



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:46 PM GMT

PDB ID : 4PSO
Title : Crystal structure of apeThermo-DBP-RP2 bound to ssDNA dT10
Authors : Gahlei, H.; von Moeller, H.; Eppers, D.; Loll, B.; Wahl, M.C.
Deposited on : 2014-03-07
Resolution : 2.90 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 ⓘ 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

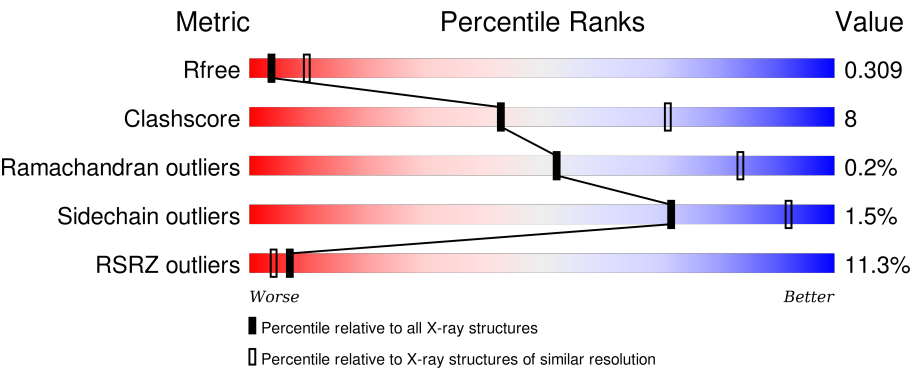
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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 $\leq 5\%$. 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	237	<div><div>9%</div><div>72%</div><div>19%</div><div>• 8%</div></div>
1	B	237	<div><div>8%</div><div>74%</div><div>18%</div><div>7%</div></div>
1	C	237	<div><div>3%</div><div>72%</div><div>20%</div><div>• 7%</div></div>
1	D	237	<div><div>9%</div><div>70%</div><div>21%</div><div>• 8%</div></div>
1	F	237	<div><div>12%</div><div>75%</div><div>17%</div><div>8%</div></div>

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Mol	Chain	Length	Quality of chain
1	G	237	<div><div></div><div>14%</div><div>72%</div><div>19%</div><div>8%</div></div>
1	H	237	<div><div></div><div>12%</div><div>78%</div><div>14%</div><div>7%</div></div>
1	I	237	<div><div></div><div>15%</div><div>78%</div><div>15%</div><div>7%</div></div>
2	E	10	<div><div></div><div>10%</div><div>20%</div><div>40%</div><div>10%</div><div>30%</div></div>
2	L	10	<div><div></div><div>10%</div><div>20%</div><div>40%</div><div>40%</div></div>
2	X	10	<div><div></div><div>20%</div><div>20%</div><div>60%</div></div>
2	Z	10	<div><div></div><div>20%</div><div>20%</div><div>60%</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14476 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 ssDNA binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	0	0	0
			1734	1097	310	324	3			
1	B	220	Total	C	N	O	S	0	0	0
			1748	1106	313	326	3			
1	C	220	Total	C	N	O	S	0	0	0
			1748	1106	313	326	3			
1	D	219	Total	C	N	O	S	0	0	0
			1739	1100	311	325	3			
1	F	219	Total	C	N	O	S	0	0	0
			1741	1102	312	324	3			
1	G	217	Total	C	N	O	S	0	0	0
			1723	1091	306	323	3			
1	H	220	Total	C	N	O	S	0	0	0
			1748	1106	313	326	3			
1	I	221	Total	C	N	O	S	0	0	0
			1752	1108	314	327	3			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q9YAS7
A	-1	ALA	-	EXPRESSION TAG	UNP Q9YAS7
A	1	LEU	-	EXPRESSION TAG	UNP Q9YAS7
B	-2	GLY	-	EXPRESSION TAG	UNP Q9YAS7
B	-1	ALA	-	EXPRESSION TAG	UNP Q9YAS7
B	1	LEU	-	EXPRESSION TAG	UNP Q9YAS7
C	-2	GLY	-	EXPRESSION TAG	UNP Q9YAS7
C	-1	ALA	-	EXPRESSION TAG	UNP Q9YAS7
C	1	LEU	-	EXPRESSION TAG	UNP Q9YAS7
D	-2	GLY	-	EXPRESSION TAG	UNP Q9YAS7
D	-1	ALA	-	EXPRESSION TAG	UNP Q9YAS7
D	1	LEU	-	EXPRESSION TAG	UNP Q9YAS7
F	-2	GLY	-	EXPRESSION TAG	UNP Q9YAS7

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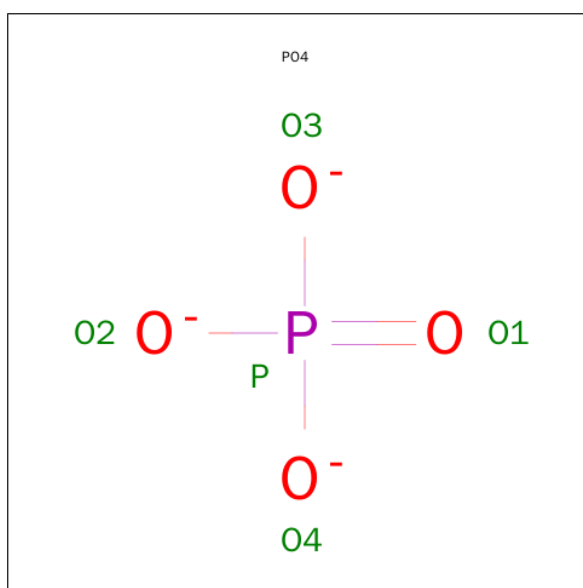
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Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	ALA	-	EXPRESSION TAG	UNP Q9YAS7
F	1	LEU	-	EXPRESSION TAG	UNP Q9YAS7
G	-2	GLY	-	EXPRESSION TAG	UNP Q9YAS7
G	-1	ALA	-	EXPRESSION TAG	UNP Q9YAS7
G	1	LEU	-	EXPRESSION TAG	UNP Q9YAS7
H	-2	GLY	-	EXPRESSION TAG	UNP Q9YAS7
H	-1	ALA	-	EXPRESSION TAG	UNP Q9YAS7
H	1	LEU	-	EXPRESSION TAG	UNP Q9YAS7
I	-2	GLY	-	EXPRESSION TAG	UNP Q9YAS7
I	-1	ALA	-	EXPRESSION TAG	UNP Q9YAS7
I	1	LEU	-	EXPRESSION TAG	UNP Q9YAS7

- Molecule 2 is a DNA chain called polydeoxyribonucleotide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	10	Total	C	N	O	P	0	0	0
			200	100	20	70	10			
2	E	7	Total	C	N	O	P	0	0	0
			140	70	14	49	7			
2	X	4	Total	C	N	O	P	0	0	0
			80	40	8	28	4			
2	Z	4	Total	C	N	O	P	0	0	0
			80	40	8	28	4			

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		

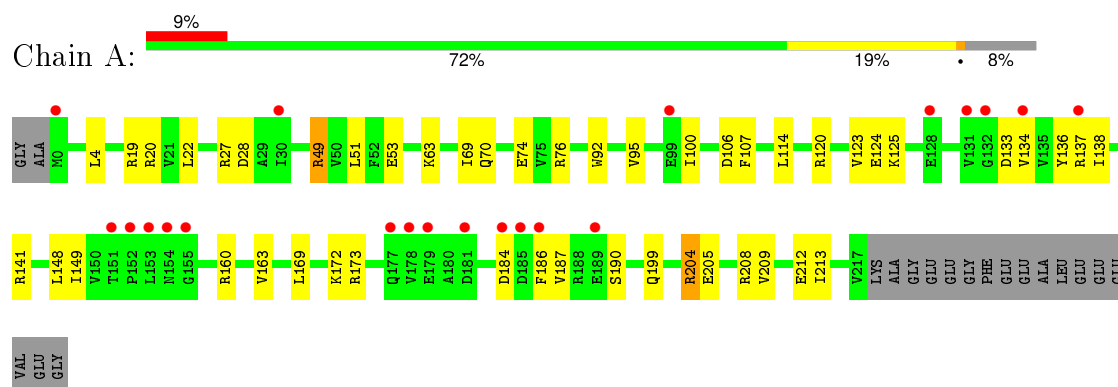
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	O	0	0
			8	8		
4	B	6	Total	O	0	0
			6	6		
4	C	6	Total	O	0	0
			6	6		
4	D	5	Total	O	0	0
			5	5		
4	F	3	Total	O	0	0
			3	3		
4	G	4	Total	O	0	0
			4	4		
4	H	1	Total	O	0	0
			1	1		
4	I	3	Total	O	0	0
			3	3		
4	E	1	Total	O	0	0
			1	1		
4	X	1	Total	O	0	0
			1	1		

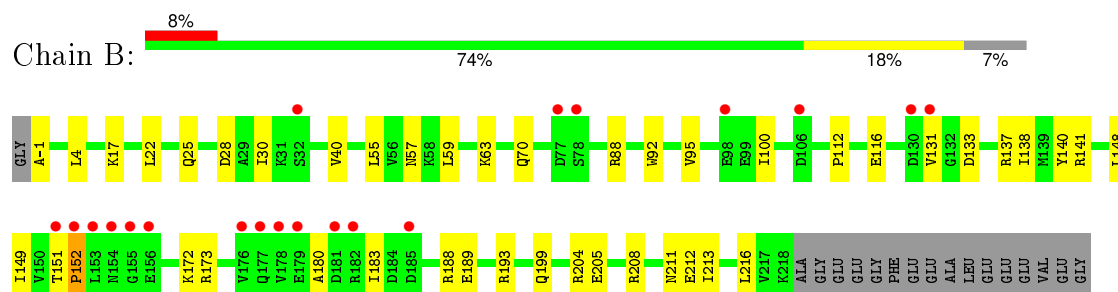
3 Residue-property plots [i](#)

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.

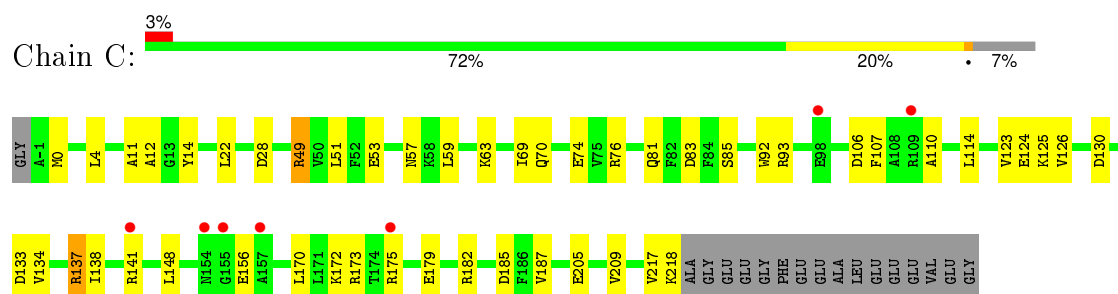
- Molecule 1: ssDNA binding protein



- Molecule 1: ssDNA binding protein

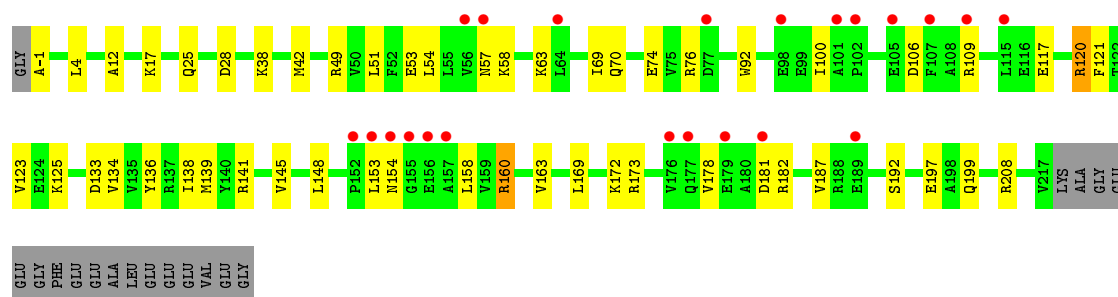


- Molecule 1: ssDNA binding protein

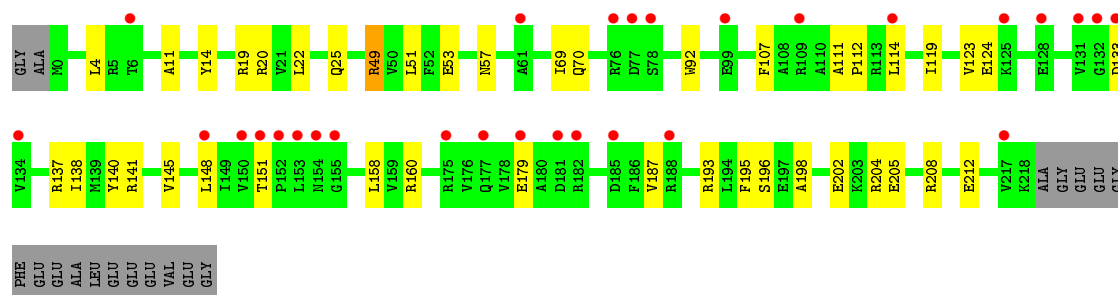
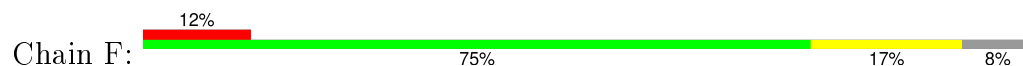


- Molecule 1: ssDNA binding protein

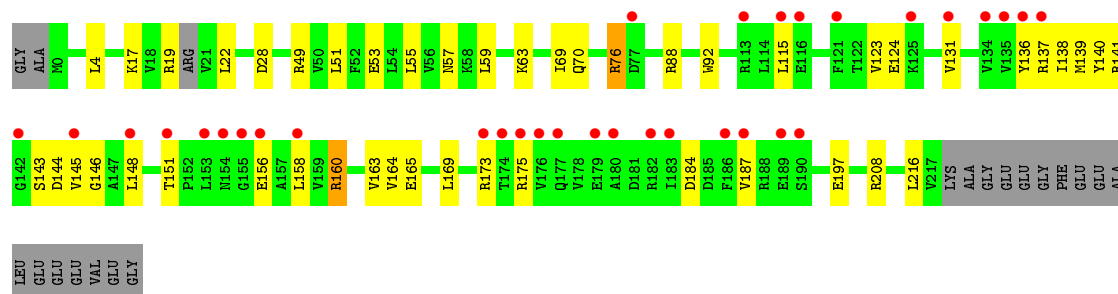
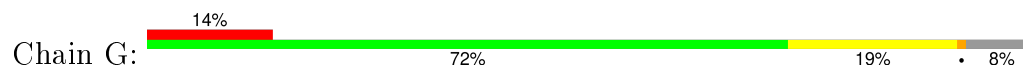




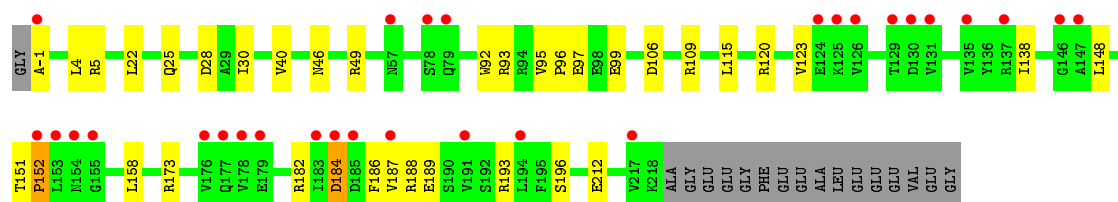
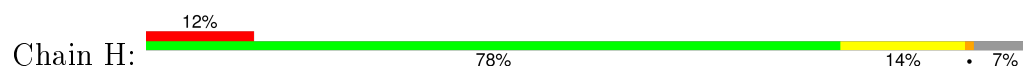
- Molecule 1: ssDNA binding protein



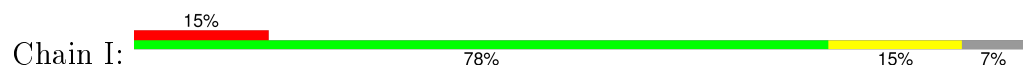
- Molecule 1: ssDNA binding protein

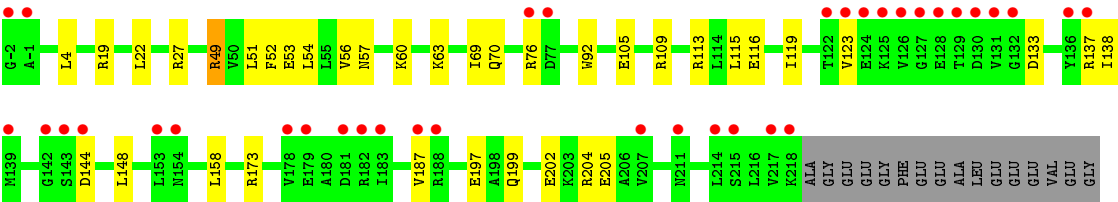


- Molecule 1: ssDNA binding protein

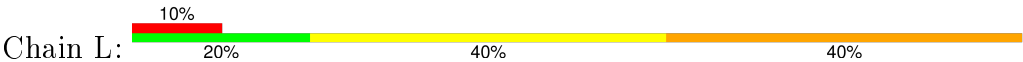


- Molecule 1: ssDNA binding protein

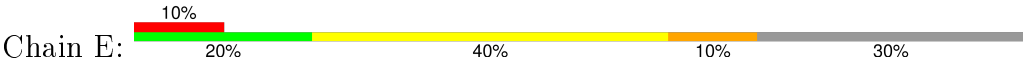




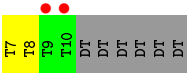
● Molecule 2: polydeoxyribonucleotide



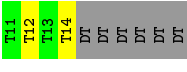
● Molecule 2: polydeoxyribonucleotide



● Molecule 2: polydeoxyribonucleotide



● Molecule 2: polydeoxyribonucleotide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	120.49 Å 190.28 Å 89.79 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.93 – 2.90 48.38 – 2.32	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.93-2.90) 90.9 (48.38-2.32)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 2.32 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1389)	Depositor
R, R_{free}	0.267 , 0.309 0.266 , 0.309	Depositor DCC
R_{free} test set	2322 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	2 of 81674 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	14476	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.63% 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 ⓘ

5.1 Standard geometry ⓘ

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

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	0.25	0/1754	0.47	0/2373
1	B	0.25	0/1768	0.48	0/2391
1	C	0.26	0/1768	0.49	0/2391
1	D	0.26	0/1759	0.51	0/2380
1	F	0.25	0/1761	0.46	0/2381
1	G	0.25	0/1742	0.47	0/2356
1	H	0.24	0/1768	0.43	0/2391
1	I	0.24	0/1772	0.45	0/2396
2	E	0.51	0/153	1.69	3/234 (1.3%)
2	L	0.52	0/219	1.72	7/336 (2.1%)
2	X	0.56	0/87	1.55	1/132 (0.8%)
2	Z	0.58	0/87	1.55	0/132
All	All	0.27	0/14638	0.57	11/19893 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	5	DT	O4'-C1'-N1	7.78	113.44	108.00
2	E	2	DT	O4'-C1'-N1	7.34	113.14	108.00
2	X	7	DT	O4'-C1'-N1	6.02	112.22	108.00
2	L	3	DT	C6-C5-C7	-5.85	119.39	122.90
2	L	5	DT	O4'-C4'-C3'	-5.63	102.25	104.50
2	L	1	DT	O4'-C1'-C2'	-5.59	101.43	105.90
2	E	0	DT	C5-C4-O4	-5.48	121.06	124.90
2	E	0	DT	N3-C4-O4	5.39	123.13	119.90
2	L	8	DT	N3-C4-O4	5.16	123.00	119.90
2	L	1	DT	N1-C1'-C2'	5.01	122.13	112.60
2	L	3	DT	C4-C5-C7	5.00	122.00	119.00

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	1734	0	1800	39	0
1	B	1748	0	1818	41	0
1	C	1748	0	1818	37	0
1	D	1739	0	1805	44	0
1	F	1741	0	1808	37	1
1	G	1723	0	1786	41	1
1	H	1748	0	1818	29	0
1	I	1752	0	1821	31	0
2	E	140	0	85	3	0
2	L	200	0	121	12	0
2	X	80	0	49	2	0
2	Z	80	0	49	2	0
3	B	5	0	0	1	0
4	A	8	0	0	1	0
4	B	6	0	0	1	0
4	C	6	0	0	0	0
4	D	5	0	0	2	0
4	E	1	0	0	0	0
4	F	3	0	0	0	0
4	G	4	0	0	0	0
4	H	1	0	0	0	0
4	I	3	0	0	2	0
4	X	1	0	0	0	0
All	All	14476	0	14778	239	1

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 8.

All (239) 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:G:137:ARG:NH2	1:G:146:GLY:O	2.03	0.92
1:H:5:ARG:HD3	1:I:27:ARG:HH12	1.41	0.83
1:C:63:LYS:HD3	2:L:3:DT:H5'	1.64	0.78
1:G:137:ARG:NH2	1:G:138:ILE:O	2.19	0.76
1:D:138:ILE:HD11	1:D:148:LEU:HB2	1.69	0.75
1:F:70:GLN:NE2	1:G:28:ASP:OD2	2.17	0.74
1:B:133:ASP:OD2	1:D:173:ARG:NH1	2.20	0.74
1:F:138:ILE:HD11	1:F:148:LEU:HB2	1.71	0.72
1:A:70:GLN:NE2	1:B:28:ASP:OD1	2.15	0.71
1:G:138:ILE:HD11	1:G:148:LEU:HB2	1.72	0.71
1:B:193:ARG:HD2	1:H:93:ARG:HD3	1.74	0.69
1:A:173:ARG:NH1	1:C:133:ASP:OD2	2.26	0.69
1:H:184:ASP:O	1:H:188:ARG:NH2	2.26	0.68
1:G:137:ARG:HH11	1:G:144:ASP:HB3	1.59	0.68
1:A:136:TYR:OH	1:A:184:ASP:OD1	2.11	0.68
1:G:123:VAL:HG21	1:G:187:VAL:HG12	1.75	0.68
1:H:138:ILE:HD11	1:H:148:LEU:HB2	1.75	0.67
1:B:17:LYS:HZ2	2:L:9:DT:H3	1.43	0.66
1:G:53:GLU:OE1	1:G:57:ASN:ND2	2.29	0.66
1:C:53:GLU:OE1	1:C:57:ASN:ND2	2.29	0.66
1:G:137:ARG:NH1	1:G:144:ASP:HB3	2.11	0.65
1:C:4:LEU:HD12	1:C:22:LEU:HD13	1.77	0.65
1:B:173:ARG:NH2	1:D:133:ASP:OD2	2.20	0.65
1:I:4:LEU:HD12	1:I:22:LEU:HD13	1.79	0.64
1:A:125:LYS:HE3	1:A:134:VAL:HG13	1.80	0.64
1:D:53:GLU:OE1	1:D:57:ASN:ND2	2.31	0.63
1:C:81:GLN:HE22	1:G:124:GLU:HB2	1.62	0.63
1:A:106:ASP:OD1	1:A:107:PHE:N	2.31	0.63
1:D:106:ASP:OD1	1:D:109:ARG:NH2	2.30	0.63
1:G:156:GLU:OE2	1:G:175:ARG:NH1	2.31	0.63
1:B:92:TRP:HB3	1:C:92:TRP:HB3	1.80	0.62
1:G:4:LEU:HD12	1:G:22:LEU:HD13	1.80	0.62
1:I:138:ILE:HD11	1:I:148:LEU:HB2	1.81	0.62
1:C:106:ASP:OD1	1:C:107:PHE:N	2.33	0.62
1:B:208:ARG:NH2	1:D:199:GLN:OE1	2.32	0.62
1:F:53:GLU:OE1	1:F:57:ASN:ND2	2.32	0.62
1:G:131:VAL:HG21	1:I:158:LEU:HD11	1.83	0.61
1:I:76:ARG:NH2	4:I:303:HOH:O	2.28	0.61
1:A:114:LEU:HD21	1:D:53:GLU:HG3	1.81	0.61
1:D:160:ARG:HB3	1:D:160:ARG:HH21	1.65	0.61
1:D:123:VAL:HG21	1:D:187:VAL:HG12	1.83	0.61
1:G:136:TYR:OH	1:G:184:ASP:OD1	2.17	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:TRP:HB3	1:D:92:TRP:HB3	1.82	0.60
1:G:160:ARG:HG3	1:G:173:ARG:HA	1.84	0.60
1:I:53:GLU:OE1	1:I:57:ASN:ND2	2.33	0.60
1:D:25:GLN:NE2	4:D:305:HOH:O	2.35	0.60
1:B:4:LEU:HD12	1:B:22:LEU:HD13	1.84	0.59
1:F:4:LEU:HD12	1:F:22:LEU:HD13	1.84	0.59
1:B:193:ARG:CD	1:H:93:ARG:HD3	2.33	0.59
1:H:25:GLN:HE21	2:Z:12:DT:H3	1.51	0.59
1:B:189:GLU:HG2	1:H:95:VAL:HG13	1.83	0.59
1:A:123:VAL:HG21	1:A:187:VAL:HG12	1.85	0.59
1:A:133:ASP:OD2	1:C:173:ARG:NH1	2.36	0.58
1:A:28:ASP:OD2	1:B:70:GLN:NE2	2.29	0.58
1:A:53:GLU:OE1	1:D:141:ARG:NH2	2.36	0.58
1:A:204:ARG:NH1	1:C:205:GLU:OE1	2.35	0.58
1:C:130:ASP:OD1	1:C:218:LYS:HD3	2.03	0.58
1:G:173:ARG:NH1	1:I:133:ASP:OD2	2.34	0.58
1:F:160:ARG:NH2	1:H:212:GLU:OE1	2.35	0.57
1:D:58:LYS:NZ	4:D:303:HOH:O	2.37	0.57
1:I:123:VAL:HG21	1:I:187:VAL:HG12	1.86	0.57
1:C:126:VAL:HG12	1:C:217:VAL:HG11	1.85	0.57
1:F:123:VAL:HG21	1:F:187:VAL:HG12	1.86	0.56
1:D:125:LYS:HE3	1:D:134:VAL:HG13	1.87	0.56
1:G:145:VAL:HG12	1:G:165:GLU:HB2	1.87	0.56
1:F:20:ARG:NH1	2:Z:14:DT:O3'	2.30	0.56
1:C:28:ASP:OD2	1:D:70:GLN:NE2	2.25	0.56
1:B:172:LYS:HD2	1:B:199:GLN:NE2	2.21	0.56
1:C:123:VAL:HG21	1:C:187:VAL:HG12	1.88	0.56
1:D:74:GLU:OE2	1:D:76:ARG:NE	2.37	0.56
1:C:12:ALA:HB3	2:L:4:DT:H1'	1.89	0.55
1:H:-1:ALA:N	4:I:303:HOH:O	2.38	0.55
1:G:115:LEU:HG	1:H:46:ASN:HB3	1.87	0.55
1:H:123:VAL:HG21	1:H:187:VAL:HG12	1.88	0.55
1:A:205:GLU:HG2	1:C:205:GLU:HG2	1.89	0.55
1:F:25:GLN:HE21	2:X:8:DT:H3	1.53	0.55
1:B:55:LEU:HD23	1:B:59:LEU:HD12	1.88	0.55
1:F:107:PHE:CE2	1:I:54:LEU:HB2	2.41	0.55
1:C:125:LYS:HE3	1:C:134:VAL:HG13	1.89	0.55
1:F:107:PHE:CD2	1:I:54:LEU:HB2	2.42	0.55
1:F:151:THR:HB	1:F:158:LEU:HB2	1.89	0.55
1:G:140:TYR:OH	1:G:141:ARG:NH2	2.39	0.55
1:C:138:ILE:HD11	1:C:148:LEU:HB2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LEU:HD12	1:A:22:LEU:HD13	1.88	0.55
1:B:17:LYS:NZ	2:L:9:DT:H3	2.04	0.55
1:F:20:ARG:HH22	1:I:204:ARG:CZ	2.19	0.54
1:H:158:LEU:HD13	1:H:173:ARG:HE	1.72	0.54
1:A:49:ARG:NH1	1:D:117:GLU:O	2.40	0.54
1:D:160:ARG:HB3	1:D:160:ARG:NH2	2.23	0.54
1:F:119:ILE:HG12	1:I:49:ARG:NH1	2.23	0.54
1:B:112:PRO:HB2	1:G:144:ASP:OD2	2.08	0.54
1:D:178:VAL:HG13	1:D:182:ARG:HB2	1.89	0.54
1:G:92:TRP:HB3	1:H:92:TRP:HB3	1.90	0.54
1:B:211:ASN:ND2	2:L:1:DT:C2	2.76	0.53
1:D:121:PHE:N	1:D:192:SER:OG	2.34	0.53
1:C:81:GLN:NE2	1:G:124:GLU:HB2	2.23	0.53
1:G:208:ARG:NH1	1:I:199:GLN:OE1	2.40	0.53
2:E:-3:DT:H5'	2:E:-3:DT:H6	1.73	0.53
1:H:4:LEU:HD12	1:H:22:LEU:HD13	1.89	0.53
1:A:120:ARG:HB2	1:A:120:ARG:CZ	2.38	0.53
1:A:141:ARG:NH2	1:D:53:GLU:OE1	2.41	0.52
1:D:120:ARG:HA	1:D:192:SER:HB3	1.92	0.52
1:H:182:ARG:O	1:H:186:PHE:N	2.36	0.52
1:H:186:PHE:O	1:H:189:GLU:HG2	2.09	0.52
1:B:216:LEU:HD11	1:D:173:ARG:HD3	1.91	0.52
1:B:193:ARG:HD2	1:H:93:ARG:HH11	1.74	0.52
1:G:124:GLU:HB3	1:G:137:ARG:HB3	1.91	0.52
1:F:202:GLU:HG2	1:I:19:ARG:HD3	1.91	0.52
1:H:106:ASP:HA	1:H:109:ARG:HH11	1.75	0.52
1:C:12:ALA:HB2	1:C:63:LYS:HB2	1.92	0.51
1:F:140:TYR:OH	1:F:141:ARG:NH2	2.44	0.51
1:B:141:ARG:NH2	1:C:53:GLU:OE2	2.38	0.51
1:B:188:ARG:HB3	1:H:99:GLU:OE1	2.10	0.51
2:E:-1:DT:H4'	2:E:0:DT:OP1	2.10	0.51
1:C:81:GLN:NE2	1:G:139:MET:SD	2.83	0.51
1:D:38:LYS:O	1:D:42:MET:HB2	2.10	0.51
1:B:57:ASN:OD1	1:C:141:ARG:NH2	2.44	0.51
1:G:163:VAL:HG22	1:G:169:LEU:HB2	1.92	0.51
1:A:51:LEU:HD13	1:A:69:ILE:HD13	1.93	0.50
1:H:106:ASP:OD1	1:H:109:ARG:NH1	2.45	0.50
1:G:145:VAL:O	1:G:164:VAL:N	2.37	0.50
1:A:76:ARG:CZ	1:B:-1:ALA:HB2	2.42	0.50
1:D:172:LYS:HD2	1:D:199:GLN:NE2	2.27	0.50
1:F:92:TRP:HB3	1:I:92:TRP:HB3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:51:LEU:HD13	1:G:69:ILE:HD13	1.93	0.50
1:F:124:GLU:OE2	1:F:137:ARG:HD2	2.12	0.50
2:L:3:DT:H2"	2:L:4:DT:H5"	1.93	0.49
1:I:113:ARG:O	1:I:116:GLU:HB3	2.13	0.49
1:D:17:LYS:HZ2	2:E:1:DT:H3	1.59	0.49
1:I:137:ARG:HD3	1:I:144:ASP:OD1	2.12	0.49
1:I:51:LEU:HD13	1:I:69:ILE:HD13	1.93	0.49
1:A:172:LYS:HE3	1:A:199:GLN:OE1	2.13	0.49
1:F:124:GLU:OE2	1:F:137:ARG:NH1	2.46	0.49
1:B:131:VAL:HB	1:D:153:LEU:HD13	1.95	0.49
1:F:140:TYR:CZ	1:F:141:ARG:NH2	2.81	0.48
1:F:114:LEU:HD21	1:I:53:GLU:HG3	1.93	0.48
1:A:163:VAL:HG22	1:A:169:LEU:HB2	1.96	0.48
1:F:205:GLU:HA	1:F:208:ARG:NH1	2.28	0.48
2:L:8:DT:O5'	2:L:9:DT:H72	2.14	0.48
1:D:12:ALA:HB2	1:D:63:LYS:HB2	1.95	0.48
1:C:156:GLU:OE2	1:C:175:ARG:NH1	2.47	0.48
1:G:137:ARG:HH21	1:G:138:ILE:N	2.11	0.48
1:G:216:LEU:HD13	1:I:173:ARG:HH21	1.79	0.48
1:G:197:GLU:HG2	1:G:197:GLU:O	2.14	0.47
1:F:160:ARG:HH21	1:H:212:GLU:CD	2.18	0.47
1:C:74:GLU:HG2	1:C:76:ARG:HG2	1.95	0.47
1:F:193:ARG:O	1:F:196:SER:OG	2.22	0.47
1:F:4:LEU:O	1:F:70:GLN:HA	2.15	0.47
1:A:4:LEU:O	1:A:70:GLN:HA	2.14	0.47
1:B:172:LYS:HD2	1:B:199:GLN:HE21	1.79	0.47
1:D:139:MET:HA	1:D:145:VAL:HG22	1.97	0.47
1:A:208:ARG:HG3	1:A:209:VAL:N	2.28	0.47
1:B:30:ILE:HD11	1:B:40:VAL:HG21	1.97	0.47
1:B:116:GLU:OE2	1:G:143:SER:OG	2.28	0.46
1:A:138:ILE:HD11	1:A:148:LEU:HB2	1.96	0.46
1:F:205:GLU:HA	1:F:208:ARG:HH12	1.80	0.46
1:H:115:LEU:HD23	1:H:115:LEU:HA	1.76	0.46
2:L:6:DT:H3'	2:L:7:DT:H5"	1.96	0.46
1:B:149:ILE:HB	1:B:213:ILE:HG23	1.96	0.46
1:F:51:LEU:HD13	1:F:69:ILE:HD13	1.98	0.46
1:H:30:ILE:HD11	1:H:40:VAL:HG21	1.97	0.46
1:B:140:TYR:OH	1:C:49:ARG:NH1	2.48	0.46
1:B:131:VAL:HG21	1:D:158:LEU:HD11	1.98	0.46
1:F:20:ARG:HH22	1:I:204:ARG:NH2	2.13	0.46
1:A:74:GLU:OE2	1:A:76:ARG:NE	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:158:LEU:HD13	1:G:173:ARG:HE	1.80	0.45
1:F:140:TYR:HB3	1:F:145:VAL:HG11	1.99	0.45
1:A:124:GLU:OE2	1:A:137:ARG:HD2	2.16	0.45
1:A:95:VAL:HG11	1:A:100:ILE:HD11	1.98	0.45
1:F:204:ARG:NH2	1:G:19:ARG:HH22	2.15	0.45
1:C:182:ARG:NH1	1:C:185:ASP:OD2	2.49	0.45
1:I:197:GLU:O	1:I:197:GLU:HG2	2.17	0.45
1:A:20:ARG:HG2	2:L:6:DT:H2''	1.99	0.45
1:G:88:ARG:NE	1:H:97:GLU:OE1	2.50	0.45
1:A:28:ASP:CG	1:B:88:ARG:HH22	2.20	0.45
1:F:208:ARG:O	1:F:212:GLU:HG3	2.17	0.45
1:C:11:ALA:O	1:C:14:TYR:HD2	2.00	0.45
4:A:308:HOH:O	1:D:100:ILE:HG21	2.17	0.44
1:I:56:VAL:O	1:I:60:LYS:HD3	2.17	0.44
1:F:19:ARG:HD3	1:I:202:GLU:HG2	1.99	0.44
1:A:212:GLU:OE1	1:C:172:LYS:HE2	2.17	0.44
2:L:5:DT:H2'	2:L:6:DT:C6	2.53	0.44
1:H:96:PRO:HB2	1:H:99:GLU:HG2	1.99	0.44
1:D:4:LEU:O	1:D:70:GLN:HA	2.18	0.44
1:A:186:PHE:O	1:A:190:SER:OG	2.21	0.44
1:C:63:LYS:HB3	1:C:63:LYS:HE2	1.70	0.44
1:C:76:ARG:HH21	1:D:-1:ALA:HB2	1.83	0.44
1:B:205:GLU:OE2	1:D:208:ARG:NE	2.48	0.44
1:G:137:ARG:HH21	1:G:138:ILE:H	1.66	0.43
1:H:28:ASP:OD2	1:I:70:GLN:NE2	2.44	0.43
1:C:51:LEU:HD13	1:C:69:ILE:HD13	2.00	0.43
1:I:115:LEU:HA	1:I:115:LEU:HD23	1.79	0.43
1:I:105:GLU:HG2	1:I:109:ARG:NH1	2.34	0.43
1:B:95:VAL:HG11	1:B:100:ILE:HD11	2.00	0.43
1:C:83:ASP:OD1	1:C:85:SER:OG	2.27	0.43
1:D:53:GLU:O	1:D:57:ASN:HB2	2.18	0.43
1:G:17:LYS:NZ	2:X:8:DT:OP1	2.47	0.43
1:I:105:GLU:HG2	1:I:109:ARG:HH12	1.83	0.43
1:F:49:ARG:CZ	1:I:119:ILE:HG12	2.49	0.43
1:A:141:ARG:HH22	1:D:57:ASN:ND2	2.16	0.43
1:F:195:PHE:HA	1:F:198:ALA:HB2	2.00	0.43
1:B:216:LEU:HD13	1:D:173:ARG:NH2	2.34	0.42
1:G:55:LEU:HD23	1:G:59:LEU:HD12	2.01	0.42
1:G:63:LYS:HE2	1:G:63:LYS:HB3	1.83	0.42
1:G:4:LEU:O	1:G:70:GLN:HA	2.20	0.42
1:B:25:GLN:NE2	4:B:406:HOH:O	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:PHE:CD1	1:D:54:LEU:HB2	2.54	0.42
1:D:136:TYR:CE1	1:D:187:VAL:HG21	2.55	0.42
1:A:28:ASP:OD1	1:B:88:ARG:NH2	2.46	0.42
1:H:151:THR:HA	1:H:152:PRO:HD3	1.91	0.42
1:B:151:THR:HA	1:B:152:PRO:HD3	1.87	0.42
1:C:59:LEU:O	1:C:93:ARG:NH2	2.52	0.42
1:F:179:GLU:N	1:F:179:GLU:OE1	2.53	0.42
1:C:110:ALA:O	1:C:114:LEU:HG	2.20	0.42
1:B:212:GLU:O	1:B:216:LEU:HG	2.20	0.42
1:B:180:ALA:HA	1:B:183:ILE:HG13	2.02	0.42
1:I:52:PHE:CE1	1:I:56:VAL:HG21	2.55	0.42
1:B:138:ILE:HD11	1:B:148:LEU:HB2	2.01	0.42
1:C:124:GLU:CD	1:C:137:ARG:HH21	2.23	0.42
2:L:4:DT:H2'	2:L:5:DT:O4'	2.20	0.41
1:G:151:THR:HB	1:G:158:LEU:HB2	2.01	0.41
1:A:149:ILE:HB	1:A:213:ILE:HG23	2.02	0.41
1:C:70:GLN:NE2	1:D:28:ASP:OD2	2.33	0.41
1:D:51:LEU:HD13	1:D:69:ILE:HD13	2.02	0.41
1:I:4:LEU:O	1:I:70:GLN:HA	2.20	0.41
1:H:193:ARG:O	1:H:196:SER:OG	2.30	0.41
1:B:137:ARG:HD2	1:B:137:ARG:HA	1.75	0.41
1:F:11:ALA:O	1:F:14:TYR:HD2	2.04	0.41
1:D:163:VAL:HG22	1:D:169:LEU:HB2	2.01	0.41
1:C:170:LEU:HD11	1:C:209:VAL:HG21	2.03	0.41
1:G:146:GLY:HA3	1:G:163:VAL:HA	2.03	0.41
1:F:111:ALA:HB3	1:F:112:PRO:HD3	2.03	0.41
1:A:19:ARG:NH1	3:B:301:PO4:O4	2.53	0.41
1:B:4:LEU:O	1:B:70:GLN:HA	2.21	0.41
1:A:27:ARG:NH2	2:L:5:DT:O3'	2.54	0.40
1:A:107:PHE:CE1	1:D:54:LEU:HB2	2.57	0.40
1:D:181:ASP:OD1	1:D:182:ARG:HG2	2.21	0.40
1:A:63:LYS:HE2	1:A:63:LYS:HB3	1.80	0.40
1:F:133:ASP:OD2	1:H:173:ARG:NH1	2.55	0.40

All (1) 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)
1:F:179:GLU:OE1	1:G:76:ARG:NH1[2_855]	2.18	0.02

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	216/237 (91%)	211 (98%)	5 (2%)	0	100	100
1	B	218/237 (92%)	213 (98%)	5 (2%)	0	100	100
1	C	218/237 (92%)	213 (98%)	4 (2%)	1 (0%)	34	71
1	D	217/237 (92%)	210 (97%)	5 (2%)	2 (1%)	21	57
1	F	217/237 (92%)	212 (98%)	5 (2%)	0	100	100
1	G	213/237 (90%)	208 (98%)	5 (2%)	0	100	100
1	H	218/237 (92%)	213 (98%)	5 (2%)	0	100	100
1	I	219/237 (92%)	214 (98%)	5 (2%)	0	100	100
All	All	1736/1896 (92%)	1694 (98%)	39 (2%)	3 (0%)	52	84

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	154	ASN
1	C	0	MET
1	D	120	ARG

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	189/201 (94%)	186 (98%)	3 (2%)	70	91
1	B	190/201 (94%)	187 (98%)	3 (2%)	70	91

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	190/201 (94%)	187 (98%)	3 (2%)	70	91
1	D	189/201 (94%)	186 (98%)	3 (2%)	70	91
1	F	189/201 (94%)	188 (100%)	1 (0%)	92	98
1	G	188/201 (94%)	185 (98%)	3 (2%)	70	91
1	H	190/201 (94%)	186 (98%)	4 (2%)	61	88
1	I	190/201 (94%)	187 (98%)	3 (2%)	70	91
All	All	1515/1608 (94%)	1492 (98%)	23 (2%)	72	92

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ARG
1	A	160	ARG
1	A	204	ARG
1	B	63	LYS
1	B	152	PRO
1	B	204	ARG
1	C	49	ARG
1	C	137	ARG
1	C	179	GLU
1	D	49	ARG
1	D	160	ARG
1	D	197	GLU
1	F	49	ARG
1	G	49	ARG
1	G	76	ARG
1	G	160	ARG
1	H	49	ARG
1	H	120	ARG
1	H	152	PRO
1	H	184	ASP
1	I	49	ARG
1	I	63	LYS
1	I	205	GLU

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

Mol	Chain	Res	Type
1	C	81	GLN

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Mol	Chain	Res	Type
1	D	57	ASN
1	I	79	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 ⓘ

1 ligand is 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	$\# Z > 2$
3	PO4	B	301	-	4,4,4	0.48	0	6,6,6	0.27	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
3	PO4	B	301	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	301	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	218/237 (91%)	0.58	21 (9%) 10 6	21, 47, 84, 101	0
1	B	220/237 (92%)	0.39	20 (9%) 11 7	21, 45, 80, 103	0
1	C	220/237 (92%)	0.22	7 (3%) 51 43	21, 43, 77, 92	0
1	D	219/237 (92%)	0.35	22 (10%) 9 5	20, 45, 93, 121	0
1	F	219/237 (92%)	0.68	29 (13%) 4 2	36, 63, 100, 113	0
1	G	217/237 (91%)	0.82	33 (15%) 3 1	35, 68, 102, 120	0
1	H	220/237 (92%)	0.81	29 (13%) 4 2	35, 66, 96, 115	0
1	I	221/237 (93%)	0.77	36 (16%) 2 1	40, 64, 98, 123	0
2	E	7/10 (70%)	1.76	1 (14%) 4 2	64, 78, 121, 131	0
2	L	10/10 (100%)	1.40	1 (10%) 9 5	58, 85, 104, 112	0
2	X	4/10 (40%)	1.94	2 (50%) 0 0	92, 100, 111, 127	0
2	Z	4/10 (40%)	1.52	0 100 100	91, 96, 104, 112	0
All	All	1779/1936 (91%)	0.59	201 (11%) 7 4	20, 56, 97, 131	0

All (201) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	185	ASP	8.2
1	I	130	ASP	7.8
1	I	-2	GLY	7.3
1	G	175	ARG	6.6
1	G	154	ASN	6.5
1	C	154	ASN	6.5
1	B	154	ASN	6.4
1	A	155	GLY	6.1
1	C	98	GLU	6.0
1	A	154	ASN	5.9
1	H	126	VAL	5.8

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Mol	Chain	Res	Type	RSRZ
1	H	154	ASN	5.8
1	G	183	ILE	5.7
1	I	137	ARG	5.5
1	D	177	GLN	5.5
1	F	154	ASN	5.4
1	G	179	GLU	5.4
1	H	79	GLN	5.4
1	I	218	LYS	5.3
1	H	153	LEU	5.3
1	D	98	GLU	5.2
1	H	129	THR	5.2
1	H	131	VAL	5.2
1	D	155	GLY	5.1
1	H	178	VAL	5.0
1	I	129	THR	5.0
1	D	154	ASN	5.0
1	G	135	VAL	4.8
1	I	77	ASP	4.8
1	G	155	GLY	4.7
1	H	135	VAL	4.7
1	G	153	LEU	4.7
1	H	-1	ALA	4.7
1	B	153	LEU	4.6
1	F	132	GLY	4.6
1	B	155	GLY	4.5
1	H	155	GLY	4.5
1	I	124	GLU	4.5
1	G	136	TYR	4.5
1	B	78	SER	4.5
1	I	126	VAL	4.5
1	G	156	GLU	4.4
1	F	177	GLN	4.3
1	B	179	GLU	4.3
1	A	178	VAL	4.3
2	E	3	DT	4.3
1	H	130	ASP	4.3
1	F	153	LEU	4.2
1	I	214	LEU	4.2
1	H	185	ASP	4.2
1	I	181	ASP	4.2
1	F	131	VAL	4.1
1	G	151	THR	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	0	MET	4.1
1	D	156	GLU	4.1
1	G	158	LEU	4.0
1	D	77	ASP	4.0
1	G	177	GLN	4.0
1	B	131	VAL	4.0
1	H	78	SER	3.9
1	H	124	GLU	3.8
1	D	152	PRO	3.8
1	A	179	GLU	3.8
1	C	155	GLY	3.7
1	H	146	GLY	3.7
1	H	176	VAL	3.6
1	B	176	VAL	3.6
1	I	217	VAL	3.6
1	G	189	GLU	3.6
1	D	181	ASP	3.6
1	G	176	VAL	3.6
1	D	109	ARG	3.6
1	A	177	GLN	3.6
1	F	181	ASP	3.6
1	A	152	PRO	3.6
1	H	152	PRO	3.5
1	F	99	GLU	3.5
1	G	180	ALA	3.5
1	D	64	LEU	3.4
1	I	128	GLU	3.4
1	H	187	VAL	3.4
1	F	134	VAL	3.4
1	F	179	GLU	3.4
1	F	77	ASP	3.4
1	G	182	ARG	3.4
1	G	142	GLY	3.3
1	I	182	ARG	3.3
1	I	132	GLY	3.3
1	H	179	GLU	3.2
1	G	186	PHE	3.2
1	D	153	LEU	3.2
1	B	156	GLU	3.2
1	D	105	GLU	3.2
1	B	181	ASP	3.2
1	B	32	SER	3.2

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Mol	Chain	Res	Type	RSRZ
1	I	143	SER	3.1
1	A	134	VAL	3.0
1	F	152	PRO	3.0
1	I	183	ILE	3.0
1	A	128	GLU	3.0
1	G	190	SER	3.0
1	A	132	GLY	2.9
1	G	125	LYS	2.9
1	I	211	ASN	2.9
1	F	76	ARG	2.9
1	B	106	ASP	2.9
1	F	128	GLU	2.9
1	D	101	ALA	2.9
1	B	152	PRO	2.9
1	F	151	THR	2.8
2	L	1	DT	2.8
1	A	151	THR	2.8
1	C	141	ARG	2.8
1	F	155	GLY	2.8
1	B	182	ARG	2.8
1	H	194	LEU	2.8
1	F	133	ASP	2.8
1	H	147	ALA	2.8
1	B	130	ASP	2.8
1	F	217	VAL	2.8
1	I	123	VAL	2.8
1	B	185	ASP	2.7
1	I	131	VAL	2.7
1	B	151	THR	2.7
1	B	77	ASP	2.7
1	I	125	LYS	2.7
1	A	131	VAL	2.7
1	F	188	ARG	2.7
1	H	125	LYS	2.7
1	F	125	LYS	2.7
1	D	176	VAL	2.7
1	I	188	ARG	2.7
2	X	9	DT	2.6
1	I	76	ARG	2.6
1	G	187	VAL	2.6
1	G	148	LEU	2.6
1	D	56	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	57	ASN	2.6
1	F	185	ASP	2.6
1	H	217	VAL	2.6
1	G	116	GLU	2.6
1	G	131	VAL	2.6
1	G	137	ARG	2.6
1	F	78	SER	2.5
1	B	178	VAL	2.5
1	G	173	ARG	2.5
1	A	99	GLU	2.5
1	I	127	GLY	2.5
1	C	157	ALA	2.5
1	I	122	THR	2.5
1	I	215	SER	2.5
1	G	121	PHE	2.4
1	F	6	THR	2.4
1	I	139	MET	2.4
1	G	115	LEU	2.4
1	I	153	LEU	2.4
1	I	-1	ALA	2.4
1	F	61	ALA	2.4
1	I	178	VAL	2.4
1	I	207	VAL	2.4
1	A	153	LEU	2.4
1	D	179	GLU	2.3
1	A	137	ARG	2.3
1	C	109	ARG	2.3
1	F	175	ARG	2.3
1	I	136	TYR	2.3
1	H	183	ILE	2.3
1	H	177	GLN	2.3
1	D	189	GLU	2.3
1	F	148	LEU	2.3
1	A	30	ILE	2.3
1	D	115	LEU	2.3
1	F	109	ARG	2.2
1	I	187	VAL	2.2
1	B	177	GLN	2.2
1	C	175	ARG	2.2
1	H	57	ASN	2.2
1	F	114	LEU	2.2
1	G	145	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	77	ASP	2.2
1	G	174	THR	2.2
1	A	186	PHE	2.1
1	A	181	ASP	2.1
1	H	191	VAL	2.1
1	D	102	PRO	2.1
1	A	184	ASP	2.1
1	I	144	ASP	2.1
1	G	113	ARG	2.1
2	X	10	DT	2.1
1	H	184	ASP	2.1
1	D	157	ALA	2.1
1	I	142	GLY	2.0
1	H	137	ARG	2.0
1	A	189	GLU	2.0
1	B	98	GLU	2.0
1	I	154	ASN	2.0
1	D	107	PHE	2.0
1	I	179	GLU	2.0
1	G	134	VAL	2.0
1	F	182	ARG	2.0
1	F	150	VAL	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(\AA^2)	Q<0.9
3	PO4	B	301	5/5	0.91	0.24	-0.25	70,84,94,106	0

6.5 Other polymers [i](#)

There are no such residues in this entry.