Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 10:56 PM GMT

PDB ID	:	4ZRU
Title	:	X-ray crystal structure of Lymnaea stagnalis acetylcholine binding protein (Ls-AChBP) in complex with 3-[2-[(2S)-pyrrolidin-2-yl]ethynyl]pyridine (TI-5180)
Authors	:	Bobango, J.; Sankaran, B.; Park, J.F.; Wu, J.; Talley, T.T.
Deposited on	:	2015-05-12
Resolution	:	1.90 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

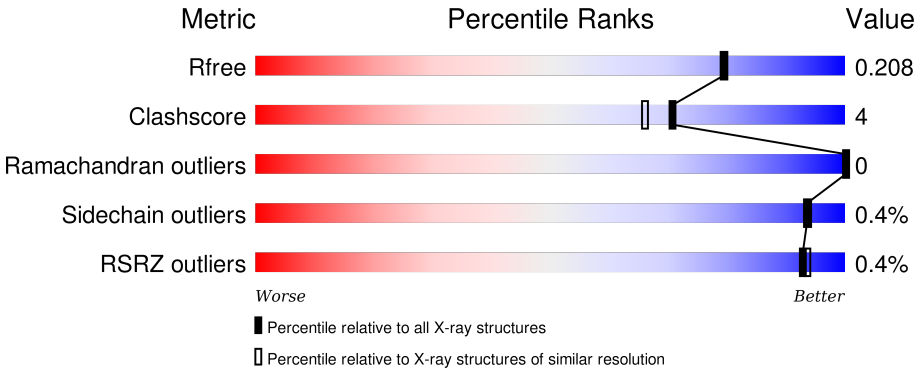
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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 <=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	218	87% 10% ·
1	B	218	89% 6% ·
1	C	218	87% 6% 7%
1	D	218	84% 11% 6%
1	E	218	% 87% 6% 6%

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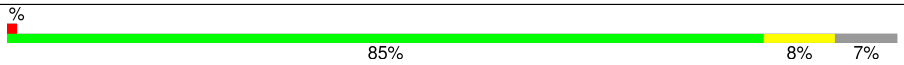
Mol	Chain	Length	Quality of chain
1	F	218	
1	G	218	
1	H	218	
1	I	218	
1	J	218	

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
3	TI9	B	302	-	-	-	X
3	TI9	G	302	-	-	-	X
3	TI9	H	301	-	-	-	X
3	TI9	I	303	-	-	-	X
3	TI9	J	301	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17808 atoms, of which 120 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	211	Total	C	N	O	S	0	0	0
			1631	1023	274	329	5			
1	B	209	Total	C	N	O	S	0	3	0
			1648	1032	277	334	5			
1	C	203	Total	C	N	O	S	0	4	0
			1599	1005	267	322	5			
1	D	206	Total	C	N	O	S	0	2	0
			1617	1016	271	325	5			
1	E	204	Total	C	N	O	S	0	3	0
			1588	1002	264	317	5			
1	F	211	Total	C	N	O	S	0	1	0
			1618	1013	267	333	5			
1	G	201	Total	C	N	O	S	0	0	0
			1561	983	261	312	5			
1	H	203	Total	C	N	O	S	0	0	0
			1564	984	259	316	5			
1	I	203	Total	C	N	O	S	0	0	0
			1569	988	263	313	5			
1	J	203	Total	C	N	O	S	0	1	0
			1563	985	261	312	5			

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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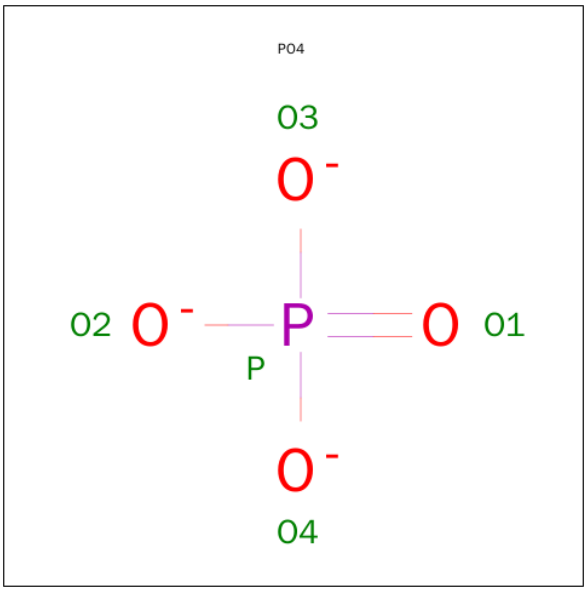
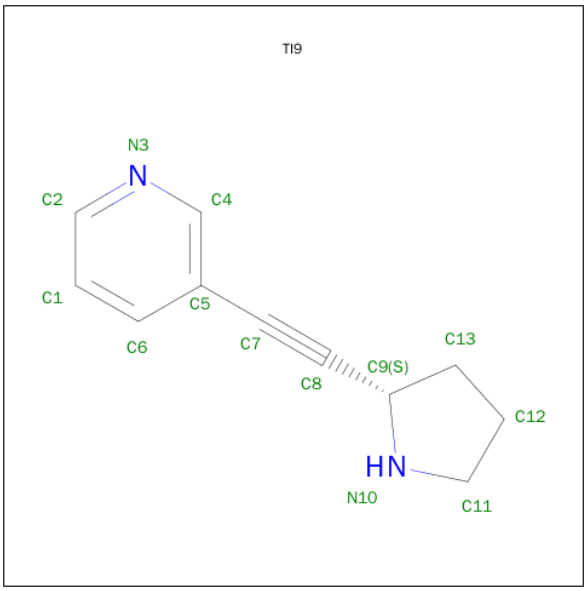
Chain	Residue	Modelled	Actual	Comment	Reference
G	-4	ASP	-	expression tag	UNP P58154
G	-3	ASP	-	expression tag	UNP P58154
G	-2	ASP	-	expression tag	UNP P58154
G	-1	ASP	-	expression tag	UNP P58154
G	0	LYS	-	expression tag	UNP P58154
H	-7	ASP	-	expression tag	UNP P58154
H	-6	TYR	-	expression tag	UNP P58154
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 PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



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

- Molecule 3 is 3-[(2S)-pyrrolidin-2-ylethynyl]pyridine (three-letter code: TI9) (formula: $C_{11}H_{12}N_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C H N 25 11 12 2	0	0
3	B	1	Total C H N 25 11 12 2	0	0
3	C	1	Total C H N 25 11 12 2	0	0
3	D	1	Total C H N 25 11 12 2	0	0
3	E	1	Total C H N 25 11 12 2	0	0
3	F	1	Total C H N 25 11 12 2	0	0
3	G	1	Total C H N 25 11 12 2	0	0
3	H	1	Total C H N 25 11 12 2	0	0
3	I	1	Total C H N 25 11 12 2	0	0
3	J	1	Total C H N 25 11 12 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	144	Total O 144 144	0	0
4	B	185	Total O 185 185	0	0

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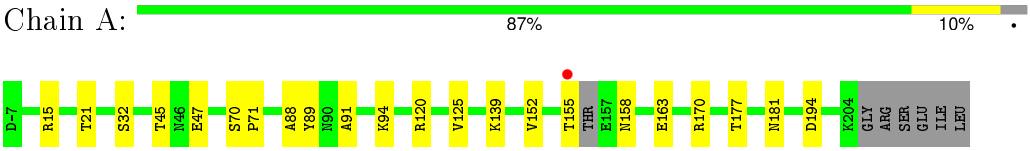
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	168	Total O 168 168	0	0
4	D	161	Total O 161 161	0	0
4	E	181	Total O 181 181	0	0
4	F	158	Total O 158 158	0	0
4	G	120	Total O 120 120	0	0
4	H	133	Total O 133 133	0	0
4	I	159	Total O 159 159	0	0
4	J	141	Total O 141 141	0	0

3 Residue-property plots (i)

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

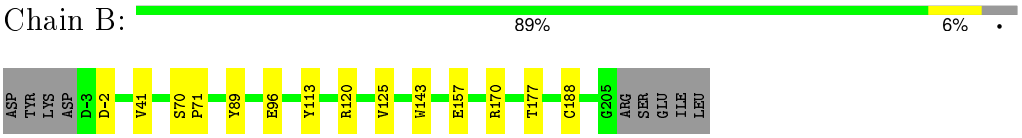
- Molecule 1: Acetylcholine-binding protein

Chain A:



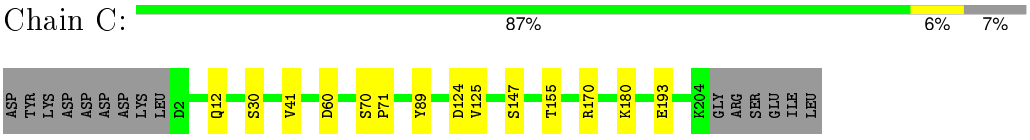
- Molecule 1: Acetylcholine-binding protein

Chain B:



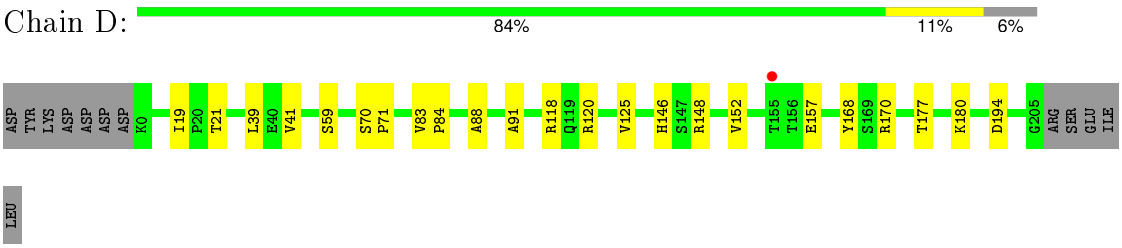
- Molecule 1: Acetylcholine-binding protein

Chain C:



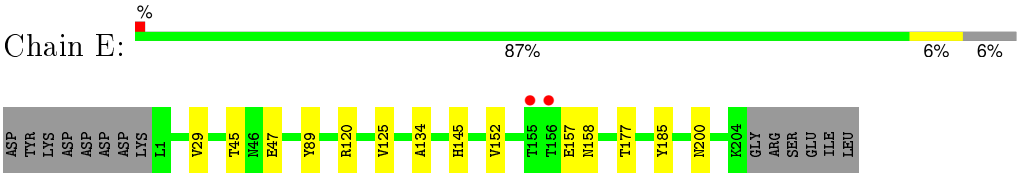
- Molecule 1: Acetylcholine-binding protein

Chain D:



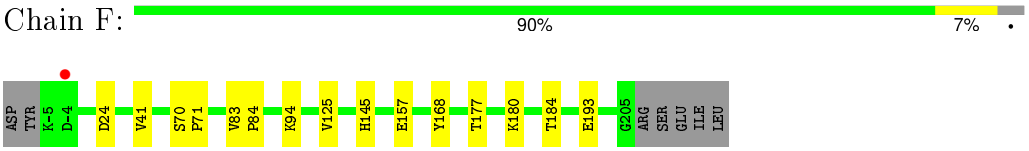
- Molecule 1: Acetylcholine-binding protein

Chain E:



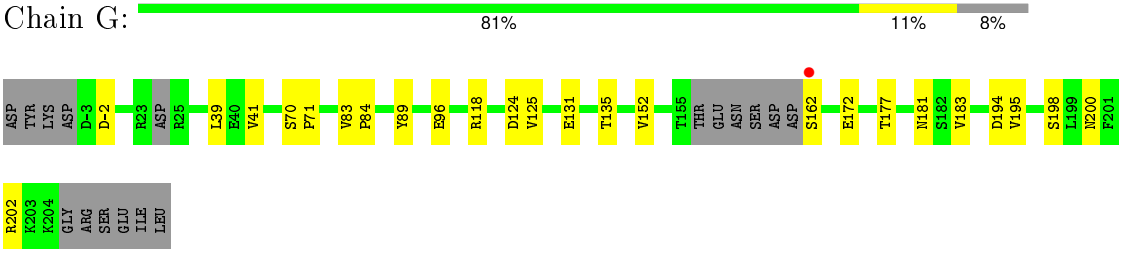
- Molecule 1: Acetylcholine-binding protein

Chain F:



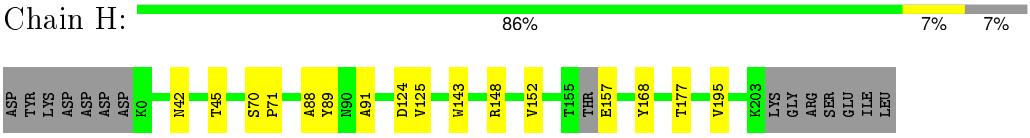
• Molecule 1: Acetylcholine-binding protein

Chain G:



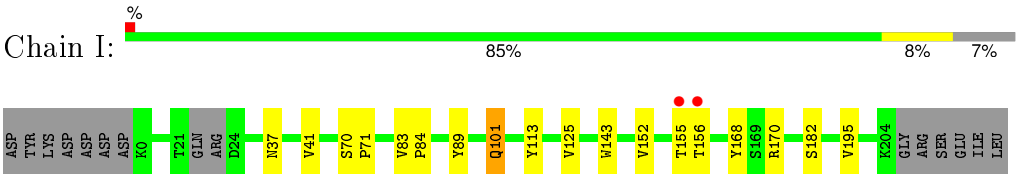
• Molecule 1: Acetylcholine-binding protein

Chain H:



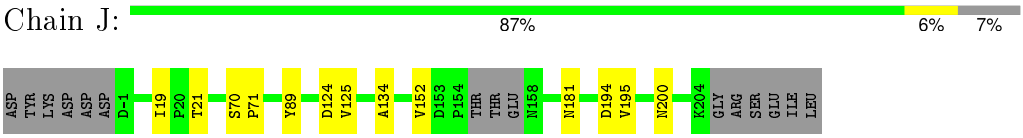
• Molecule 1: Acetylcholine-binding protein

Chain I:



• Molecule 1: Acetylcholine-binding protein

Chain J:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	92.91Å 108.87Å 122.70Å 90.00° 95.31° 90.00°	Depositor
Resolution (Å)	90.00 – 1.90 92.51 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.1 (90.00-1.90) 94.6 (92.51-1.85)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.28 (at 1.86Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.171 , 0.204 0.176 , 0.208	Depositor DCC
R_{free} test set	9553 reflections (5.47%)	DCC
Wilson B-factor (Å^2)	22.5	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 48.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 201803 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	17808	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, TI9

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	0.39	0/1667	0.52	0/2284
1	B	0.38	0/1693	0.54	0/2315
1	C	0.38	0/1647	0.54	0/2256
1	D	0.40	0/1659	0.55	0/2271
1	E	0.39	0/1633	0.52	0/2240
1	F	0.36	0/1657	0.52	0/2273
1	G	0.35	0/1595	0.51	0/2182
1	H	0.34	0/1599	0.50	0/2193
1	I	0.36	0/1604	0.60	3/2198 (0.1%)
1	J	0.36	0/1601	0.53	0/2197
All	All	0.37	0/16355	0.53	3/22409 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	I	155	THR	N-CA-C	9.42	136.45	111.00
1	I	156	THR	N-CA-C	-8.24	88.76	111.00
1	I	156	THR	N-CA-CB	-5.12	100.58	110.30

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	1631	0	1518	18	0
1	B	1648	0	1574	13	0
1	C	1599	0	1526	13	0
1	D	1617	0	1542	18	0
1	E	1588	0	1514	11	0
1	F	1618	0	1496	15	0
1	G	1561	0	1474	21	1
1	H	1564	0	1462	15	1
1	I	1569	0	1483	16	0
1	J	1563	0	1465	11	0
2	A	10	0	0	1	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	F	10	0	0	1	0
2	G	5	0	0	0	0
2	I	10	0	0	0	0
3	A	13	12	12	1	0
3	B	13	12	12	2	0
3	C	13	12	12	1	0
3	D	13	12	12	0	0
3	E	13	12	12	1	0
3	F	13	12	12	0	0
3	G	13	12	12	2	0
3	H	13	12	12	3	0
3	I	13	12	12	3	0
3	J	13	12	12	1	0
4	A	144	0	0	3	1
4	B	185	0	0	2	1
4	C	168	0	0	4	0
4	D	161	0	0	3	0
4	E	181	0	0	0	0
4	F	158	0	0	0	0
4	G	120	0	0	2	0
4	H	133	0	0	0	0
4	I	159	0	0	0	0
4	J	141	0	0	0	0
All	All	17688	120	15174	138	2

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 (138) 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:J:152:VAL:HG12	1:J:195:VAL:HG23	1.43	1.00
1:A:94:LYS:HE3	1:B:96:GLU:HG3	1.62	0.81
1:I:152:VAL:CG1	1:I:195:VAL:HG23	2.11	0.80
1:I:101:GLN:NE2	1:I:113:TYR:OH	2.17	0.78
1:J:152:VAL:CG1	1:J:195:VAL:HG23	2.15	0.76
1:A:47:GLU:OE1	1:A:120:ARG:NH2	2.18	0.76
1:H:152:VAL:CG1	1:H:195:VAL:HG23	2.20	0.71
1:F:94:LYS:HE3	1:G:96:GLU:HG3	1.71	0.70
1:I:152:VAL:HG11	1:I:195:VAL:HG23	1.75	0.67
1:G:131:GLU:OE1	1:G:202:ARG:NH1	2.27	0.67
1:I:152:VAL:HG12	1:I:195:VAL:HG23	1.76	0.67
1:F:70:SER:HB2	1:F:71:PRO:HD2	1.79	0.64
1:H:70:SER:HB2	1:H:71:PRO:HD2	1.81	0.63
1:E:47:GLU:OE1	1:E:120:ARG:NH2	2.32	0.62
1:I:101:GLN:HA	1:I:101:GLN:HE21	1.63	0.62
1:D:170:ARG:NH2	4:D:403:HOH:O	2.30	0.62
1:F:83:VAL:HG13	1:F:84:PRO:HD2	1.82	0.61
1:G:83:VAL:HG13	1:G:84:PRO:HD2	1.84	0.59
1:D:70:SER:HB2	1:D:71:PRO:HD2	1.85	0.59
1:H:89:TYR:CE2	3:H:301:TI9:H112	2.38	0.59
1:G:89:TYR:CE2	3:G:302:TI9:H112	2.38	0.59
1:C:124:ASP:HB2	1:D:168:TYR:CE1	2.38	0.58
1:F:180:LYS:NZ	1:F:193:GLU:OE1	2.36	0.58
1:H:89:TYR:CZ	3:H:301:TI9:H112	2.39	0.58
1:G:172:GLU:OE2	1:G:202:ARG:NE	2.31	0.57
1:H:124:ASP:HB2	1:I:168:TYR:CE1	2.41	0.56
1:C:70:SER:HB2	1:C:71:PRO:HD2	1.88	0.55
1:D:41:VAL:HG13	1:D:125:VAL:HG11	1.89	0.55
1:C:155:THR:HG21	4:C:538:HOH:O	2.07	0.55
1:A:158:ASN:OD1	1:A:177:THR:HG22	2.08	0.54
1:C:60:ASP:OD2	4:C:401:HOH:O	2.19	0.53
1:G:70:SER:HB2	1:G:71:PRO:HD2	1.90	0.53
1:H:152:VAL:HG12	1:H:195:VAL:HG23	1.91	0.53
1:D:83:VAL:HG13	1:D:84:PRO:HD2	1.89	0.53
1:F:41:VAL:CG1	1:F:125:VAL:HG11	2.39	0.53
1:E:158:ASN:OD1	1:E:177[A]:THR:HG22	2.09	0.52
1:A:45:THR:HG22	1:B:170:ARG:NH1	2.24	0.52
1:F:70:SER:HB2	1:F:71:PRO:CD	2.39	0.52
1:B:125:VAL:O	1:B:125:VAL:HG12	2.09	0.52
1:F:24:ASP:HA	1:G:-2:ASP:HB2	1.92	0.51
1:C:41:VAL:CG1	1:C:125:VAL:HG11	2.40	0.51
1:B:143:TRP:O	3:B:302:TI9:H111	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:70:SER:HB2	1:I:71:PRO:HD2	1.92	0.51
1:C:170:ARG:NH2	4:C:407:HOH:O	2.42	0.51
1:J:70:SER:HB2	1:J:71:PRO:CD	2.40	0.51
1:J:125:VAL:O	1:J:125:VAL:HG12	2.11	0.51
1:A:32:SER:HB2	1:A:155:THR:HB	1.93	0.51
1:I:125:VAL:HG12	1:I:125:VAL:O	2.11	0.51
1:H:70:SER:HB2	1:H:71:PRO:CD	2.40	0.51
1:C:125:VAL:HG12	1:C:125:VAL:O	2.09	0.51
1:H:125:VAL:HG12	1:H:125:VAL:O	2.12	0.50
1:I:152:VAL:HG12	1:I:195:VAL:CG2	2.41	0.50
1:B:120:ARG:NH1	4:B:408:HOH:O	2.43	0.50
1:A:70:SER:HB2	1:A:71:PRO:CD	2.41	0.50
1:G:200:ASN:ND2	4:G:403:HOH:O	2.39	0.50
1:G:125:VAL:O	1:G:125:VAL:HG12	2.12	0.50
1:B:70:SER:HB2	1:B:71:PRO:HD2	1.93	0.49
1:E:157:GLU:O	1:E:177[A]:THR:HA	2.12	0.49
1:E:157:GLU:O	1:E:177[B]:THR:HA	2.13	0.49
1:F:157:GLU:O	1:F:177:THR:HA	2.13	0.49
1:D:120:ARG:HD2	4:D:435:HOH:O	2.13	0.49
1:D:157:GLU:O	1:D:177:THR:HA	2.14	0.48
1:J:70:SER:HB2	1:J:71:PRO:HD2	1.96	0.48
1:F:125:VAL:HG12	1:F:125:VAL:O	2.12	0.48
1:I:89:TYR:CE2	3:I:303:TI9:H112	2.48	0.48
1:J:152:VAL:HG12	1:J:195:VAL:CG2	2.29	0.48
1:I:41:VAL:HG12	1:I:125:VAL:HG11	1.96	0.48
2:A:302:PO4:O4	1:E:145:HIS:HE1	1.96	0.47
1:C:70:SER:HB2	1:C:71:PRO:CD	2.44	0.47
1:D:70:SER:HB2	1:D:71:PRO:CD	2.45	0.47
1:G:152:VAL:HG12	1:G:195:VAL:HG23	1.96	0.47
1:G:162:SER:HA	4:G:502:HOH:O	2.16	0.46
1:G:89:TYR:CZ	3:G:302:TI9:H112	2.50	0.46
1:I:143:TRP:CE3	3:I:303:TI9:H111	2.51	0.46
1:C:89:TYR:CE2	3:C:302:TI9:H112	2.50	0.46
1:J:89:TYR:CE2	3:J:301:TI9:H112	2.51	0.46
1:B:113:TYR:OH	4:B:401:HOH:O	2.20	0.45
1:E:134:ALA:O	1:E:200:ASN:HA	2.16	0.45
1:A:70:SER:HB2	1:A:71:PRO:HD2	1.97	0.45
1:H:45:THR:HG22	1:I:170:ARG:NH1	2.31	0.45
1:I:83:VAL:HG13	1:I:84:PRO:HD2	1.98	0.45
1:D:146:HIS:CE1	1:D:148:ARG:HB2	2.51	0.45
1:G:124:ASP:HB2	1:H:168:TYR:CE1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:ARG:NH1	1:E:45:THR:HG22	2.32	0.45
1:C:41:VAL:CG1	1:C:125:VAL:CG1	2.95	0.44
1:D:152:VAL:HG21	1:D:194:ASP:HA	1.99	0.44
1:D:39:LEU:CD1	1:D:118:ARG:CZ	2.96	0.44
1:D:41:VAL:CG1	1:D:125:VAL:HG11	2.47	0.44
1:I:70:SER:HB2	1:I:71:PRO:CD	2.47	0.44
1:H:143:TRP:CE3	3:H:301:TI9:H111	2.53	0.44
1:J:181:ASN:ND2	1:J:194:ASP:OD2	2.49	0.44
1:F:94:LYS:HE3	1:G:96:GLU:CG	2.46	0.43
1:G:181:ASN:HB3	1:G:194:ASP:OD1	2.17	0.43
1:D:180:LYS:HB3	1:D:180:LYS:HE3	1.86	0.43
1:B:89:TYR:OH	3:B:302:TI9:H112	2.17	0.43
1:A:163:GLU:OE2	1:E:185:TYR:HA	2.19	0.43
1:G:177:THR:OG1	1:G:198:SER:HB2	2.18	0.43
1:G:83:VAL:CG1	1:G:84:PRO:HD2	2.49	0.43
1:E:125:VAL:HG12	1:E:125:VAL:O	2.18	0.43
1:F:83:VAL:CG1	1:F:84:PRO:HD2	2.48	0.43
1:F:41:VAL:CG1	1:F:125:VAL:CG1	2.97	0.43
1:A:125:VAL:O	1:A:125:VAL:HG12	2.18	0.43
1:F:41:VAL:HG13	1:F:125:VAL:HG11	2.01	0.42
1:B:41:VAL:CG1	1:B:125:VAL:HG11	2.49	0.42
1:B:41:VAL:HG13	1:B:125:VAL:HG11	2.01	0.42
1:C:12:GLN:HG2	4:C:523:HOH:O	2.18	0.42
1:F:168:TYR:CE1	1:J:124:ASP:HB2	2.54	0.42
1:B:157:GLU:O	1:B:177:THR:HA	2.20	0.42
1:D:41:VAL:HG13	1:D:125:VAL:CG1	2.49	0.42
1:A:88:ALA:HB3	1:A:91:ALA:HB2	2.02	0.42
1:D:19:ILE:HD12	1:D:21:THR:HG23	2.02	0.42
1:H:88:ALA:HB3	1:H:91:ALA:HB2	2.01	0.42
1:D:88:ALA:HB3	1:D:91:ALA:HB2	2.01	0.42
1:H:42:ASN:HB3	1:H:45:THR:OG1	2.20	0.42
1:F:145:HIS:HE1	2:F:302:PO4:O3	2.03	0.42
1:J:134:ALA:O	1:J:200:ASN:HA	2.20	0.41
1:G:183:VAL:HG13	1:G:194:ASP:OD2	2.20	0.41
1:A:139:LYS:NZ	1:A:181:ASN:OD1	2.45	0.41
1:E:29:VAL:O	1:E:152:VAL:HA	2.20	0.41
1:D:125:VAL:HG12	1:D:125:VAL:O	2.20	0.41
1:B:70:SER:HB2	1:B:71:PRO:CD	2.49	0.41
1:A:15:ARG:HD3	4:A:1919:HOH:O	2.19	0.41
1:H:152:VAL:HG11	1:H:195:VAL:HG23	1.99	0.41
1:A:170:ARG:NH2	4:A:1806:HOH:O	2.43	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:147[B]:SER:OG	1:C:193:GLU:HG3	2.20	0.41
1:A:170:ARG:NH1	4:A:1814:HOH:O	2.54	0.41
1:J:19:ILE:HD12	1:J:21:THR:HG23	2.03	0.41
1:A:152:VAL:HG21	1:A:194:ASP:HA	2.02	0.41
1:D:170:ARG:NH1	4:D:411:HOH:O	2.53	0.41
1:A:89:TYR:CE2	3:A:303:TI9:H112	2.55	0.41
1:G:39:LEU:CD1	1:G:118:ARG:CZ	2.99	0.41
1:E:89:TYR:CE2	3:E:301:TI9:H112	2.56	0.41
1:I:89:TYR:CZ	3:I:303:TI9:H112	2.56	0.40
1:C:30:SER:HB3	1:C:155:THR:HG22	2.04	0.40
1:H:157:GLU:O	1:H:177:THR:HA	2.21	0.40
1:G:41:VAL:CG1	1:G:125:VAL:HG11	2.51	0.40
1:G:39:LEU:HD12	1:G:118:ARG:NE	2.37	0.40
1:A:21:THR:OG1	1:B:-2:ASP:HB3	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1872:HOH:O	4:B:446:HOH:O[2_646]	2.12	0.08
1:G:135:THR:OG1	1:H:148:ARG:NH2[2_657]	2.14	0.06

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	207/218 (95%)	207 (100%)	0	0	100	100
1	B	210/218 (96%)	210 (100%)	0	0	100	100
1	C	205/218 (94%)	204 (100%)	1 (0%)	0	100	100
1	D	206/218 (94%)	206 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	205/218 (94%)	203 (99%)	2 (1%)	0	100	100
1	F	210/218 (96%)	209 (100%)	1 (0%)	0	100	100
1	G	195/218 (89%)	194 (100%)	1 (0%)	0	100	100
1	H	199/218 (91%)	199 (100%)	0	0	100	100
1	I	199/218 (91%)	197 (99%)	2 (1%)	0	100	100
1	J	200/218 (92%)	200 (100%)	0	0	100	100
All	All	2036/2180 (93%)	2029 (100%)	7 (0%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	183/204 (90%)	183 (100%)	0	100	100
1	B	191/204 (94%)	190 (100%)	1 (0%)	92	92
1	C	186/204 (91%)	185 (100%)	1 (0%)	92	92
1	D	186/204 (91%)	185 (100%)	1 (0%)	92	92
1	E	182/204 (89%)	182 (100%)	0	100	100
1	F	182/204 (89%)	181 (100%)	1 (0%)	92	92
1	G	177/204 (87%)	177 (100%)	0	100	100
1	H	177/204 (87%)	177 (100%)	0	100	100
1	I	178/204 (87%)	175 (98%)	3 (2%)	68	64
1	J	176/204 (86%)	176 (100%)	0	100	100
All	All	1818/2040 (89%)	1811 (100%)	7 (0%)	93	94

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	188	CYS

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Mol	Chain	Res	Type
1	C	180	LYS
1	D	59	SER
1	F	184	THR
1	I	37	ASN
1	I	101	GLN
1	I	182	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	158	ASN
1	I	101	GLN

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

20 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	PO4	A	301	-	4,4,4	0.37	0	6,6,6	0.26	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	PO4	A	302	-	4,4,4	0.60	0	6,6,6	0.29	0
3	TI9	A	303	-	12,14,14	1.72	3 (25%)	15,17,17	1.60	4 (26%)
2	PO4	B	301	-	4,4,4	0.39	0	6,6,6	0.26	0
3	TI9	B	302	-	12,14,14	1.39	2 (16%)	15,17,17	1.54	4 (26%)
2	PO4	C	301	-	4,4,4	0.45	0	6,6,6	0.26	0
3	TI9	C	302	-	12,14,14	1.48	2 (16%)	15,17,17	1.77	5 (33%)
2	PO4	D	301	-	4,4,4	0.38	0	6,6,6	0.27	0
3	TI9	D	302	-	12,14,14	1.69	2 (16%)	15,17,17	1.79	4 (26%)
3	TI9	E	301	-	12,14,14	1.59	2 (16%)	15,17,17	1.56	4 (26%)
2	PO4	F	301	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	F	302	-	4,4,4	0.31	0	6,6,6	0.28	0
3	TI9	F	303	-	12,14,14	1.51	2 (16%)	15,17,17	1.58	4 (26%)
2	PO4	G	301	-	4,4,4	0.44	0	6,6,6	0.27	0
3	TI9	G	302	-	12,14,14	1.48	1 (8%)	15,17,17	2.00	6 (40%)
3	TI9	H	301	-	12,14,14	1.38	2 (16%)	15,17,17	1.82	4 (26%)
2	PO4	I	301	-	4,4,4	0.32	0	6,6,6	0.29	0
2	PO4	I	302	-	4,4,4	0.37	0	6,6,6	0.30	0
3	TI9	I	303	-	12,14,14	1.51	2 (16%)	15,17,17	1.61	4 (26%)
3	TI9	J	301	-	12,14,14	1.54	2 (16%)	15,17,17	1.46	3 (20%)

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	PO4	A	301	-	-	0/0/0/0	0/0/0/0
2	PO4	A	302	-	-	0/0/0/0	0/0/0/0
3	TI9	A	303	-	-	0/3/12/12	0/2/2/2
2	PO4	B	301	-	-	0/0/0/0	0/0/0/0
3	TI9	B	302	-	-	0/3/12/12	0/2/2/2
2	PO4	C	301	-	-	0/0/0/0	0/0/0/0
3	TI9	C	302	-	-	0/3/12/12	0/2/2/2
2	PO4	D	301	-	-	0/0/0/0	0/0/0/0
3	TI9	D	302	-	-	0/3/12/12	0/2/2/2
3	TI9	E	301	-	-	0/3/12/12	0/2/2/2
2	PO4	F	301	-	-	0/0/0/0	0/0/0/0
2	PO4	F	302	-	-	0/0/0/0	0/0/0/0
3	TI9	F	303	-	-	0/3/12/12	0/2/2/2
2	PO4	G	301	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TI9	G	302	-	-	0/3/12/12	0/2/2/2
3	TI9	H	301	-	-	0/3/12/12	0/2/2/2
2	PO4	I	301	-	-	0/0/0/0	0/0/0/0
2	PO4	I	302	-	-	0/0/0/0	0/0/0/0
3	TI9	I	303	-	-	0/3/12/12	0/2/2/2
3	TI9	J	301	-	-	0/3/12/12	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	303	TI9	C13-C9	-4.84	1.47	1.53
3	D	302	TI9	C13-C9	-4.80	1.47	1.53
3	J	301	TI9	C13-C9	-4.35	1.48	1.53
3	E	301	TI9	C13-C9	-4.26	1.48	1.53
3	C	302	TI9	C13-C9	-4.12	1.48	1.53
3	F	303	TI9	C13-C9	-4.02	1.48	1.53
3	I	303	TI9	C13-C9	-3.92	1.48	1.53
3	G	302	TI9	C13-C9	-3.61	1.48	1.53
3	B	302	TI9	C13-C9	-3.40	1.49	1.53
3	H	301	TI9	C13-C9	-3.33	1.49	1.53
3	A	303	TI9	C4-C5	-2.04	1.36	1.39
3	H	301	TI9	C4-C5	-2.03	1.36	1.39
3	C	302	TI9	C5-C7	2.11	1.49	1.44
3	D	302	TI9	C5-C7	2.12	1.49	1.44
3	A	303	TI9	C5-C7	2.24	1.49	1.44
3	B	302	TI9	C5-C7	2.37	1.50	1.44
3	I	303	TI9	C5-C7	2.37	1.50	1.44
3	J	301	TI9	C5-C7	2.38	1.50	1.44
3	F	303	TI9	C5-C7	2.40	1.50	1.44
3	E	301	TI9	C5-C7	2.47	1.50	1.44

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	G	302	TI9	C5-C4-N3	-3.69	119.09	123.63
3	D	302	TI9	C5-C4-N3	-3.52	119.29	123.63
3	C	302	TI9	C5-C4-N3	-3.47	119.36	123.63
3	H	301	TI9	C5-C4-N3	-3.23	119.65	123.63
3	B	302	TI9	C5-C4-N3	-3.06	119.86	123.63
3	J	301	TI9	C5-C4-N3	-3.01	119.92	123.63
3	I	303	TI9	C5-C4-N3	-2.99	119.95	123.63
3	A	303	TI9	C5-C4-N3	-2.99	119.95	123.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	302	TI9	C4-C5-C7	-2.94	115.26	120.77
3	F	303	TI9	C5-C4-N3	-2.78	120.20	123.63
3	E	301	TI9	C5-C4-N3	-2.75	120.24	123.63
3	C	302	TI9	C5-C7-C8	-2.74	173.25	177.38
3	G	302	TI9	C5-C7-C8	-2.66	173.37	177.38
3	E	301	TI9	C4-C5-C7	-2.56	115.97	120.77
3	H	301	TI9	C4-C5-C7	-2.48	116.11	120.77
3	C	302	TI9	C4-C5-C7	-2.38	116.30	120.77
3	I	303	TI9	C4-C5-C7	-2.34	116.38	120.77
3	F	303	TI9	C4-C5-C7	-2.33	116.39	120.77
3	A	303	TI9	C4-C5-C7	-2.21	116.61	120.77
3	G	302	TI9	C4-C5-C7	-2.18	116.67	120.77
3	B	302	TI9	C11-N10-C9	-2.16	101.95	107.20
3	E	301	TI9	C2-N3-C4	2.02	120.57	116.84
3	G	302	TI9	C13-C9-N10	2.12	109.39	105.01
3	J	301	TI9	C2-N3-C4	2.27	121.01	116.84
3	F	303	TI9	C2-N3-C4	2.35	121.18	116.84
3	A	303	TI9	C2-N3-C4	2.41	121.28	116.84
3	B	302	TI9	C6-C5-C4	2.42	120.33	117.10
3	C	302	TI9	C2-N3-C4	2.46	121.38	116.84
3	B	302	TI9	C2-N3-C4	2.59	121.60	116.84
3	H	301	TI9	C2-N3-C4	2.63	121.69	116.84
3	I	303	TI9	C2-N3-C4	2.64	121.71	116.84
3	D	302	TI9	C2-N3-C4	2.65	121.72	116.84
3	G	302	TI9	C2-N3-C4	2.67	121.75	116.84
3	I	303	TI9	C6-C5-C4	2.71	120.70	117.10
3	F	303	TI9	C6-C5-C4	2.81	120.84	117.10
3	J	301	TI9	C6-C5-C4	2.84	120.89	117.10
3	E	301	TI9	C6-C5-C4	3.01	121.10	117.10
3	A	303	TI9	C6-C5-C4	3.08	121.20	117.10
3	H	301	TI9	C6-C5-C4	3.22	121.38	117.10
3	C	302	TI9	C6-C5-C4	3.22	121.38	117.10
3	D	302	TI9	C6-C5-C4	3.33	121.53	117.10
3	G	302	TI9	C6-C5-C4	3.47	121.72	117.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	302	PO4	1	0
3	A	303	TI9	1	0
3	B	302	TI9	2	0
3	C	302	TI9	1	0
3	E	301	TI9	1	0
2	F	302	PO4	1	0
3	G	302	TI9	2	0
3	H	301	TI9	3	0
3	I	303	TI9	3	0
3	J	301	TI9	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^{th} 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	211/218 (96%)	-0.18	1 (0%) 91 92	14, 24, 40, 52	0
1	B	209/218 (95%)	-0.26	0 100 100	14, 22, 36, 48	0
1	C	203/218 (93%)	-0.24	0 100 100	14, 22, 34, 40	0
1	D	206/218 (94%)	-0.23	1 (0%) 91 92	14, 22, 38, 50	0
1	E	204/218 (93%)	-0.22	2 (0%) 84 86	13, 21, 38, 45	0
1	F	211/218 (96%)	-0.24	1 (0%) 91 92	16, 25, 41, 57	0
1	G	201/218 (92%)	-0.17	1 (0%) 91 92	17, 27, 41, 52	0
1	H	203/218 (93%)	-0.23	0 100 100	16, 27, 44, 51	0
1	I	203/218 (93%)	-0.22	2 (0%) 84 86	16, 27, 39, 53	0
1	J	203/218 (93%)	-0.22	0 100 100	17, 26, 40, 51	0
All	All	2054/2180 (94%)	-0.22	8 (0%) 93 93	13, 24, 40, 57	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	155	THR	4.9
1	I	156	THR	4.0
1	G	162	SER	3.0
1	A	155	THR	2.7
1	E	155	THR	2.7
1	F	-4	ASP	2.5
1	E	156	THR	2.3
1	D	155	THR	2.1

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

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	TI9	H	301	13/13	0.95	0.13	7.58	16,25,30,32	0
3	TI9	G	302	13/13	0.94	0.16	4.26	22,30,37,38	0
3	TI9	B	302	13/13	0.94	0.12	4.01	17,24,33,40	0
3	TI9	J	301	13/13	0.97	0.12	3.15	20,25,33,35	0
3	TI9	I	303	13/13	0.97	0.11	2.22	21,26,34,37	0
3	TI9	F	303	13/13	0.95	0.12	1.31	18,27,33,35	0
3	TI9	A	303	13/13	0.97	0.10	0.48	18,23,28,29	0
2	PO4	B	301	5/5	0.98	0.09	-0.04	23,26,28,28	0
2	PO4	I	302	5/5	0.97	0.09	-0.08	22,25,27,27	0
3	TI9	D	302	13/13	0.97	0.09	-0.39	15,21,27,28	0
2	PO4	D	301	5/5	0.96	0.09	-0.46	20,22,26,27	0
3	TI9	C	302	13/13	0.97	0.08	-0.77	16,23,30,32	0
2	PO4	I	301	5/5	0.98	0.07	-0.82	19,23,23,24	0
2	PO4	G	301	5/5	0.98	0.08	-0.99	18,21,23,24	0
2	PO4	C	301	5/5	0.99	0.08	-1.26	20,22,22,23	0
2	PO4	A	302	5/5	0.99	0.08	-1.30	15,15,16,17	0
2	PO4	A	301	5/5	0.98	0.08	-1.32	18,20,22,22	0
2	PO4	F	302	5/5	0.99	0.07	-1.55	22,22,24,24	0
3	TI9	E	301	13/13	0.97	0.08	-1.91	17,20,24,28	0
2	PO4	F	301	5/5	0.99	0.06	-3.20	20,21,22,23	0

6.5 Other polymers (i)

There are no such residues in this entry.