LEU:HD21	1:F:202:ARG:HG2	1.90	0.53
1:A:174:LEU:HD21	1:A:202:ARG:HG2	1.91	0.53
1:A:154:PRO:HB3	1:A:195:VAL:HG11	1.89	0.53
1:I:37:ASN:HD21	1:I:169:SER:HB2	1.73	0.52
1:A:157:GLU:CG	1:A:178:GLN:O	2.56	0.52
1:I:146:HIS:CE1	1:I:149:GLU:OE2	2.62	0.52
1:A:20:PRO:HA	1:B:0:LYS:HE3	1.92	0.52
1:J:55:GLN:HG2	1:J:114:MET:SD	2.49	0.52
1:F:39:LEU:HD12	1:F:40:GLU:H	1.74	0.52
1:C:23:ARG:HB2	1:C:25:ARG:HD2	1.90	0.52
1:E:129:ASP:OD1	1:E:203:LYS:HD3	2.10	0.52
1:H:43:GLU:HA	1:H:125:VAL:HG23	1.89	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:112:LEU:O	2:J:301:KK1:H24	2.09	0.52
1:I:37:ASN:HD22	1:I:166:SER:HB3	1.74	0.52
1:G:101:GLN:OE1	1:G:113:TYR:OH	2.27	0.52
1:J:172:GLU:HG2	1:J:204:LYS:HG2	1.90	0.52
1:H:175:ASP:OD1	1:H:176:VAL:N	2.41	0.52
1:A:190:GLU:N	1:A:190:GLU:OE1	2.42	0.52
1:B:159:SER:O	1:B:160:ASP:CB	2.58	0.52
1:I:73:GLN:HG3	1:I:104:ARG:HH21	1.75	0.52
1:F:145:HIS:HB3	1:F:149:GLU:HB2	1.92	0.51
1:F:183:VAL:HG11	2:F:301:KK1:C02	2.23	0.51
1:J:22:GLN:NE2	1:J:61:ARG:CD	2.69	0.51
1:F:67:SER:O	1:F:70:SER:O	2.28	0.51
1:C:89:TYR:OH	2:C:301:KK1:N15	2.41	0.51
1:F:159:SER:O	1:F:176:VAL:CG2	2.54	0.51
1:B:158:ASN:HD21	1:B:178:GLN:HB3	1.75	0.51
1:J:104:ARG:HB2	1:J:112:LEU:HD12	1.93	0.51
1:D:34:LYS:CD	1:D:160:ASP:HB2	2.40	0.51
1:F:114:MET:H	2:J:301:KK1:C22	2.16	0.51
1:I:-6:TYR:HD2	1:I:-6:TYR:H	1.58	0.51
1:F:181:ASN:ND2	2:F:301:KK1:H1	2.23	0.51
1:D:55:GLN:HA	1:D:114:MET:HG3	1.92	0.51
1:I:101:GLN:OE1	1:I:113:TYR:OH	2.27	0.51
1:A:131:GLU:O	1:A:202:ARG:NH1	2.37	0.51
1:I:55:GLN:HG2	1:I:114:MET:SD	2.50	0.51
1:J:152:VAL:HG12	1:J:195:VAL:HG23	1.93	0.51
1:B:73:GLN:HG3	1:B:104:ARG:HH21	1.76	0.51
1:F:172:GLU:HB3	1:F:204:LYS:HD2	1.92	0.51
1:E:41:VAL:HG22	1:E:48:VAL:HG22	1.92	0.51
1:J:48:VAL:HG22	1:J:50:VAL:HG13	1.92	0.50
1:E:-1:ASP:HB3	1:E:2:ASP:HB2	1.93	0.50
1:D:161:ASP:OD2	1:D:164:TYR:HD2	1.95	0.50
1:J:73:GLN:HB3	1:J:106:VAL:HA	1.93	0.50
1:D:94:LYS:HE3	1:E:96:GLU:HG3	1.92	0.50
1:D:68:SER:C	1:D:69:HIS:CG	2.85	0.50
1:F:102:LEU:CD1	1:J:144:THR:OG1	2.59	0.50
1:G:157:GLU:O	1:G:177:THR:HA	2.13	0.49
1:A:160:ASP:HB3	1:A:161:ASP:CB	2.42	0.49
1:F:102:LEU:O	1:F:113:TYR:CD1	2.65	0.49
1:I:146:HIS:CD2	1:I:149:GLU:CG	2.94	0.49
1:H:143:TRP:CE2	1:I:99:THR:HG21	2.47	0.49
1:G:23:ARG:HB3	1:G:25:ARG:HH12	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:22:GLN:OE1	1:E:61:ARG:HG3	2.13	0.49
1:I:147:SER:HA	1:I:150:ILE:O	2.13	0.49
1:H:6:ILE:HD13	1:H:65:TRP:CZ2	2.48	0.49
1:J:147:SER:HA	1:J:150:ILE:O	2.13	0.49
4:A:303:PO4:O2	1:B:3:ARG:NH2	2.37	0.49
1:B:-5:LYS:HB3	1:B:-4:ASP:H	1.52	0.49
1:D:137:ARG:NH1	1:D:196:GLU:OE2	2.45	0.48
1:G:131:GLU:HG2	1:G:202:ARG:HH12	1.69	0.48
1:H:53:TRP:NE1	1:H:116:SER:HB2	2.27	0.48
1:D:203:LYS:HE3	1:D:204:LYS:O	2.13	0.48
1:G:131:GLU:HG2	1:G:202:ARG:HH11	1.67	0.48
1:D:69:HIS:CE1	1:F:69:HIS:CD2	3.01	0.48
1:D:30:SER:HB3	1:D:155:THR:HG22	1.94	0.48
1:A:157:GLU:HA	1:A:158:ASN:HA	1.65	0.48
1:J:-3:ASP:HA	1:J:71:PRO:CB	2.37	0.48
1:H:-1:ASP:HB3	1:H:2:ASP:HB2	1.94	0.48
1:D:124:ASP:OD1	1:D:126:SER:OG	2.31	0.48
1:C:148:ARG:HG3	1:C:148:ARG:O	2.13	0.48
1:F:113:TYR:HA	2:J:301:KK1:H24	1.94	0.48
1:E:170:ARG:O	1:E:204:LYS:HB2	2.14	0.48
1:C:-3:ASP:C	1:C:71:PRO:HG3	2.34	0.48
1:A:160:ASP:HB3	1:A:161:ASP:OD2	2.14	0.48
1:F:183:VAL:HG12	2:F:301:KK1:H4	1.91	0.47
1:B:158:ASN:ND2	1:B:178:GLN:HB3	2.28	0.47
1:F:82:TRP:O	5:F:407:HOH:O	2.20	0.47
1:D:161:ASP:OD2	1:D:164:TYR:HB2	2.13	0.47
1:F:39:LEU:HD11	1:J:45:THR:HB	1.97	0.47
1:H:203:LYS:HG2	1:H:204:LYS:H	1.80	0.47
1:H:152:VAL:HG21	1:H:194:ASP:HA	1.96	0.47
1:H:20:PRO:HA	1:I:0:LYS:HE3	1.96	0.47
1:I:41:VAL:HG13	1:I:125:VAL:HG11	1.97	0.47
1:C:158:ASN:ND2	1:C:159:SER:N	2.63	0.47
1:B:17:ASP:OD2	1:C:11:ARG:NH2	2.39	0.47
1:H:144:THR:HG21	1:I:102:LEU:HB2	1.97	0.47
1:B:57:THR:HG22	1:B:112:LEU:HB3	1.96	0.46
1:J:143:TRP:O	2:J:301:KK1:C24	2.63	0.46
1:I:44:ILE:HG22	1:J:170:ARG:HH11	1.81	0.46
1:D:165:PHE:CE2	1:D:169:SER:HB3	2.50	0.46
1:D:86:LEU:HA	1:D:141:GLY:O	2.15	0.46
1:I:124:ASP:HB2	1:J:168:TYR:CE1	2.50	0.46
1:D:160:ASP:HA	1:D:161:ASP:HA	1.62	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:ASN:OD1	3:C:302:NAG:O5	2.33	0.46
1:H:98:LEU:HD12	1:H:116:SER:OG	2.15	0.46
1:A:112:LEU:O	2:E:301:KK1:H24	2.15	0.46
1:B:158:ASN:OD1	1:B:178:GLN:HB2	2.15	0.46
1:F:152:VAL:O	1:F:180:LYS:HD3	2.16	0.46
1:I:37:ASN:ND2	1:I:169:SER:HB2	2.30	0.46
2:H:301:KK1:H24	1:I:112:LEU:O	2.15	0.46
1:B:158:ASN:O	1:B:161:ASP:OD1	2.34	0.46
1:I:22:GLN:OE1	1:I:61:ARG:HB2	2.16	0.46
1:H:10:ILE:O	1:H:14:SER:HB3	2.16	0.46
1:A:78:ILE:HD11	1:A:103:ALA:HB2	1.98	0.46
1:D:165:PHE:HE2	1:D:169:SER:HB3	1.81	0.46
1:A:73:GLN:HG2	1:A:106:VAL:HG13	1.98	0.46
1:F:24:ASP:HA	1:G:-2:ASP:HB2	1.97	0.46
1:A:160:ASP:CB	1:A:161:ASP:CG	2.85	0.45
1:F:41:VAL:HG13	1:F:125:VAL:HG11	1.98	0.45
1:H:41:VAL:HG22	1:H:48:VAL:HG12	1.97	0.45
1:D:-1:ASP:HB3	1:D:2:ASP:HB2	1.98	0.45
1:I:24:ASP:HA	1:J:-2:ASP:HB2	1.98	0.45
1:J:23:ARG:O	1:J:24:ASP:HB2	2.16	0.45
1:C:30:SER:HB3	1:C:155:THR:HG22	1.99	0.45
1:H:91:ALA:HB1	1:H:119:GLN:NE2	2.32	0.45
1:D:149:GLU:OE2	1:E:3:ARG:NH2	2.48	0.45
1:D:184:THR:HG22	1:D:185:TYR:H	1.81	0.45
1:B:174:LEU:HD21	1:B:202:ARG:HG2	1.99	0.45
1:I:41:VAL:HG21	1:I:201:PHE:CE2	2.50	0.45
1:I:54:GLN:O	1:I:114:MET:HA	2.17	0.45
1:D:34:LYS:NZ	1:D:160:ASP:CB	2.67	0.45
1:A:155:THR:HA	1:A:156:THR:HA	1.64	0.45
1:C:144:THR:HB	1:D:104:ARG:HD3	1.98	0.45
1:D:123:CYS:O	1:D:125:VAL:HG23	2.17	0.45
1:J:146:HIS:NE2	1:J:148:ARG:HB2	2.32	0.45
1:I:73:GLN:HG3	1:I:104:ARG:NH2	2.31	0.45
1:H:111:VAL:C	1:H:112:LEU:HD23	2.37	0.45
1:J:178:GLN:HA	1:J:196:GLU:O	2.17	0.45
1:I:66:ASN:ND2	3:I:302:NAG:O7	2.50	0.45
1:F:183:VAL:HG11	2:F:301:KK1:H8	1.99	0.45
1:D:33:LEU:HD21	1:D:140:ILE:HD12	1.98	0.45
1:J:10:ILE:O	1:J:14:SER:HB3	2.16	0.45
5:F:405:HOH:O	2:J:301:KK1:H25	2.18	0.44
1:B:55:GLN:HA	1:B:114:MET:HG3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:143:TRP:CE2	2:C:301:KK1:H28	2.52	0.44
1:J:40:GLU:O	1:J:41:VAL:HG23	2.18	0.44
1:A:24:ASP:HA	1:B:-2:ASP:HB3	1.99	0.44
1:I:30:SER:HB2	1:I:57:THR:OG1	2.17	0.44
1:F:48:VAL:HG22	1:F:49:ASP:H	1.83	0.44
1:A:149:GLU:CD	1:B:104:ARG:HH22	2.21	0.44
1:G:158:ASN:ND2	1:G:177:THR:HG22	2.32	0.44
1:A:154:PRO:HG3	1:A:195:VAL:HG22	1.99	0.44
1:I:139:LYS:HG3	1:I:196:GLU:HG2	1.99	0.44
1:C:67:SER:HB2	1:C:107:SER:OG	2.16	0.44
1:D:185:TYR:CE1	2:D:301:KK1:H14	2.53	0.44
1:I:36:ILE:HD13	1:I:36:ILE:N	2.33	0.44
1:A:139:LYS:HZ1	2:A:301:KK1:H4	1.82	0.44
1:H:143:TRP:O	2:H:301:KK1:N13	2.51	0.43
1:J:32:SER:HB3	1:J:155:THR:HG22	1.99	0.43
1:B:124:ASP:HB2	1:C:168:TYR:CE1	2.53	0.43
1:C:72:ASP:OD2	1:C:73:GLN:HG2	2.18	0.43
1:E:67:SER:HA	1:E:70:SER:OG	2.18	0.43
1:E:30:SER:O	1:E:56:THR:HA	2.19	0.43
1:D:24:ASP:HA	1:E:-2:ASP:HB2	2.01	0.43
1:J:123:CYS:CB	1:J:136:CYS:SG	3.01	0.43
1:J:43:GLU:HA	1:J:125:VAL:HG12	2.00	0.43
1:E:89:TYR:HD2	1:E:139:LYS:O	2.01	0.43
1:G:13:THR:HG21	1:G:62:THR:O	2.18	0.43
1:D:34:LYS:CD	1:D:160:ASP:HB3	2.39	0.43
1:J:25:ARG:HG3	1:J:25:ARG:NH1	2.33	0.43
1:F:102:LEU:HB2	1:J:144:THR:CG2	2.44	0.43
1:F:102:LEU:HB3	1:J:144:THR:HG21	1.92	0.43
1:I:-7:ASP:HB2	1:I:-6:TYR:H	1.65	0.43
1:C:144:THR:HG21	1:D:102:LEU:HB2	2.00	0.43
1:F:102:LEU:CD1	1:J:144:THR:CG2	2.81	0.43
1:B:-4:ASP:O	1:B:-2:ASP:N	2.51	0.43
1:D:14:SER:O	1:D:16:PRO:HD3	2.18	0.43
1:J:23:ARG:O	1:J:24:ASP:CB	2.66	0.43
1:A:154:PRO:HB3	1:A:195:VAL:HG21	2.00	0.43
1:C:17:ASP:OD2	1:D:11:ARG:NH2	2.35	0.43
1:I:-5:LYS:NZ	1:I:-5:LYS:HB2	2.33	0.43
1:D:85:ASP:N	1:D:85:ASP:OD1	2.47	0.43
1:D:68:SER:C	1:D:69:HIS:ND1	2.71	0.43
1:I:-6:TYR:CD2	1:I:-6:TYR:N	2.87	0.43
1:B:33:LEU:HD22	1:B:52:PHE:CD1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:177:THR:HG23	1:J:198:SER:HB2	2.00	0.43
1:A:99:THR:HG21	1:E:143:TRP:CE2	2.53	0.43
1:B:101:GLN:OE1	1:B:113:TYR:OH	2.36	0.43
1:D:65:TRP:CD1	1:D:66:ASN:O	2.72	0.43
1:F:54:GLN:O	1:F:114:MET:HA	2.18	0.43
1:I:33:LEU:O	1:I:161:ASP:HB3	2.19	0.43
2:I:301:KK1:H25	1:J:104:ARG:HG2	1.99	0.42
1:A:174:LEU:HD12	1:A:200:ASN:OD1	2.18	0.42
1:J:19:ILE:HG13	1:J:21:THR:HG23	2.00	0.42
1:C:20:PRO:HA	1:D:0:LYS:HE3	2.00	0.42
1:H:123:CYS:CB	1:H:136:CYS:SG	3.06	0.42
1:F:75:SER:HB3	1:F:102:LEU:HD13	2.01	0.42
2:I:301:KK1:H16	1:J:53:TRP:CZ3	2.55	0.42
1:A:154:PRO:HB3	1:A:195:VAL:CG1	2.48	0.42
1:D:165:PHE:CE2	1:D:169:SER:CB	3.02	0.42
1:H:133:GLY:HA2	1:H:202:ARG:HB3	2.01	0.42
1:A:160:ASP:HA	1:A:161:ASP:HA	1.80	0.42
1:F:157:GLU:HB3	1:F:158:ASN:H	1.60	0.42
1:E:-7:ASP:N	1:E:72:ASP:OD1	2.49	0.42
1:C:158:ASN:ND2	1:C:158:ASN:C	2.73	0.42
1:A:24:ASP:HA	1:B:-2:ASP:CB	2.49	0.42
1:I:124:ASP:HB2	1:J:168:TYR:CD1	2.54	0.42
1:E:101:GLN:NE2	1:E:113:TYR:OH	2.51	0.42
1:D:88:ALA:HB3	1:D:91:ALA:HB2	2.02	0.42
1:J:36:ILE:HD13	1:J:164:TYR:HB3	2.01	0.42
2:H:301:KK1:H25	1:I:104:ARG:HG2	2.00	0.42
1:F:92:ILE:O	1:G:118:ARG:HD2	2.20	0.42
1:D:37:ASN:C	1:D:165:PHE:HE1	2.23	0.42
1:I:125:VAL:O	1:I:128:VAL:HG12	2.20	0.42
1:G:158:ASN:O	1:G:159:SER:HB2	2.20	0.42
1:J:41:VAL:HG21	1:J:201:PHE:HE2	1.85	0.42
1:G:31:VAL:O	1:G:154:PRO:HA	2.20	0.42
1:A:-7:ASP:OD2	1:A:-7:ASP:N	2.35	0.42
1:I:37:ASN:HD21	1:I:166:SER:HB3	1.82	0.42
1:C:147:SER:CB	1:C:193:GLU:HG3	2.50	0.42
1:J:143:TRP:O	2:J:301:KK1:H28	2.20	0.41
1:J:106:VAL:HG23	1:J:110:GLU:O	2.20	0.41
1:E:47:GLU:HB3	1:E:120:ARG:NH2	2.34	0.41
1:F:112:LEU:N	1:F:112:LEU:HD12	2.35	0.41
1:B:158:ASN:OD1	1:B:178:GLN:CB	2.68	0.41
1:B:144:THR:HG21	1:C:102:LEU:HB2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:57:THR:HG22	1:J:112:LEU:HD23	2.00	0.41
1:D:143:TRP:CZ2	1:E:99:THR:HG21	2.55	0.41
1:I:88:ALA:HB3	1:I:91:ALA:HB2	2.02	0.41
1:F:55:GLN:HA	1:F:114:MET:HG3	2.01	0.41
1:A:139:LYS:HG3	1:A:196:GLU:HG3	2.02	0.41
1:C:147:SER:N	1:C:191:ALA:O	2.48	0.41
1:G:45:THR:HG22	1:H:170:ARG:NH1	2.35	0.41
1:J:173:ILE:HG12	1:J:199:LEU:HD11	2.03	0.41
1:A:157:GLU:HB2	1:A:158:ASN:HD22	1.84	0.41
2:B:301:KK1:H13	1:C:53:TRP:CH2	2.52	0.41
1:A:152:VAL:O	1:A:180:LYS:NZ	2.54	0.41
1:C:88:ALA:HB2	1:C:140:ILE:HD13	2.02	0.41
1:E:88:ALA:HB3	1:E:91:ALA:HB2	2.01	0.41
3:F:302:NAG:H2	3:F:302:NAG:H82	1.98	0.41
1:C:67:SER:O	1:C:67:SER:OG	2.36	0.41
1:A:149:GLU:OE2	1:B:104:ARG:NH2	2.52	0.41
1:D:88:ALA:HA	1:D:139:LYS:O	2.20	0.41
1:H:7:LEU:HA	1:H:7:LEU:HD12	1.73	0.41
1:F:102:LEU:HA	1:F:102:LEU:HD23	1.92	0.41
1:H:10:ILE:HG23	1:H:63:LEU:HD22	2.03	0.41
1:D:185:TYR:HA	1:D:185:TYR:HD2	1.65	0.41
1:A:32:SER:HB2	1:A:155:THR:O	2.21	0.41
1:H:81:LEU:HD23	1:H:81:LEU:HA	1.88	0.41
1:G:143:TRP:CE2	1:H:99:THR:HG21	2.56	0.41
1:D:69:HIS:ND1	1:F:69:HIS:CD2	2.89	0.41
1:F:65:TRP:CZ3	1:F:107:SER:HA	2.56	0.41
1:I:6:ILE:HG22	1:I:10:ILE:HD12	2.02	0.41
1:D:160:ASP:HB3	1:D:161:ASP:HA	2.03	0.40
1:E:36:ILE:HB	1:E:51:VAL:HG12	2.02	0.40
1:F:48:VAL:HG12	1:F:121:PHE:HB2	2.03	0.40
2:H:301:KK1:H25	1:I:104:ARG:HG3	2.03	0.40
1:B:73:GLN:HG3	1:B:104:ARG:NH2	2.36	0.40
1:C:30:SER:HA	1:C:153:ASP:O	2.21	0.40
1:H:89:TYR:HE1	2:H:301:KK1:N16	2.19	0.40
1:J:78:ILE:HD11	1:J:101:GLN:HG3	2.04	0.40
1:A:152:VAL:CG1	1:A:195:VAL:HG23	2.51	0.40
1:J:178:GLN:HB3	1:J:197:VAL:HG22	2.04	0.40
1:I:162:SER:HB3	1:I:173:ILE:HG21	2.04	0.40
1:G:19:ILE:HG22	1:G:82:TRP:CZ2	2.57	0.40
1:H:180:LYS:H	1:H:180:LYS:HG2	1.43	0.40
1:F:102:LEU:HD13	1:J:144:THR:OG1	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:ILE:HA	1:A:78:ILE:HD13	1.72	0.40
1:I:92:ILE:HD11	1:I:120:ARG:HG2	2.04	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	203/217 (94%)	201 (99%)	1 (0%)	1 (0%)	34	63
1	B	203/217 (94%)	200 (98%)	3 (2%)	0	100	100
1	C	206/217 (95%)	202 (98%)	4 (2%)	0	100	100
1	D	211/217 (97%)	208 (99%)	3 (1%)	0	100	100
1	E	199/217 (92%)	198 (100%)	1 (0%)	0	100	100
1	F	203/217 (94%)	195 (96%)	8 (4%)	0	100	100
1	G	202/217 (93%)	198 (98%)	4 (2%)	0	100	100
1	H	201/217 (93%)	193 (96%)	7 (4%)	1 (0%)	34	63
1	I	201/217 (93%)	195 (97%)	6 (3%)	0	100	100
1	J	202/217 (93%)	193 (96%)	8 (4%)	1 (0%)	34	63
All	All	2031/2170 (94%)	1983 (98%)	45 (2%)	3 (0%)	56	83

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	PRO
1	J	154	PRO
1	H	154	PRO

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	193/203 (95%)	181 (94%)	12 (6%)	23	49
1	B	193/203 (95%)	179 (93%)	14 (7%)	17	39
1	C	196/203 (97%)	180 (92%)	16 (8%)	14	32
1	D	199/203 (98%)	182 (92%)	17 (8%)	13	30
1	E	191/203 (94%)	177 (93%)	14 (7%)	17	39
1	F	195/203 (96%)	172 (88%)	23 (12%)	6	15
1	G	194/203 (96%)	184 (95%)	10 (5%)	29	58
1	H	193/203 (95%)	173 (90%)	20 (10%)	9	20
1	I	193/203 (95%)	174 (90%)	19 (10%)	10	23
1	J	192/203 (95%)	173 (90%)	19 (10%)	10	22
All	All	1939/2030 (96%)	1775 (92%)	164 (8%)	13	30

All (164) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-3	ASP
1	A	14	SER
1	A	25	ARG
1	A	112	LEU
1	A	122	SER
1	A	125	VAL
1	A	129	ASP
1	A	156	THR
1	A	157	GLU
1	A	158	ASN
1	A	159	SER
1	A	161	ASP
1	B	-3	ASP
1	B	14	SER
1	B	31	VAL
1	B	59	SER

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Mol	Chain	Res	Type
1	B	68	SER
1	B	79	SER
1	B	112	LEU
1	B	136	CYS
1	B	160	ASP
1	B	167	GLN
1	B	178	GLN
1	B	182	SER
1	B	184	THR
1	B	202	ARG
1	C	-4	ASP
1	C	-3	ASP
1	C	7	LEU
1	C	13	THR
1	C	21	THR
1	C	24	ASP
1	C	59	SER
1	C	67	SER
1	C	73	GLN
1	C	148	ARG
1	C	153	ASP
1	C	158	ASN
1	C	160	ASP
1	C	161	ASP
1	C	162	SER
1	C	196	GLU
1	D	24	ASP
1	D	48	VAL
1	D	59	SER
1	D	67	SER
1	D	68	SER
1	D	112	LEU
1	D	123	CYS
1	D	126	SER
1	D	129	ASP
1	D	136	CYS
1	D	158	ASN
1	D	159	SER
1	D	160	ASP
1	D	183	VAL
1	D	185	TYR
1	D	187	CYS

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Mol	Chain	Res	Type
1	D	195	VAL
1	E	13	THR
1	E	14	SER
1	E	24	ASP
1	E	32	SER
1	E	48	VAL
1	E	50	VAL
1	E	61	ARG
1	E	99	THR
1	E	112	LEU
1	E	132	SER
1	E	151	SER
1	E	152	VAL
1	E	170	ARG
1	E	183	VAL
1	F	-7	ASP
1	F	-6	TYR
1	F	-3	ASP
1	F	2	ASP
1	F	24	ASP
1	F	31	VAL
1	F	39	LEU
1	F	56	THR
1	F	68	SER
1	F	72	ASP
1	F	83	VAL
1	F	118	ARG
1	F	122	SER
1	F	149	GLU
1	F	155	THR
1	F	163	GLU
1	F	167	GLN
1	F	170	ARG
1	F	176	VAL
1	F	178	GLN
1	F	179	LYS
1	F	180	LYS
1	F	194	ASP
1	G	-7	ASP
1	G	32	SER
1	G	56	THR
1	G	59	SER

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Mol	Chain	Res	Type
1	G	110	GLU
1	G	129	ASP
1	G	138	ILE
1	G	155	THR
1	G	157	GLU
1	G	184	THR
1	H	14	SER
1	H	18	VAL
1	H	22	GLN
1	H	23	ARG
1	H	24	ASP
1	H	31	VAL
1	H	48	VAL
1	H	59	SER
1	H	61	ARG
1	H	112	LEU
1	H	116	SER
1	H	128	VAL
1	H	136	CYS
1	H	155	THR
1	H	164	TYR
1	H	180	LYS
1	H	185	TYR
1	H	190	GLU
1	H	194	ASP
1	H	196	GLU
1	I	-7	ASP
1	I	-5	LYS
1	I	37	ASN
1	I	39	LEU
1	I	41	VAL
1	I	49	ASP
1	I	99	THR
1	I	110	GLU
1	I	112	LEU
1	I	122	SER
1	I	128	VAL
1	I	129	ASP
1	I	148	ARG
1	I	152	VAL
1	I	163	GLU
1	I	170	ARG

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Mol	Chain	Res	Type
1	I	176	VAL
1	I	200	ASN
1	I	202	ARG
1	J	-5	LYS
1	J	13	THR
1	J	14	SER
1	J	23	ARG
1	J	25	ARG
1	J	31	VAL
1	J	56	THR
1	J	72	ASP
1	J	73	GLN
1	J	112	LEU
1	J	129	ASP
1	J	137	ARG
1	J	157	GLU
1	J	159	SER
1	J	163	GLU
1	J	170	ARG
1	J	177	THR
1	J	178	GLN
1	J	190	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	C	158	ASN
1	D	158	ASN
1	E	181	ASN
1	F	181	ASN
1	G	9	ASN
1	G	158	ASN
1	G	200	ASN
1	I	37	ASN
1	J	146	HIS

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 ⓘ

39 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	KK1	A	301	-	25,25,25	3.06	11 (44%)	29,31,31	1.98	7 (24%)
3	NAG	A	302	1	14,14,15	0.44	0	15,19,21	1.18	2 (13%)
4	PO4	A	303	-	4,4,4	0.58	0	6,6,6	0.29	0
4	PO4	A	304	-	4,4,4	0.48	0	6,6,6	0.29	0
4	PO4	A	305	-	4,4,4	0.27	0	6,6,6	0.28	0
2	KK1	B	301	-	25,25,25	3.30	10 (40%)	29,31,31	2.15	8 (27%)
3	NAG	B	302	1	14,14,15	0.56	0	15,19,21	0.87	0
4	PO4	B	303	-	4,4,4	0.33	0	6,6,6	0.27	0
2	KK1	C	301	-	25,25,25	3.59	12 (48%)	29,31,31	2.03	8 (27%)
3	NAG	C	302	1	14,14,15	1.46	3 (21%)	15,19,21	2.70	8 (53%)
4	PO4	C	303	-	4,4,4	0.69	0	6,6,6	0.26	0
4	PO4	C	304	-	4,4,4	0.58	0	6,6,6	0.27	0
4	PO4	C	305	-	4,4,4	0.33	0	6,6,6	0.27	0
2	KK1	D	301	-	25,25,25	3.57	9 (36%)	29,31,31	2.17	9 (31%)
3	NAG	D	302	1	14,14,15	0.53	0	15,19,21	1.11	2 (13%)
4	PO4	D	303	-	4,4,4	0.38	0	6,6,6	0.29	0
4	PO4	D	304	-	4,4,4	0.48	0	6,6,6	0.27	0
2	KK1	E	301	-	25,25,25	3.56	12 (48%)	29,31,31	1.80	9 (31%)
3	NAG	E	302	1	14,14,15	0.44	0	15,19,21	1.06	1 (6%)
4	PO4	E	303	-	4,4,4	0.69	0	6,6,6	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PO4	E	304	-	4,4,4	0.30	0	6,6,6	0.26	0
4	PO4	E	305	-	4,4,4	0.45	0	6,6,6	0.30	0
2	KK1	F	301	-	25,25,25	3.37	11 (44%)	29,31,31	1.79	9 (31%)
3	NAG	F	302	1	14,14,15	0.48	0	15,19,21	1.07	0
4	PO4	F	303	-	4,4,4	0.55	0	6,6,6	0.28	0
4	PO4	F	304	-	4,4,4	0.29	0	6,6,6	0.29	0
2	KK1	G	301	-	17,17,25	3.47	10 (58%)	21,23,31	1.89	7 (33%)
3	NAG	G	302	1	14,14,15	0.43	0	15,19,21	0.99	1 (6%)
4	PO4	G	303	-	4,4,4	0.50	0	6,6,6	0.28	0
4	PO4	G	304	-	4,4,4	0.34	0	6,6,6	0.28	0
2	KK1	H	301	-	17,17,25	3.46	10 (58%)	21,23,31	1.89	7 (33%)
3	NAG	H	302	1	14,14,15	0.54	0	15,19,21	0.60	0
4	PO4	H	303	-	4,4,4	0.65	0	6,6,6	0.27	0
4	PO4	H	304	-	4,4,4	0.69	0	6,6,6	0.31	0
2	KK1	I	301	-	25,25,25	3.51	12 (48%)	29,31,31	1.81	9 (31%)
3	NAG	I	302	1	14,14,15	1.03	1 (7%)	15,19,21	3.24	8 (53%)
2	KK1	J	301	-	17,17,25	3.46	10 (58%)	21,23,31	1.89	7 (33%)
3	NAG	J	302	1	14,14,15	0.44	0	15,19,21	1.70	2 (13%)
4	PO4	J	303	-	4,4,4	0.60	0	6,6,6	0.29	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	KK1	A	301	-	-	0/15/15/15	0/2/2/2
3	NAG	A	302	1	-	0/6/23/26	0/1/1/1
4	PO4	A	303	-	-	0/0/0/0	0/0/0/0
4	PO4	A	304	-	-	0/0/0/0	0/0/0/0
4	PO4	A	305	-	-	0/0/0/0	0/0/0/0
2	KK1	B	301	-	-	0/15/15/15	0/2/2/2
3	NAG	B	302	1	-	0/6/23/26	0/1/1/1
4	PO4	B	303	-	-	0/0/0/0	0/0/0/0
2	KK1	C	301	-	-	0/15/15/15	0/2/2/2
3	NAG	C	302	1	-	0/6/23/26	0/1/1/1
4	PO4	C	303	-	-	0/0/0/0	0/0/0/0
4	PO4	C	304	-	-	0/0/0/0	0/0/0/0
4	PO4	C	305	-	-	0/0/0/0	0/0/0/0
2	KK1	D	301	-	-	0/15/15/15	0/2/2/2
3	NAG	D	302	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PO4	D	303	-	-	0/0/0/0	0/0/0/0
4	PO4	D	304	-	-	0/0/0/0	0/0/0/0
2	KK1	E	301	-	-	0/15/15/15	0/2/2/2
3	NAG	E	302	1	-	0/6/23/26	0/1/1/1
4	PO4	E	303	-	-	0/0/0/0	0/0/0/0
4	PO4	E	304	-	-	0/0/0/0	0/0/0/0
4	PO4	E	305	-	-	0/0/0/0	0/0/0/0
2	KK1	F	301	-	-	0/15/15/15	0/2/2/2
3	NAG	F	302	1	-	0/6/23/26	0/1/1/1
4	PO4	F	303	-	-	0/0/0/0	0/0/0/0
4	PO4	F	304	-	-	0/0/0/0	0/0/0/0
2	KK1	G	301	-	-	0/6/6/15	0/2/2/2
3	NAG	G	302	1	-	0/6/23/26	0/1/1/1
4	PO4	G	303	-	-	0/0/0/0	0/0/0/0
4	PO4	G	304	-	-	0/0/0/0	0/0/0/0
2	KK1	H	301	-	-	0/6/6/15	0/2/2/2
3	NAG	H	302	1	-	0/6/23/26	0/1/1/1
4	PO4	H	303	-	-	0/0/0/0	0/0/0/0
4	PO4	H	304	-	-	0/0/0/0	0/0/0/0
2	KK1	I	301	-	-	0/15/15/15	0/2/2/2
3	NAG	I	302	1	-	0/6/23/26	0/1/1/1
2	KK1	J	301	-	-	0/6/6/15	0/2/2/2
3	NAG	J	302	1	-	0/6/23/26	0/1/1/1
4	PO4	J	303	-	-	0/0/0/0	0/0/0/0

All (111) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	302	NAG	O7-C7	-2.37	1.17	1.23
3	C	302	NAG	O5-C1	-2.31	1.39	1.43
3	I	302	NAG	C4-C5	-2.26	1.48	1.53
3	C	302	NAG	C3-C2	-2.02	1.47	1.52
2	F	301	KK1	C14-N16	2.04	1.39	1.35
2	C	301	KK1	C10-N16	2.06	1.38	1.34
2	E	301	KK1	C07-C08	2.15	1.60	1.51
2	E	301	KK1	C12-N13	2.16	1.37	1.34
2	C	301	KK1	C06-C07	2.21	1.64	1.51
2	G	301	KK1	C12-N13	2.33	1.38	1.34
2	I	301	KK1	C12-N13	2.35	1.38	1.34
2	H	301	KK1	C12-N13	2.36	1.38	1.34
2	E	301	KK1	C14-N16	2.40	1.39	1.35
2	J	301	KK1	C12-N13	2.42	1.38	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	KK1	C07-C08	2.53	1.62	1.51
2	A	301	KK1	C07-C08	2.54	1.62	1.51
2	D	301	KK1	C07-C08	2.56	1.62	1.51
2	F	301	KK1	C07-C08	2.56	1.62	1.51
2	A	301	KK1	C14-N16	2.57	1.40	1.35
2	C	301	KK1	C14-N16	2.58	1.40	1.35
2	F	301	KK1	C10-N16	2.64	1.39	1.34
2	A	301	KK1	C10-N16	2.67	1.39	1.34
2	E	301	KK1	C10-N16	2.73	1.39	1.34
2	B	301	KK1	C08-N09	2.75	1.52	1.46
2	E	301	KK1	C08-N09	2.82	1.52	1.46
2	G	301	KK1	C14-N16	2.83	1.40	1.35
2	H	301	KK1	C14-N16	2.83	1.40	1.35
2	D	301	KK1	C24-C23	2.84	1.43	1.38
2	J	301	KK1	C14-N16	2.86	1.40	1.35
2	I	301	KK1	C07-C08	2.86	1.63	1.51
2	B	301	KK1	C24-C23	2.93	1.44	1.38
2	I	301	KK1	C10-N16	2.96	1.39	1.34
2	J	301	KK1	C14-N13	2.99	1.40	1.35
2	A	301	KK1	C24-C23	3.00	1.44	1.38
2	B	301	KK1	C10-N16	3.00	1.39	1.34
2	G	301	KK1	C14-N13	3.01	1.40	1.35
2	H	301	KK1	C14-N13	3.03	1.40	1.35
2	J	301	KK1	C11-C12	3.09	1.44	1.39
2	G	301	KK1	C11-C12	3.10	1.44	1.39
2	H	301	KK1	C11-C12	3.10	1.44	1.39
2	C	301	KK1	C07-C08	3.11	1.64	1.51
2	I	301	KK1	C14-N16	3.16	1.41	1.35
2	A	301	KK1	C08-N09	3.16	1.53	1.46
2	F	301	KK1	C24-C23	3.17	1.44	1.38
2	H	301	KK1	C24-C23	3.33	1.44	1.38
2	J	301	KK1	C24-C23	3.33	1.44	1.38
2	G	301	KK1	C24-C23	3.34	1.44	1.38
2	J	301	KK1	C10-N09	3.35	1.45	1.35
2	H	301	KK1	C10-N09	3.35	1.45	1.35
2	G	301	KK1	C10-N09	3.38	1.45	1.35
2	A	301	KK1	C11-C10	3.41	1.47	1.39
2	F	301	KK1	C11-C10	3.42	1.47	1.39
2	E	301	KK1	C24-C23	3.45	1.45	1.38
2	F	301	KK1	C08-N09	3.54	1.54	1.46
2	G	301	KK1	C11-C10	3.56	1.46	1.40
2	J	301	KK1	C11-C10	3.57	1.46	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	301	KK1	C11-C10	3.59	1.46	1.40
2	I	301	KK1	C24-C23	3.60	1.45	1.38
2	I	301	KK1	C08-N09	3.60	1.54	1.46
2	A	301	KK1	C11-C12	3.65	1.45	1.39
2	E	301	KK1	C11-C10	3.68	1.48	1.39
2	D	301	KK1	C11-C10	3.73	1.48	1.39
2	I	301	KK1	C11-C10	3.75	1.48	1.39
2	C	301	KK1	C11-C10	3.79	1.48	1.39
2	J	301	KK1	C18-C19	3.82	1.45	1.38
2	G	301	KK1	C18-C19	3.85	1.45	1.38
2	H	301	KK1	C18-C19	3.86	1.45	1.38
2	B	301	KK1	C11-C12	3.89	1.45	1.39
2	F	301	KK1	C11-C12	3.93	1.45	1.39
2	B	301	KK1	C11-C10	3.95	1.49	1.39
2	F	301	KK1	C14-N15	4.04	1.42	1.34
2	F	301	KK1	C18-C19	4.11	1.46	1.38
2	I	301	KK1	C18-C19	4.13	1.46	1.38
2	E	301	KK1	C11-C12	4.18	1.46	1.39
2	I	301	KK1	C11-C12	4.19	1.46	1.39
2	B	301	KK1	C14-N15	4.21	1.42	1.34
2	C	301	KK1	C24-C23	4.22	1.46	1.38
2	E	301	KK1	C14-N15	4.33	1.42	1.34
2	D	301	KK1	C11-C12	4.33	1.46	1.39
2	I	301	KK1	C14-N15	4.34	1.42	1.34
2	C	301	KK1	C18-C19	4.47	1.46	1.38
2	E	301	KK1	C18-C19	4.47	1.46	1.38
2	C	301	KK1	C11-C12	4.47	1.46	1.39
2	A	301	KK1	C18-C19	4.47	1.46	1.38
2	A	301	KK1	C14-N15	4.50	1.43	1.34
2	H	301	KK1	C14-N15	4.53	1.43	1.34
2	J	301	KK1	C14-N15	4.53	1.43	1.34
2	G	301	KK1	C14-N15	4.56	1.43	1.34
2	D	301	KK1	C18-C19	4.57	1.47	1.38
2	C	301	KK1	C08-N09	4.66	1.57	1.46
2	C	301	KK1	C14-N15	4.70	1.43	1.34
2	D	301	KK1	C08-N09	4.70	1.57	1.46
2	D	301	KK1	C14-N15	4.80	1.43	1.34
2	B	301	KK1	C18-C19	4.90	1.47	1.38
2	A	301	KK1	C10-N09	6.27	1.45	1.36
2	D	301	KK1	C10-N09	7.22	1.46	1.36
2	F	301	KK1	C10-N09	7.64	1.47	1.36
2	C	301	KK1	C10-N09	7.81	1.47	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	KK1	C17-C12	8.28	1.61	1.48
2	I	301	KK1	C10-N09	8.34	1.48	1.36
2	B	301	KK1	C17-C12	8.44	1.61	1.48
2	B	301	KK1	C10-N09	8.66	1.49	1.36
2	E	301	KK1	C10-N09	9.05	1.49	1.36
2	I	301	KK1	C17-C12	9.28	1.63	1.48
2	H	301	KK1	C17-C12	9.38	1.63	1.48
2	J	301	KK1	C17-C12	9.41	1.63	1.48
2	G	301	KK1	C17-C12	9.44	1.63	1.48
2	C	301	KK1	C17-C12	9.84	1.63	1.48
2	E	301	KK1	C17-C12	9.91	1.63	1.48
2	F	301	KK1	C17-C12	10.09	1.64	1.48
2	D	301	KK1	C17-C12	11.27	1.66	1.48

All (104) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	302	NAG	C2-N2-C7	-4.84	116.83	123.04
3	I	302	NAG	O3-C3-C4	-4.36	100.53	110.34
3	I	302	NAG	C4-C3-C2	-4.30	104.54	111.23
3	C	302	NAG	C3-C2-N2	-4.26	100.36	110.56
3	I	302	NAG	C3-C4-C5	-3.42	104.24	110.20
2	D	301	KK1	C24-C17-C18	-3.28	111.35	117.55
2	C	301	KK1	C24-C17-C18	-3.13	111.63	117.55
2	D	301	KK1	C11-C12-N13	-2.96	118.94	122.39
3	C	302	NAG	C6-C5-C4	-2.94	105.76	113.02
3	C	302	NAG	O4-C4-C3	-2.92	103.77	110.34
2	E	301	KK1	C24-C17-C18	-2.89	112.08	117.55
2	H	301	KK1	C24-C17-C18	-2.77	112.31	117.55
2	J	301	KK1	C24-C17-C18	-2.77	112.31	117.55
2	F	301	KK1	C24-C17-C18	-2.77	112.31	117.55
2	G	301	KK1	C24-C17-C18	-2.77	112.31	117.55
2	A	301	KK1	C24-C17-C18	-2.77	112.32	117.55
2	I	301	KK1	C24-C17-C18	-2.75	112.36	117.55
3	C	302	NAG	O7-C7-C8	-2.70	117.11	122.06
3	A	302	NAG	C2-N2-C7	-2.68	119.59	123.04
2	B	301	KK1	C24-C17-C18	-2.67	112.51	117.55
2	J	301	KK1	C11-C12-N13	-2.57	119.39	122.39
2	G	301	KK1	C11-C12-N13	-2.57	119.39	122.39
2	H	301	KK1	C11-C12-N13	-2.57	119.40	122.39
2	A	301	KK1	C11-C12-N13	-2.54	119.43	122.39
2	B	301	KK1	C17-C12-N13	-2.51	112.79	116.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	301	KK1	C11-C12-N13	-2.50	119.48	122.39
2	I	301	KK1	C11-C12-N13	-2.45	119.53	122.39
2	E	301	KK1	C11-C12-N13	-2.40	119.59	122.39
2	J	301	KK1	C23-C20-C19	-2.39	116.31	120.20
2	C	301	KK1	C11-C12-N13	-2.39	119.60	122.39
3	J	302	NAG	C2-N2-C7	-2.39	119.97	123.04
2	H	301	KK1	C23-C20-C19	-2.38	116.33	120.20
2	G	301	KK1	C23-C20-C19	-2.37	116.36	120.20
3	A	302	NAG	C4-C3-C2	-2.36	107.56	111.23
2	C	301	KK1	C23-C20-C19	-2.25	116.56	120.20
2	B	301	KK1	C11-C12-N13	-2.24	119.78	122.39
2	I	301	KK1	C23-C20-C19	-2.24	116.57	120.20
2	E	301	KK1	C23-C20-C19	-2.20	116.63	120.20
2	D	301	KK1	C23-C20-C19	-2.18	116.66	120.20
3	D	302	NAG	C1-O5-C5	-2.09	109.59	112.25
2	F	301	KK1	C23-C20-C19	-2.07	116.84	120.20
2	A	301	KK1	C23-C20-C19	-2.02	116.92	120.20
2	B	301	KK1	C23-C20-C19	-2.01	116.93	120.20
2	F	301	KK1	C07-C08-N09	2.00	116.88	111.46
3	C	302	NAG	O4-C4-C5	2.04	114.65	109.24
2	A	301	KK1	C24-C17-C12	2.06	124.60	121.24
3	I	302	NAG	O5-C5-C6	2.11	111.91	107.35
2	D	301	KK1	C07-C08-N09	2.13	117.21	111.46
2	I	301	KK1	C24-C17-C12	2.13	124.71	121.24
2	F	301	KK1	C19-C18-C17	2.13	124.25	121.14
2	D	301	KK1	C24-C17-C12	2.14	124.73	121.24
2	E	301	KK1	C23-C24-C17	2.14	124.27	121.14
2	F	301	KK1	C11-C12-C17	2.17	124.77	121.89
2	J	301	KK1	C23-C24-C17	2.17	124.31	121.14
2	H	301	KK1	C23-C24-C17	2.19	124.33	121.14
2	G	301	KK1	C23-C24-C17	2.19	124.34	121.14
2	I	301	KK1	C19-C18-C17	2.25	124.42	121.14
2	I	301	KK1	C07-C08-N09	2.26	117.57	111.46
2	E	301	KK1	C11-C12-C17	2.31	124.96	121.89
2	C	301	KK1	C23-C24-C17	2.37	124.59	121.14
2	F	301	KK1	C24-C17-C12	2.41	125.17	121.24
2	B	301	KK1	C19-C18-C17	2.44	124.70	121.14
2	I	301	KK1	C11-C12-C17	2.48	125.19	121.89
2	F	301	KK1	C23-C24-C17	2.51	124.80	121.14
2	I	301	KK1	C23-C24-C17	2.54	124.84	121.14
3	G	302	NAG	C1-O5-C5	2.55	115.48	112.25
2	B	301	KK1	C23-C24-C17	2.60	124.94	121.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	302	NAG	O5-C5-C6	2.64	113.07	107.35
2	E	301	KK1	C07-C08-N09	2.65	118.65	111.46
2	G	301	KK1	C19-C18-C17	2.67	125.04	121.14
2	E	301	KK1	C24-C17-C12	2.68	125.61	121.24
2	H	301	KK1	C19-C18-C17	2.69	125.06	121.14
2	J	301	KK1	C19-C18-C17	2.70	125.08	121.14
2	D	301	KK1	C23-C24-C17	2.85	125.31	121.14
2	E	301	KK1	C19-C18-C17	2.86	125.32	121.14
2	C	301	KK1	C19-C18-C17	2.89	125.35	121.14
2	A	301	KK1	C23-C24-C17	2.95	125.44	121.14
3	E	302	NAG	C1-O5-C5	3.04	116.10	112.25
2	D	301	KK1	C19-C18-C17	3.06	125.61	121.14
3	C	302	NAG	C2-N2-C7	3.12	127.05	123.04
2	J	301	KK1	C24-C17-C12	3.25	126.54	121.24
2	C	301	KK1	C24-C17-C12	3.25	126.55	121.24
2	H	301	KK1	C24-C17-C12	3.27	126.57	121.24
2	G	301	KK1	C24-C17-C12	3.28	126.59	121.24
2	A	301	KK1	C11-C12-C17	3.28	126.24	121.89
2	C	301	KK1	C11-C12-C17	3.38	126.38	121.89
3	I	302	NAG	C3-C2-N2	3.42	118.75	110.56
2	D	301	KK1	C11-C12-C17	3.44	126.46	121.89
2	B	301	KK1	C11-C12-C17	4.18	127.43	121.89
2	J	301	KK1	C12-N13-C14	4.32	118.68	116.34
2	G	301	KK1	C12-N13-C14	4.33	118.69	116.34
2	H	301	KK1	C12-N13-C14	4.35	118.70	116.34
3	I	302	NAG	O3-C3-C2	4.54	118.10	109.11
3	C	302	NAG	C1-O5-C5	4.62	118.11	112.25
3	C	302	NAG	C8-C7-N2	4.64	125.00	116.11
3	J	302	NAG	C1-O5-C5	5.13	118.75	112.25
2	E	301	KK1	C12-N13-C14	5.34	119.23	116.34
3	I	302	NAG	C1-O5-C5	6.13	120.03	112.25
2	I	301	KK1	C12-N13-C14	6.18	119.69	116.34
2	F	301	KK1	C12-N13-C14	6.29	119.75	116.34
2	C	301	KK1	C12-N13-C14	6.70	119.97	116.34
2	A	301	KK1	C12-N13-C14	7.45	120.38	116.34
2	D	301	KK1	C12-N13-C14	7.59	120.45	116.34
2	B	301	KK1	C12-N13-C14	8.15	120.76	116.34

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

15 monomers are involved in 51 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	KK1	4	0
4	A	303	PO4	1	0
2	B	301	KK1	6	0
2	C	301	KK1	2	0
3	C	302	NAG	6	0
4	C	305	PO4	1	0
2	D	301	KK1	1	0
2	E	301	KK1	1	0
2	F	301	KK1	8	0
3	F	302	NAG	1	0
2	H	301	KK1	6	0
2	I	301	KK1	4	0
3	I	302	NAG	1	0
2	J	301	KK1	8	0
4	J	303	PO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	207/217 (95%)	0.09	9 (4%) 39 38	9, 26, 53, 75	0
1	B	207/217 (95%)	-0.24	0 100 100	8, 20, 39, 55	0
1	C	209/217 (96%)	-0.19	2 (0%) 84 85	7, 21, 52, 67	0
1	D	213/217 (98%)	-0.02	9 (4%) 40 39	9, 21, 53, 82	0
1	E	205/217 (94%)	-0.00	3 (1%) 76 76	9, 26, 50, 66	0
1	F	209/217 (96%)	0.10	4 (1%) 70 70	10, 33, 63, 83	0
1	G	208/217 (95%)	-0.07	4 (1%) 70 70	9, 31, 59, 76	0
1	H	207/217 (95%)	0.11	9 (4%) 39 38	15, 34, 61, 72	0
1	I	207/217 (95%)	0.11	5 (2%) 62 62	13, 30, 56, 81	0
1	J	206/217 (94%)	0.14	8 (3%) 43 43	12, 34, 62, 75	0
All	All	2078/2170 (95%)	0.00	53 (2%) 59 59	7, 27, 57, 83	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	160	ASP	7.1
1	A	156	THR	7.1
1	A	158	ASN	6.9
1	A	159	SER	6.7
1	D	158	ASN	5.5
1	A	157	GLU	5.0
1	D	161	ASP	4.9
1	A	161	ASP	4.8
1	A	-6	TYR	4.6
1	F	189	PRO	4.5
1	H	205	GLY	3.7
1	G	-6	TYR	3.4
1	I	185	TYR	3.4

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Mol	Chain	Res	Type	RSRZ
1	J	160	ASP	3.3
1	J	205	GLY	3.3
1	I	-6	TYR	3.2
1	H	22	GLN	3.2
1	C	72	ASP	3.1
1	F	188	CYS	2.9
1	I	205	GLY	2.9
1	J	44	ILE	2.9
1	E	136	CYS	2.9
1	E	190	GLU	2.8
1	J	25	ARG	2.8
1	J	132	SER	2.8
1	H	-4	ASP	2.8
1	H	108	ASP	2.8
1	J	23	ARG	2.8
1	E	-6	TYR	2.7
1	D	68	SER	2.6
1	G	157	GLU	2.6
1	A	-7	ASP	2.6
1	H	-7	ASP	2.5
1	H	23	ARG	2.5
1	I	23	ARG	2.5
1	J	128	VAL	2.4
1	A	155	THR	2.4
1	H	24	ASP	2.4
1	G	179	LYS	2.4
1	D	160	ASP	2.4
1	D	159	SER	2.4
1	C	71	PRO	2.3
1	H	43	GLU	2.3
1	D	162	SER	2.2
1	J	43	GLU	2.2
1	F	68	SER	2.1
1	D	185	TYR	2.1
1	D	67	SER	2.1
1	G	190	GLU	2.0
1	I	129	ASP	2.0
1	H	89	TYR	2.0
1	D	163	GLU	2.0
1	F	137	ARG	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
4	PO4	B	303	5/5	0.90	0.45	8.83	37,48,63,72	0
4	PO4	G	304	5/5	0.96	0.21	4.35	37,40,51,67	0
3	NAG	I	302	14/15	0.77	0.25	4.32	38,57,62,65	0
2	KK1	B	301	24/24	0.91	0.21	4.06	16,30,45,48	0
3	NAG	J	302	14/15	0.84	0.31	3.51	45,57,68,70	0
4	PO4	F	304	5/5	0.92	0.27	3.21	45,46,48,71	0
4	PO4	E	305	5/5	0.64	0.32	3.03	78,80,87,103	0
2	KK1	E	301	24/24	0.86	0.26	2.50	23,32,55,69	0
4	PO4	C	305	5/5	0.92	0.27	2.45	47,51,63,66	0
2	KK1	C	301	24/24	0.87	0.26	2.40	14,22,44,51	0
2	KK1	F	301	24/24	0.88	0.31	2.19	24,34,59,63	0
2	KK1	G	301	16/24	0.87	0.27	2.15	29,36,51,53	0
3	NAG	D	302	14/15	0.71	0.32	1.64	55,65,75,77	0
2	KK1	I	301	24/24	0.88	0.29	1.52	27,37,62,66	0
2	KK1	H	301	16/24	0.86	0.24	1.32	29,37,53,57	0
2	KK1	J	301	16/24	0.80	0.26	1.30	22,38,51,56	0
4	PO4	E	303	5/5	0.99	0.16	0.96	8,12,14,20	0
2	KK1	A	301	24/24	0.93	0.18	0.91	14,21,51,58	0
4	PO4	H	304	5/5	0.98	0.17	0.83	14,16,20,28	0
4	PO4	C	303	5/5	0.99	0.16	0.63	5,7,9,12	0
2	KK1	D	301	24/24	0.90	0.19	0.57	23,34,46,54	0
4	PO4	H	303	5/5	0.98	0.17	0.54	16,17,20,23	0
4	PO4	J	303	5/5	0.98	0.16	0.50	15,16,20,21	0
4	PO4	A	303	5/5	0.97	0.16	0.46	12,14,17,20	0
4	PO4	A	304	5/5	0.97	0.16	0.17	11,15,21,25	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	PO4	G	303	5/5	0.99	0.15	-0.09	12,16,19,21	0
3	NAG	C	302	14/15	0.82	0.16	-0.30	20,31,43,44	0
4	PO4	F	303	5/5	0.98	0.14	-0.46	17,17,20,21	0
3	NAG	E	302	14/15	0.88	0.12	-0.98	55,61,66,66	0
4	PO4	C	304	5/5	0.98	0.13	-1.35	7,12,14,15	0
4	PO4	A	305	5/5	0.85	0.22	-	41,56,64,70	0
3	NAG	B	302	14/15	0.89	0.29	-	59,66,74,76	0
4	PO4	D	304	5/5	0.57	0.39	-	63,76,80,107	0
4	PO4	E	304	5/5	0.62	0.29	-	72,76,79,97	0
3	NAG	H	302	14/15	0.76	0.32	-	61,67,74,81	0
3	NAG	F	302	14/15	0.81	0.19	-	47,58,63,64	0
3	NAG	G	302	14/15	0.77	0.16	-	45,52,63,63	0
3	NAG	A	302	14/15	0.89	0.18	-	48,55,61,68	0
4	PO4	D	303	5/5	0.85	0.25	-	49,56,73,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.