



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:36 AM UTC

PDB ID : 4UMK / pdb_00004umk
Title : The complex of Spo0J and parS DNA in chromosomal partition system
Authors : Chen, B.W.; Chu, C.H.; Tung, J.Y.; Hsu, C.E.; Hsiao, C.D.; Sun, Y.J.
Deposited on : 2014-05-19
Resolution : 3.10 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

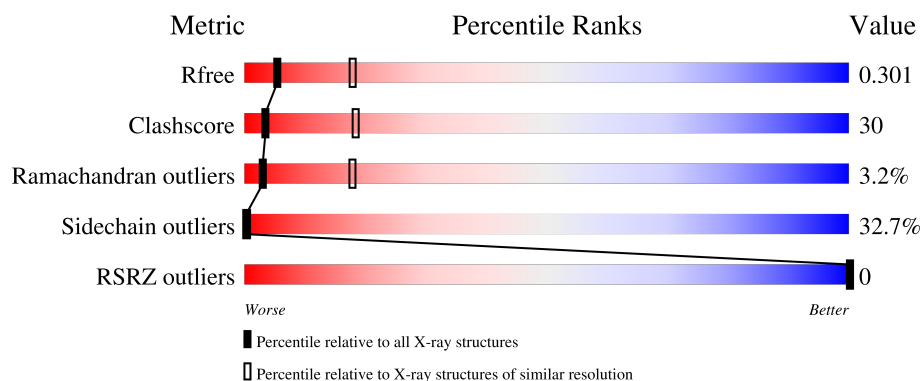
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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.



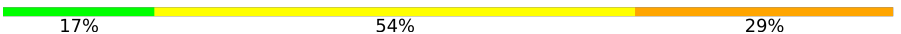
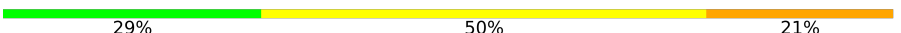
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	240	
1	B	240	
1	C	240	
1	D	240	
2	W	24	

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Mol	Chain	Length	Quality of chain
2	Z	24	 38% 46% 17%
3	X	24	 17% 54% 29%
3	Y	24	 29% 50% 21%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7994 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 PROBABLE CHROMOSOME-PARTITIONING PROTEIN PARB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	172	Total	C	N	O	S	0	0	0
			1367	858	244	261	4			
1	B	183	Total	C	N	O	S	0	0	0
			1463	920	261	278	4			
1	C	193	Total	C	N	O	S	0	0	0
			1535	964	271	295	5			
1	D	176	Total	C	N	O	S	0	0	0
			1400	880	248	268	4			

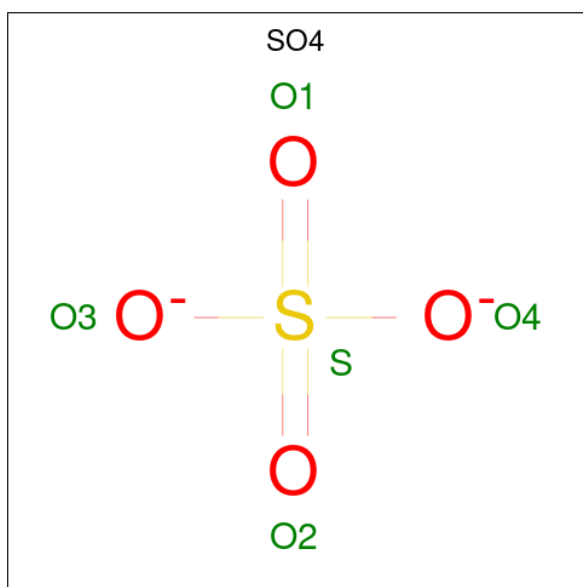
- Molecule 2 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	W	24	Total	C	N	O	P	0	0	0
			501	236	100	141	24			
2	Z	24	Total	C	N	O	P	0	0	0
			501	236	100	141	24			

- Molecule 3 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	X	24	Total	C	N	O	P	0	0	0
			483	231	81	147	24			
3	Y	24	Total	C	N	O	P	0	0	0
			483	231	81	147	24			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	W	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	53	Total	O	0	0
			53	53		
5	B	47	Total	O	0	0
			47	47		
5	C	45	Total	O	0	0
			45	45		
5	D	31	Total	O	0	0
			31	31		
5	W	16	Total	O	0	0
			16	16		
5	X	7	Total	O	0	0
			7	7		

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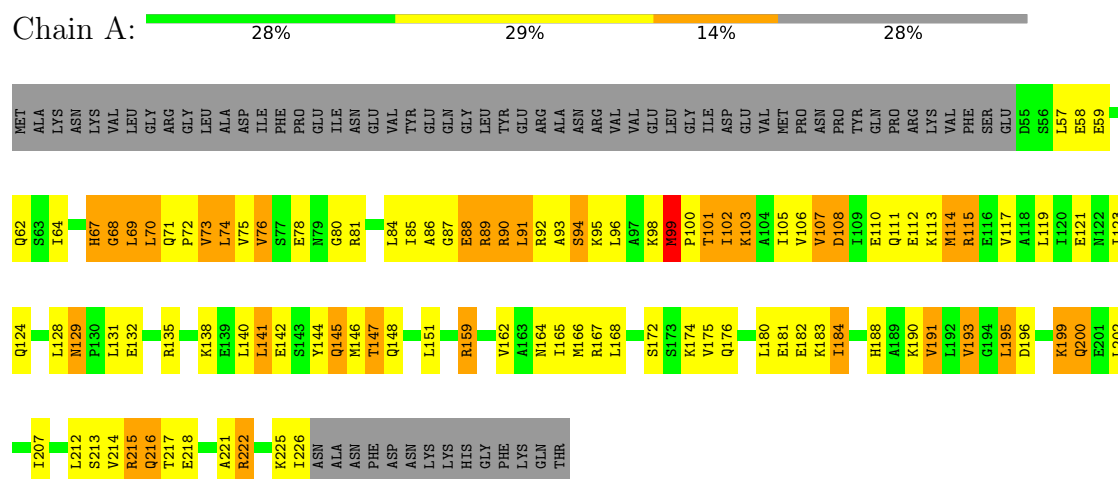
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	Y	16	Total	O	0	0
			16	16		
5	Z	16	Total	O	0	0
			16	16		

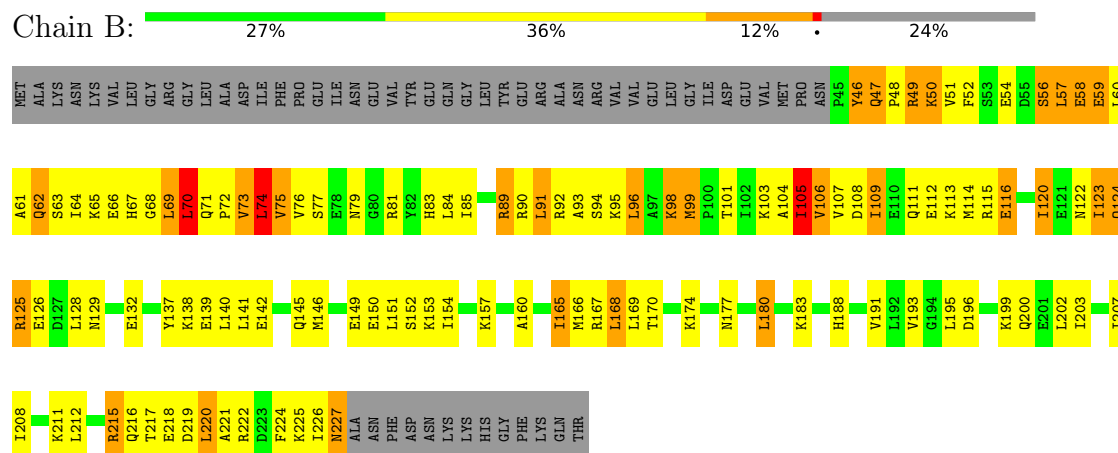
3 Residue-property plots

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

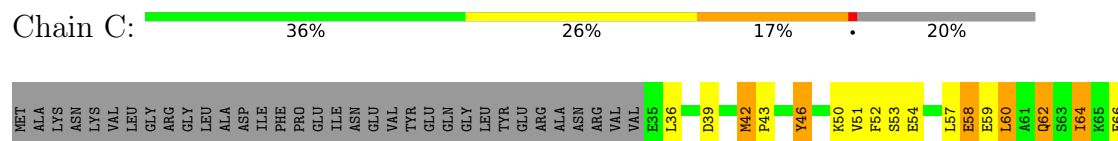
• Molecule 1: PROBABLE CHROMOSOME-PARTITIONING PROTEIN PARB

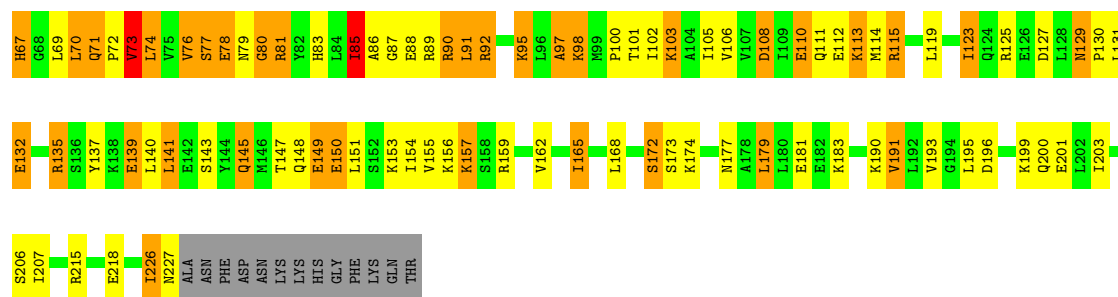


• Molecule 1: PROBABLE CHROMOSOME-PARTITIONING PROTEIN PARB



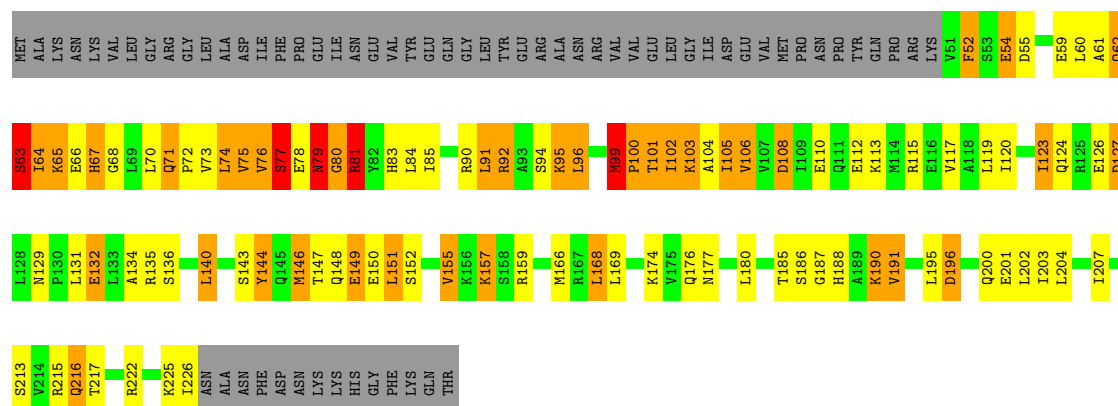
• Molecule 1: PROBABLE CHROMOSOME-PARTITIONING PROTEIN PARB





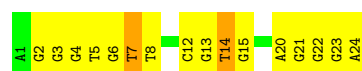
• Molecule 1: PROBABLE CHROMOSOME-PARTITIONING PROTEIN PARB

Chain D: 31% 25% 15% • 27%



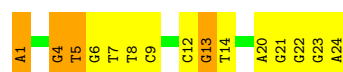
• Molecule 2: DNA

Chain W: 33% 58% 8%



• Molecule 2: DNA

Chain Z: 38% 46% 17%



• Molecule 3: DNA

Chain X: 17% 54% 29%



• Molecule 3: DNA

Chain Y: 29% 50% 21%

T1	C2	C3	C4	T5	G6	T7	T8		G12	G13	T14	G15	G16		C21	C22	C23	T24
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.48Å 232.73Å 78.34Å 90.00° 109.15° 90.00°	Depositor
Resolution (Å)	23.57 – 3.10 23.57 – 3.10	Depositor EDS
% Data completeness (in resolution range)	96.4 (23.57-3.10) 91.4 (23.57-3.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 3.10Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.251 , 0.299 0.252 , 0.301	Depositor DCC
R_{free} test set	1994 reflections (5.99%)	wwPDB-VP
Wilson B-factor (Å ²)	73.0	Xtriage
Anisotropy	0.540	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 71.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.119 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7994	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.70% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.70	1/1377 (0.1%)	1.15	5/1847 (0.3%)
1	B	0.62	0/1477	1.09	5/1982 (0.3%)
1	C	0.68	0/1550	1.13	8/2084 (0.4%)
1	D	0.70	1/1411 (0.1%)	1.05	4/1893 (0.2%)
2	W	0.51	0/564	1.37	5/870 (0.6%)
2	Z	0.59	0/564	1.44	10/870 (1.1%)
3	X	0.61	0/538	1.62	12/826 (1.5%)
3	Y	0.59	0/538	1.50	6/826 (0.7%)
All	All	0.65	2/8019 (0.0%)	1.23	55/11198 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	D	0	3
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	102	ILE	CA-CB	8.80	1.66	1.54
1	A	123	ILE	CA-CB	5.21	1.60	1.53

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	X	24	DT	O5'-C5'-C4'	-13.34	90.80	110.80
1	A	99	MET	CA-C-N	-8.78	108.87	119.84
1	A	99	MET	C-N-CA	-8.78	108.87	119.84
2	W	14	DT	N1-C1'-C2'	8.08	125.61	113.50
1	B	75	VAL	N-CA-C	7.70	119.33	108.93
3	Y	14	DT	C2'-C3'-O3'	7.64	122.97	111.50
1	B	105	ILE	N-CA-C	7.51	118.69	108.17
2	Z	13	DG	O4'-C1'-N9	-7.48	97.19	108.40
1	C	106	VAL	N-CA-C	7.44	119.04	109.30
1	B	74	LEU	N-CA-C	7.13	120.28	109.23
2	Z	7	DT	P-O5'-C5'	-7.12	109.33	120.00
3	Y	15	DG	C4'-C3'-O3'	-7.02	99.47	110.00
2	W	8	DT	P-O5'-C5'	-7.01	109.49	120.00
2	W	14	DT	O4'-C1'-N1	-6.95	97.97	108.40
3	Y	5	DT	C5'-C4'-C3'	-6.93	104.50	114.90
3	X	8	DT	C4'-C3'-O3'	-6.86	99.71	110.00
2	Z	1	DA	C5-C6-N6	6.80	133.60	123.40
3	X	24	DT	C5'-C4'-C3'	6.60	124.81	114.90
3	Y	4	DC	O4'-C1'-C2'	-6.50	96.65	106.40
3	X	9	DT	C4'-C3'-O3'	-6.45	100.32	110.00
1	C	42	MET	CA-C-N	-6.29	111.98	119.84
1	C	42	MET	C-N-CA	-6.29	111.98	119.84
1	A	129	ASN	CA-C-N	-6.25	113.37	119.87
1	A	129	ASN	C-N-CA	-6.25	113.37	119.87
3	X	24	DT	C4'-C3'-O3'	-6.20	100.71	110.00
1	C	73	VAL	N-CA-C	6.16	116.74	107.75
3	X	1	DT	N1-C1'-C2'	6.05	122.58	113.50
2	Z	1	DA	N1-C6-N6	-6.02	109.97	119.00
3	X	13	DG	O4'-C1'-N9	-6.02	99.37	108.40
2	Z	7	DT	O4'-C1'-N1	-5.96	99.46	108.40
1	A	181	GLU	N-CA-C	-5.89	103.16	110.41
2	W	14	DT	C5'-C4'-O4'	-5.81	100.68	109.40
1	D	71	GLN	CA-C-N	5.75	126.07	119.92
1	D	71	GLN	C-N-CA	5.75	126.07	119.92
3	X	15	DG	C4'-C3'-O3'	-5.64	101.54	110.00
1	C	108	ASP	N-CA-C	-5.64	107.05	114.04
2	W	7	DT	C4-C5-C7	-5.61	113.99	122.40
1	D	99	MET	CA-C-N	5.58	125.25	119.56
1	D	99	MET	C-N-CA	5.58	125.25	119.56
1	C	73	VAL	CB-CA-C	-5.57	102.86	110.77
1	C	154	ILE	CB-CA-C	-5.53	104.90	111.97
2	Z	7	DT	N1-C1'-C2'	5.40	121.59	113.50
1	B	47	GLN	N-CA-C	-5.32	102.08	108.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	X	23	DC	N1-C1'-C2'	5.31	121.47	113.50
3	X	11	DA	C4'-C3'-O3'	-5.27	102.09	110.00
3	X	13	DG	O5'-C5'-C4'	-5.18	103.03	110.80
3	Y	5	DT	P-O5'-C5'	-5.17	112.24	120.00
1	C	174	LYS	N-CA-C	-5.13	105.58	111.07
2	Z	4	DG	C5'-C4'-O4'	5.11	117.06	109.40
2	Z	13	DG	O5'-C5'-C4'	-5.09	103.16	110.80
3	X	24	DT	O4'-C1'-N1	-5.09	100.76	108.40
3	Y	13	DG	C2'-C3'-O3'	5.09	119.13	111.50
2	Z	1	DA	P-O5'-C5'	5.07	127.61	120.00
1	B	222	ARG	N-CA-C	-5.07	105.84	111.36
2	Z	5	DT	C5'-C4'-C3'	-5.04	107.34	114.90

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	67	HIS	Peptide
1	B	74	LEU	Peptide
1	D	101	THR	Peptide
1	D	77	SER	Peptide
1	D	81	ARG	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1367	0	1447	126	0
1	B	1463	0	1540	109	0
1	C	1535	0	1600	75	0
1	D	1400	0	1476	83	0
2	W	501	0	269	30	0
2	Z	501	0	269	23	0
3	X	483	0	272	21	0
3	Y	483	0	272	27	0
4	A	10	0	0	0	0
4	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	5	0	0	0	0
4	D	5	0	0	0	0
4	W	5	0	0	0	0
5	A	53	0	0	5	0
5	B	47	0	0	9	0
5	C	45	0	0	7	0
5	D	31	0	0	4	0
5	W	16	0	0	3	0
5	X	7	0	0	1	0
5	Y	16	0	0	1	0
5	Z	16	0	0	2	0
All	All	7994	0	7145	441	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 30.

All (441) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Y:24:DT:H4'	3:Y:24:DT:OP1	1.49	1.10
2:Z:4:DG:H2''	2:Z:5:DT:H5''	1.36	1.07
2:W:3:DG:H1	3:X:22:DC:H42	1.02	1.00
1:A:67:HIS:HD1	1:A:70:LEU:HD12	1.29	0.95
2:W:3:DG:H1	3:X:22:DC:N4	1.68	0.91
1:A:102:ILE:HD12	1:A:106:VAL:HG11	1.53	0.90
1:D:123:ILE:HD11	1:D:155:VAL:HG12	1.55	0.88
1:B:74:LEU:H	1:B:90:ARG:HH21	1.24	0.86
1:B:62:GLN:HA	1:B:65:LYS:HG2	1.59	0.85
1:A:75:VAL:HG12	1:A:84:LEU:HA	1.59	0.84
1:A:67:HIS:HB2	1:A:70:LEU:HA	1.60	0.84
1:C:42:MET:HG2	1:C:145:GLN:HB2	1.60	0.83
2:W:20:DA:H61	3:X:5:DT:H3	1.27	0.82
1:C:60:LEU:HG	1:C:62:GLN:HB3	1.63	0.79
1:A:89:ARG:HH21	1:B:89:ARG:HE	1.30	0.78
1:C:119:LEU:HD21	1:C:139:GLU:HB3	1.64	0.77
1:A:148:GLN:NE2	2:W:5:DT:H3'	1.99	0.77
1:B:51:VAL:HG12	1:B:91:LEU:HD11	1.66	0.76
2:W:21:DG:H1	3:X:4:DC:H42	1.34	0.75
1:D:140:LEU:HA	1:D:144:TYR:CE1	2.22	0.74
1:B:74:LEU:O	1:B:90:ARG:NE	2.20	0.74
1:A:159:ARG:HH21	2:W:5:DT:H2'	1.53	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:LEU:H	1:A:69:LEU:HD22	1.53	0.73
1:B:79:ASN:HB2	1:B:111:GLN:HB2	1.69	0.73
1:C:42:MET:O	1:C:145:GLN:NE2	2.22	0.73
2:W:20:DA:N6	3:X:5:DT:H3	1.86	0.73
1:D:103:LYS:HD2	1:D:105:ILE:HG23	1.71	0.73
1:A:74:LEU:HD22	1:A:85:ILE:HD12	1.70	0.72
1:D:78:GLU:O	1:D:80:GLY:N	2.22	0.72
2:W:23:DG:N7	5:W:2014:HOH:O	2.22	0.72
1:D:65:LYS:NZ	1:D:65:LYS:O	2.23	0.72
1:D:103:LYS:HE2	1:D:106:VAL:HG13	1.71	0.71
1:A:78:GLU:HB3	1:A:81:ARG:H	1.54	0.70
1:D:78:GLU:N	1:D:78:GLU:OE1	2.24	0.70
3:Y:3:DC:H42	2:Z:22:DG:H1	1.38	0.70
1:C:149:GLU:OE1	1:C:150:GLU:N	2.25	0.70
1:C:53:SER:OG	1:C:90:ARG:NH2	2.23	0.69
1:B:85:ILE:HD11	1:B:90:ARG:HB3	1.74	0.69
1:B:74:LEU:HD13	1:B:90:ARG:CZ	2.22	0.69
1:C:70:LEU:HD22	1:C:71:GLN:H	1.58	0.69
1:D:140:LEU:HA	1:D:144:TYR:HE1	1.55	0.69
1:D:78:GLU:HG2	1:D:101:THR:HB	1.74	0.69
1:D:75:VAL:HG11	1:D:95:LYS:HE2	1.75	0.68
1:D:196:ASP:OD1	1:D:196:ASP:N	2.25	0.68
5:Y:2004:HOH:O	2:Z:23:DG:N2	2.26	0.68
1:A:103:LYS:NZ	5:A:2021:HOH:O	2.27	0.68
1:A:164:ASN:HA	1:A:167:ARG:HD3	1.75	0.67
2:Z:22:DG:H2'	2:Z:23:DG:C8	2.29	0.67
1:A:100:PRO:HD2	1:A:101:THR:H	1.60	0.66
1:C:73:VAL:HG22	1:C:108:ASP:HB3	1.76	0.66
1:A:191:VAL:HG11	1:A:217:THR:HG22	1.77	0.66
1:A:62:GLN:HE22	1:A:90:ARG:HA	1.60	0.66
1:C:112:GLU:OE1	1:C:115:ARG:NH1	2.29	0.65
1:A:114:MET:HG3	1:C:89:ARG:HH21	1.62	0.65
2:W:21:DG:H1	3:X:4:DC:N4	1.93	0.65
1:A:67:HIS:CD2	1:A:68:GLY:O	2.48	0.65
1:B:124:GLN:NE2	5:B:2023:HOH:O	2.29	0.65
3:Y:14:DT:H2''	3:Y:15:DG:O5'	1.97	0.65
1:A:90:ARG:HG2	1:A:90:ARG:HH11	1.62	0.65
1:A:214:VAL:O	1:A:218:GLU:HG3	1.97	0.65
2:W:3:DG:N2	3:X:22:DC:N3	2.42	0.64
1:B:195:LEU:O	1:B:200:GLN:NE2	2.24	0.64
1:B:177:ASN:HA	1:B:180:LEU:HD22	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:ASN:OD1	1:B:227:ASN:N	2.31	0.64
2:W:24:DA:C2	3:X:1:DT:N3	2.61	0.64
1:C:137:TYR:CE2	1:C:162:VAL:HG22	2.33	0.64
3:X:22:DC:H2'	3:X:23:DC:C6	2.32	0.64
1:A:212:LEU:HB3	1:A:216:GLN:HB2	1.81	0.63
1:C:85:ILE:HG12	1:C:86:ALA:H	1.64	0.62
2:Z:4:DG:H2''	2:Z:5:DT:C5'	2.23	0.62
1:B:46:TYR:CZ	1:B:51:VAL:HG13	2.35	0.62
1:D:159:ARG:NH2	2:Z:6:DG:O6	2.29	0.62
1:B:126:GLU:OE1	1:B:126:GLU:N	2.33	0.62
1:C:42:MET:HB3	1:C:43:PRO:HD3	1.81	0.62
3:Y:4:DC:H2''	3:Y:5:DT:H5''	1.81	0.62
3:Y:12:DC:H2''	3:Y:13:DG:C8	2.35	0.62
3:X:8:DT:H1'	3:X:9:DT:H5'	1.82	0.61
1:B:74:LEU:C	1:B:90:ARG:HE	2.08	0.61
1:C:125:ARG:HH21	1:C:127:ASP:CG	2.09	0.61
1:A:213:SER:OG	1:A:216:GLN:HG3	2.01	0.61
1:C:78:GLU:CB	1:C:87:GLY:H	2.13	0.61
1:D:222:ARG:HA	1:D:225:LYS:HD3	1.83	0.61
1:A:88:GLU:OE1	1:A:92:ARG:NH1	2.33	0.60
1:C:72:PRO:HB2	1:C:108:ASP:HB2	1.83	0.60
1:D:99:MET:SD	1:D:100:PRO:HD2	2.40	0.60
1:D:144:TYR:HB2	1:D:146:MET:HG2	1.83	0.60
3:Y:1:DT:H2'	3:Y:2:DC:C6	2.35	0.60
1:A:103:LYS:HD2	5:A:2022:HOH:O	2.00	0.60
1:A:73:VAL:HG12	1:A:74:LEU:H	1.65	0.60
1:C:129:ASN:ND2	1:C:132:GLU:OE1	2.35	0.60
1:B:69:LEU:HD12	1:B:69:LEU:O	2.00	0.60
1:A:100:PRO:CD	1:A:101:THR:H	2.15	0.59
1:C:165:ILE:O	1:C:168:LEU:HB2	2.02	0.59
1:B:154:ILE:HG21	1:D:91:LEU:HD12	1.84	0.59
1:A:87:GLY:HA2	1:A:90:ARG:HD3	1.85	0.59
2:W:12:DC:H2''	2:W:13:DG:C8	2.37	0.59
1:B:62:GLN:HG2	1:B:96:LEU:HB2	1.85	0.58
2:W:22:DG:H1'	2:W:23:DG:C8	2.37	0.58
1:B:141:LEU:HD21	1:B:151:LEU:HD22	1.84	0.58
1:A:222:ARG:HB2	1:A:222:ARG:HH11	1.68	0.58
1:D:62:GLN:O	1:D:64:ILE:N	2.37	0.58
1:A:207:ILE:HA	1:A:212:LEU:HD12	1.85	0.58
1:A:215:ARG:NH1	1:A:218:GLU:OE1	2.37	0.58
1:A:86:ALA:H	1:C:83:HIS:CE1	2.21	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:PRO:HB2	1:A:103:LYS:NZ	2.19	0.57
1:C:54:GLU:OE1	1:C:54:GLU:N	2.36	0.57
1:C:203:ILE:O	1:C:207:ILE:HG13	2.03	0.57
1:C:76:VAL:HG12	1:C:77:SER:H	1.69	0.57
1:D:204:LEU:HA	1:D:207:ILE:HD12	1.85	0.57
2:W:14:DT:H2''	2:W:15:DG:O4'	2.05	0.57
1:B:66:GLU:OE2	1:B:69:LEU:HD22	2.04	0.57
1:B:66:GLU:HA	1:B:69:LEU:HB2	1.85	0.57
1:C:91:LEU:HG	1:C:92:ARG:N	2.18	0.57
1:D:67:HIS:HB3	5:D:2004:HOH:O	2.05	0.56
1:A:69:LEU:HB2	1:A:72:PRO:HA	1.86	0.56
1:A:144:TYR:CD1	1:C:89:ARG:HB2	2.41	0.56
3:Y:5:DT:H2''	3:Y:6:DG:H8	1.70	0.56
3:Y:24:DT:OP1	3:Y:24:DT:C4'	2.37	0.56
1:C:129:ASN:OD1	1:C:129:ASN:N	2.30	0.56
1:D:92:ARG:O	1:D:96:LEU:HG	2.05	0.56
1:B:74:LEU:HD13	1:B:90:ARG:NE	2.21	0.56
1:B:125:ARG:HA	1:B:125:ARG:HH11	1.71	0.56
1:B:165:ILE:HD13	1:B:168:LEU:HD21	1.87	0.56
2:W:2:DG:H8	5:W:2001:HOH:O	1.88	0.56
1:A:144:TYR:CG	1:C:89:ARG:HD2	2.40	0.55
3:Y:3:DC:N4	2:Z:22:DG:H1	2.03	0.55
1:A:99:MET:HE3	1:A:100:PRO:HD3	1.89	0.55
1:A:69:LEU:N	1:A:69:LEU:HD13	2.22	0.55
1:A:141:LEU:HD21	1:A:151:LEU:HD22	1.88	0.55
1:C:110:GLU:HG3	1:C:113:LYS:HZ1	1.72	0.55
1:A:84:LEU:HD11	1:A:87:GLY:HA3	1.88	0.55
1:C:123:ILE:HD12	1:C:140:LEU:HD12	1.88	0.55
1:A:89:ARG:HH21	1:B:89:ARG:NE	2.03	0.55
1:A:172:SER:OG	1:A:174:LYS:NZ	2.40	0.55
1:D:152:SER:HB3	1:D:157:LYS:O	2.07	0.55
1:C:191:VAL:HG11	1:C:218:GLU:HA	1.88	0.55
1:D:131:LEU:HD12	1:D:180:LEU:HD21	1.88	0.55
1:B:46:TYR:CE1	1:B:51:VAL:HG13	2.42	0.54
3:Y:22:DC:H2'	3:Y:23:DC:C6	2.41	0.54
1:A:90:ARG:HH11	1:A:90:ARG:CG	2.20	0.54
3:Y:5:DT:H2''	3:Y:6:DG:H5'	1.89	0.54
2:Z:4:DG:C2'	2:Z:5:DT:H5''	2.25	0.54
1:A:148:GLN:OE1	1:A:159:ARG:NH1	2.40	0.54
1:D:185:THR:OG1	1:D:188:HIS:ND1	2.32	0.54
1:A:67:HIS:HB2	1:A:70:LEU:CA	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:117:VAL:HA	1:D:120:ILE:HG22	1.90	0.54
3:Y:14:DT:H2'	3:Y:15:DG:C8	2.42	0.54
1:A:165:ILE:O	1:A:168:LEU:HB2	2.07	0.54
1:C:130:PRO:HG2	1:C:179:LEU:HD11	1.90	0.54
1:A:117:VAL:HG11	5:A:2034:HOH:O	2.07	0.54
1:C:98:LYS:HE3	1:C:101:THR:HA	1.88	0.54
1:B:50:LYS:HG2	1:D:110:GLU:HG2	1.89	0.53
1:B:63:SER:N	1:B:99:MET:HE1	2.24	0.53
1:B:115:ARG:HH12	1:D:75:VAL:HG22	1.72	0.53
1:A:111:GLN:O	1:A:115:ARG:HG2	2.09	0.53
2:W:24:DA:N1	3:X:1:DT:C4	2.76	0.53
1:B:89:ARG:CZ	1:B:89:ARG:HB2	2.35	0.53
1:B:202:LEU:HD12	5:B:2039:HOH:O	2.07	0.53
1:B:124:GLN:CD	1:B:124:GLN:H	2.17	0.53
1:C:181:GLU:HB3	1:C:183:LYS:HG2	1.90	0.53
1:B:74:LEU:HG	1:B:92:ARG:HH22	1.74	0.53
1:C:155:VAL:HG13	1:C:157:LYS:HG3	1.91	0.52
1:D:191:VAL:HG21	1:D:217:THR:HG22	1.90	0.52
1:B:84:LEU:HD23	1:B:84:LEU:H	1.75	0.52
1:A:195:LEU:O	1:A:200:GLN:NE2	2.42	0.52
2:W:22:DG:OP2	2:W:22:DG:H2'	2.08	0.52
1:A:59:GLU:HG2	1:A:93:ALA:HA	1.91	0.52
1:B:92:ARG:HD2	1:B:93:ALA:N	2.25	0.52
1:C:85:ILE:HG12	1:C:86:ALA:N	2.24	0.52
1:A:191:VAL:HG21	1:A:218:GLU:HA	1.91	0.52
1:A:74:LEU:HA	1:A:107:VAL:O	2.10	0.52
1:A:172:SER:HB3	1:A:175:VAL:HG23	1.91	0.52
1:D:91:LEU:HD13	1:D:92:ARG:H	1.75	0.52
1:D:68:GLY:HA2	1:D:104:ALA:C	2.35	0.51
1:D:65:LYS:HD3	1:D:67:HIS:HB2	1.91	0.51
3:X:17:DA:H1'	3:X:18:DA:H5'	1.92	0.51
1:A:188:HIS:O	1:A:191:VAL:HG12	2.11	0.51
1:B:47:GLN:HG3	1:B:48:PRO:HD2	1.92	0.51
3:Y:4:DC:C2'	3:Y:5:DT:H5''	2.41	0.51
1:B:116:GLU:O	1:B:120:ILE:HG23	2.10	0.51
1:B:123:ILE:HD12	1:D:92:ARG:NH2	2.25	0.51
1:C:57:LEU:HG	1:C:58:GLU:H	1.75	0.51
1:C:149:GLU:O	1:C:153:LYS:HG2	2.10	0.51
1:A:90:ARG:O	1:A:94:SER:HB2	2.10	0.51
1:A:102:ILE:HG22	1:A:103:LYS:H	1.75	0.51
1:C:73:VAL:O	1:C:74:LEU:HD22	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:190:LYS:HE2	2:Z:9:DC:H41	1.74	0.51
2:Z:6:DG:H5'	2:Z:6:DG:C8	2.46	0.51
1:D:65:LYS:HG3	1:D:67:HIS:H	1.76	0.50
2:W:3:DG:H2''	2:W:4:DG:H8	1.75	0.50
1:C:132:GLU:OE2	1:C:135:ARG:NH2	2.45	0.50
1:D:155:VAL:HG23	1:D:157:LYS:HB2	1.91	0.50
3:Y:15:DG:H2''	3:Y:16:DG:H5''	1.93	0.50
1:A:73:VAL:N	5:A:2004:HOH:O	2.41	0.50
1:B:74:LEU:N	1:B:90:ARG:HH21	2.02	0.50
1:D:144:TYR:N	1:D:144:TYR:CD1	2.79	0.50
1:A:76:VAL:HG11	1:C:86:ALA:HB3	1.94	0.50
1:C:98:LYS:NZ	1:C:101:THR:HB	2.27	0.50
1:D:136:SER:O	1:D:140:LEU:HB2	2.12	0.50
1:A:167:ARG:NH2	2:W:7:DT:H2'	2.26	0.50
1:A:159:ARG:NH2	2:W:5:DT:H2'	2.23	0.50
1:C:145:GLN:HA	1:C:145:GLN:OE1	2.12	0.50
1:A:184:ILE:HD11	1:A:207:ILE:HG21	1.94	0.50
3:Y:24:DT:O4	2:Z:1:DA:N1	2.44	0.50
1:A:102:ILE:HG22	1:A:103:LYS:N	2.27	0.49
1:B:50:LYS:HE2	1:B:51:VAL:HG22	1.95	0.49
1:B:92:ARG:HH11	1:B:93:ALA:HB2	1.76	0.49
1:B:104:ALA:HB1	5:B:2017:HOH:O	2.13	0.49
2:Z:21:DG:H2'	2:Z:22:DG:C8	2.47	0.49
1:A:114:MET:HB3	1:A:115:ARG:HH21	1.77	0.49
1:B:92:ARG:HD2	1:B:93:ALA:H	1.77	0.49
3:Y:3:DC:N3	2:Z:22:DG:N2	2.53	0.49
1:A:129:ASN:ND2	5:A:2030:HOH:O	2.46	0.49
1:B:199:LYS:HD3	1:B:224:PHE:HB3	1.94	0.49
1:C:76:VAL:HG12	1:C:77:SER:N	2.27	0.49
3:Y:14:DT:H2''	3:Y:15:DG:O4'	2.12	0.49
1:A:90:ARG:HD2	1:B:89:ARG:NH2	2.27	0.49
1:A:99:MET:HB3	1:A:100:PRO:HD3	1.95	0.49
1:D:174:LYS:HD2	1:D:201:GLU:HG3	1.94	0.49
3:X:14:DT:H2''	3:X:15:DG:O5'	2.13	0.49
1:B:167:ARG:CZ	1:B:193:VAL:HG21	2.43	0.49
1:D:108:ASP:HB3	5:D:2011:HOH:O	2.13	0.49
2:Z:8:DT:H2''	2:Z:9:DC:H5'	1.93	0.49
1:B:49:ARG:NH1	5:B:2004:HOH:O	2.46	0.49
1:A:114:MET:HE3	1:A:114:MET:HA	1.95	0.48
1:C:36:LEU:HD13	1:C:79:ASN:HB2	1.95	0.48
1:C:98:LYS:HZ2	1:C:101:THR:HB	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:140:LEU:HD13	1:D:144:TYR:CE1	2.48	0.48
1:A:69:LEU:HD22	1:A:69:LEU:N	2.26	0.48
1:A:90:ARG:HG3	1:B:89:ARG:CZ	2.43	0.48
1:B:83:HIS:O	1:B:83:HIS:ND1	2.46	0.48
1:A:96:LEU:H	1:A:96:LEU:HD22	1.77	0.48
1:A:115:ARG:NH1	1:A:115:ARG:HA	2.29	0.48
1:A:89:ARG:NH2	1:B:92:ARG:HE	2.11	0.48
5:W:2015:HOH:O	3:X:2:DC:N4	2.40	0.48
1:B:90:ARG:HB2	5:B:2012:HOH:O	2.13	0.48
1:C:92:ARG:NH2	5:C:2021:HOH:O	2.39	0.48
1:C:111:GLN:O	1:C:115:ARG:HB2	2.13	0.48
1:D:71:GLN:HG2	1:D:72:PRO:HD2	1.95	0.48
1:B:74:LEU:HG	1:B:92:ARG:NH2	2.30	0.47
2:Z:6:DG:H5'	2:Z:6:DG:H8	1.79	0.47
1:D:61:ALA:C	1:D:63:SER:H	2.21	0.47
1:B:73:VAL:HG11	1:B:106:VAL:HB	1.97	0.47
1:A:74:LEU:HD12	1:B:72:PRO:HB3	1.95	0.47
1:B:141:LEU:O	1:B:145:GLN:HA	2.14	0.47
1:B:191:VAL:HG11	1:B:217:THR:HG22	1.97	0.47
1:D:52:PHE:N	1:D:52:PHE:CD1	2.80	0.47
1:C:64:ILE:N	5:C:2008:HOH:O	2.47	0.47
1:A:146:MET:HG3	1:A:147:THR:N	2.30	0.47
1:B:56:SER:O	1:B:60:LEU:HB3	2.15	0.47
1:D:77:SER:O	1:D:77:SER:OG	2.28	0.47
1:D:94:SER:C	1:D:95:LYS:HE3	2.39	0.47
1:D:119:LEU:O	1:D:123:ILE:HG22	2.14	0.47
1:B:203:ILE:O	1:B:207:ILE:HG13	2.14	0.47
1:D:127:ASP:OD1	1:D:129:ASN:N	2.48	0.47
1:A:58:GLU:HG3	1:B:52:PHE:CD2	2.50	0.47
1:A:73:VAL:HG22	1:A:90:ARG:HG2	1.96	0.47
1:A:215:ARG:HD3	1:A:215:ARG:HA	1.43	0.47
1:B:218:GLU:OE2	5:B:2037:HOH:O	2.21	0.47
2:Z:20:DA:H2'	5:Z:2012:HOH:O	2.15	0.47
1:B:146:MET:HE2	1:B:151:LEU:N	2.29	0.46
2:Z:23:DG:H2''	5:Z:2013:HOH:O	2.15	0.46
1:A:71:GLN:CD	1:B:74:LEU:HB2	2.40	0.46
1:A:162:VAL:O	1:A:166:MET:HG3	2.15	0.46
1:C:113:LYS:HZ3	1:C:114:MET:HB2	1.80	0.46
1:A:99:MET:HE3	1:A:99:MET:HB3	1.69	0.46
1:A:164:ASN:OD1	2:W:7:DT:H72	2.15	0.46
1:D:101:THR:O	1:D:103:LYS:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:146:MET:HE3	1:D:146:MET:HB2	1.68	0.46
1:A:110:GLU:OE1	1:A:112:GLU:N	2.48	0.46
1:B:149:GLU:O	1:B:153:LYS:HG2	2.16	0.46
1:B:166:MET:O	1:B:169:LEU:HB2	2.15	0.46
1:B:188:HIS:O	1:B:191:VAL:HG12	2.16	0.46
1:C:103:LYS:HE2	1:C:103:LYS:HB3	1.52	0.46
1:A:117:VAL:O	1:A:121:GLU:HG2	2.15	0.46
1:D:195:LEU:O	1:D:200:GLN:NE2	2.42	0.46
1:D:187:GLY:O	1:D:191:VAL:HG13	2.15	0.46
1:A:71:GLN:OE1	1:B:74:LEU:HB2	2.15	0.46
1:B:212:LEU:HD13	1:B:216:GLN:HB3	1.98	0.46
1:D:101:THR:HG23	5:D:2010:HOH:O	2.15	0.46
1:A:148:GLN:HE21	2:W:5:DT:H3'	1.78	0.46
1:C:196:ASP:O	1:C:200:GLN:HG3	2.15	0.46
1:C:226:ILE:HG12	1:C:227:ASN:N	2.31	0.46
3:X:10:DC:H1'	3:X:11:DA:N7	2.31	0.46
1:A:102:ILE:CD1	1:A:106:VAL:HG11	2.35	0.45
1:B:215:ARG:HA	1:B:215:ARG:HD3	1.75	0.45
1:D:174:LYS:O	1:D:177:ASN:HB3	2.16	0.45
1:A:67:HIS:ND1	1:A:70:LEU:HD12	2.13	0.45
1:A:73:VAL:HG11	1:A:84:LEU:HD11	1.97	0.45
1:A:141:LEU:C	1:A:144:TYR:O	2.59	0.45
1:B:62:GLN:O	1:B:66:GLU:HG2	2.17	0.45
1:A:131:LEU:HD23	1:A:131:LEU:HA	1.77	0.45
1:B:167:ARG:NH1	1:B:193:VAL:HG21	2.31	0.45
1:B:216:GLN:O	1:B:219:ASP:N	2.50	0.45
1:D:132:GLU:OE1	1:D:135:ARG:NH2	2.48	0.45
2:Z:13:DG:H2''	2:Z:14:DT:OP2	2.16	0.45
1:A:69:LEU:HD12	1:A:90:ARG:HB3	1.99	0.45
1:A:69:LEU:H	1:A:69:LEU:CD2	2.21	0.45
1:A:74:LEU:HD23	1:A:108:ASP:HA	1.98	0.45
1:A:91:LEU:O	1:A:94:SER:N	2.50	0.45
1:D:157:LYS:NZ	3:Y:15:DG:OP2	2.49	0.45
2:W:2:DG:H2''	2:W:3:DG:C8	2.52	0.45
1:A:175:VAL:HG21	1:A:200:GLN:HG2	1.98	0.45
1:A:72:PRO:HB3	1:A:106:VAL:HB	1.99	0.45
1:B:95:LYS:O	1:B:98:LYS:HB3	2.16	0.45
1:B:129:ASN:HB3	1:B:132:GLU:HB2	1.98	0.45
1:C:42:MET:HE2	5:C:2036:HOH:O	2.17	0.45
1:D:103:LYS:O	1:D:106:VAL:N	2.50	0.45
1:D:149:GLU:H	1:D:149:GLU:HG2	1.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:X:13:DG:H2''	3:X:14:DT:O5'	2.16	0.45
1:C:165:ILE:HD12	1:C:165:ILE:HA	1.67	0.44
2:W:3:DG:H2''	2:W:4:DG:C8	2.52	0.44
1:D:101:THR:C	1:D:103:LYS:N	2.75	0.44
1:A:75:VAL:HG23	1:A:108:ASP:HB2	2.00	0.44
1:C:71:GLN:OE1	1:C:74:LEU:HD21	2.18	0.44
1:C:132:GLU:HG3	1:C:135:ARG:NH2	2.32	0.44
1:D:79:ASN:HD22	1:D:79:ASN:HA	1.48	0.44
3:Y:16:DG:H2'	3:Y:16:DG:H5'	1.71	0.44
1:A:89:ARG:H	1:A:89:ARG:HG3	1.24	0.44
1:A:114:MET:HB3	1:A:115:ARG:NH2	2.33	0.44
1:B:165:ILE:O	1:B:168:LEU:HG	2.17	0.44
1:D:225:LYS:HD2	5:D:2024:HOH:O	2.17	0.44
1:B:74:LEU:H	1:B:90:ARG:NH2	2.04	0.44
1:A:99:MET:CE	1:A:100:PRO:HD3	2.48	0.44
1:B:174:LYS:HD2	1:B:174:LYS:HA	1.84	0.44
1:D:61:ALA:C	1:D:63:SER:N	2.75	0.44
3:Y:23:DC:H2''	3:Y:24:DT:C6	2.53	0.44
1:B:58:GLU:HG2	1:B:62:GLN:HE22	1.81	0.44
1:D:203:ILE:HG22	1:D:207:ILE:HD11	1.99	0.44
3:Y:1:DT:H3	2:Z:24:DA:N6	2.16	0.44
1:A:144:TYR:CE2	1:A:146:MET:HB2	2.53	0.43
1:B:92:ARG:H	1:B:92:ARG:HG3	1.53	0.43
1:B:169:LEU:HD23	1:B:169:LEU:HA	1.78	0.43
1:B:105:ILE:HG12	1:B:106:VAL:N	2.33	0.43
2:W:21:DG:N2	3:X:4:DC:N3	2.61	0.43
1:A:89:ARG:NH1	1:B:92:ARG:HG2	2.33	0.43
1:A:144:TYR:CZ	1:A:145:GLN:O	2.71	0.43
1:C:98:LYS:HB2	1:C:98:LYS:HE2	1.33	0.43
1:A:73:VAL:HG11	1:A:87:GLY:HA3	1.99	0.43
1:D:92:ARG:HD3	1:D:92:ARG:HA	1.76	0.43
1:D:134:ALA:HB1	1:D:169:LEU:HG	1.99	0.43
1:A:89:ARG:CZ	1:B:89:ARG:HH21	2.31	0.43
1:B:77:SER:HA	1:B:109:ILE:HD13	2.00	0.43
2:W:5:DT:H2''	2:W:6:DG:O5'	2.18	0.43
2:W:22:DG:H1	3:X:3:DC:H42	1.67	0.43
1:A:199:LYS:H	1:A:199:LYS:HG2	1.56	0.43
1:B:68:GLY:O	1:B:70:LEU:HB2	2.19	0.43
2:W:21:DG:H2''	2:W:22:DG:C8	2.54	0.43
1:A:222:ARG:HB2	1:A:222:ARG:NH1	2.32	0.43
1:D:168:LEU:HD12	1:D:168:LEU:HA	1.78	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Y:7:DT:C2	3:Y:8:DT:C5	3.06	0.43
3:Y:21:DC:H2'	3:Y:22:DC:C6	2.54	0.43
1:A:144:TYR:CE1	1:C:89:ARG:HB2	2.54	0.43
1:B:115:ARG:HD3	1:B:115:ARG:HA	1.65	0.43
1:B:56:SER:HA	1:B:59:GLU:CD	2.43	0.43
1:B:85:ILE:CD1	1:B:90:ARG:HB3	2.47	0.43
1:B:115:ARG:NH1	1:D:75:VAL:HG22	2.34	0.43
1:C:97:ALA:C	1:C:98:LYS:HG3	2.44	0.43
1:C:70:LEU:HD13	1:C:71:GLN:HG3	2.00	0.43
1:C:141:LEU:HD21	1:C:148:GLN:HB3	2.01	0.43
3:X:20:DA:H2'	5:X:2005:HOH:O	2.18	0.43
1:A:148:GLN:HE22	2:W:5:DT:H3'	1.83	0.42
1:C:70:LEU:HD22	1:C:71:GLN:N	2.31	0.42
3:Y:15:DG:H2''	3:Y:16:DG:C8	2.53	0.42
2:Z:5:DT:H2''	2:Z:6:DG:H5'	2.00	0.42
1:C:129:ASN:HB2	5:C:2031:HOH:O	2.19	0.42
1:C:110:GLU:HG3	1:C:113:LYS:NZ	2.34	0.42
1:A:78:GLU:OE1	1:A:81:ARG:HB2	2.19	0.42
1:B:61:ALA:HB3	1:B:65:LYS:HZ1	1.83	0.42
3:Y:5:DT:H2''	3:Y:6:DG:C8	2.51	0.42
2:W:4:DG:H2'	2:W:5:DT:H71	2.02	0.42
2:Z:12:DC:H2'	2:Z:13:DG:C8	2.54	0.42
1:B:89:ARG:O	1:B:92:ARG:NE	2.52	0.42
3:Y:1:DT:H2'	3:Y:2:DC:C5	2.55	0.42
1:A:141:LEU:HD22	1:A:146:MET:HG2	2.00	0.42
1:C:42:MET:CG	1:C:145:GLN:HB2	2.38	0.42
1:D:74:LEU:HB3	1:D:83:HIS:CD2	2.55	0.42
1:A:71:GLN:HA	1:A:72:PRO:HD3	1.84	0.42
1:A:221:ALA:O	1:A:225:LYS:HD3	2.19	0.42
1:C:64:ILE:HG12	5:C:2008:HOH:O	2.19	0.42
1:D:110:GLU:C	1:D:112:GLU:N	2.77	0.42
1:B:160:ALA:N	5:B:2030:HOH:O	2.50	0.41
1:C:67:HIS:HA	5:C:2010:HOH:O	2.20	0.41
1:D:120:ILE:HD12	1:D:123:ILE:HG21	2.02	0.41
1:D:213:SER:OG	1:D:216:GLN:HG3	2.19	0.41
1:A:131:LEU:HB3	1:A:135:ARG:NH2	2.35	0.41
1:B:137:TYR:O	1:B:141:LEU:HG	2.20	0.41
1:B:152:SER:HB2	1:B:157:LYS:O	2.20	0.41
1:A:69:LEU:C	1:A:71:GLN:N	2.78	0.41
1:C:95:LYS:HB2	1:C:95:LYS:HE2	1.90	0.41
1:A:100:PRO:CD	1:A:101:THR:N	2.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:52:PHE:HB2	1:D:54:GLU:OE2	2.21	0.41
3:X:21:DC:H2''	3:X:22:DC:O4'	2.20	0.41
1:B:96:LEU:HA	5:B:2015:HOH:O	2.19	0.41
1:B:112:GLU:O	1:B:116:GLU:HB2	2.20	0.41
1:C:172:SER:HB3	1:C:200:GLN:OE1	2.20	0.41
1:D:103:LYS:O	1:D:104:ALA:C	2.63	0.41
1:D:151:LEU:O	1:D:155:VAL:HG13	2.20	0.41
1:A:69:LEU:HG	1:A:73:VAL:HG23	2.03	0.41
1:B:106:VAL:HG22	5:B:2018:HOH:O	2.21	0.41
1:D:65:LYS:NZ	1:D:67:HIS:CG	2.89	0.41
1:D:77:SER:C	1:D:78:GLU:OE1	2.63	0.41
1:D:176:GLN:O	1:D:180:LEU:HG	2.20	0.41
2:Z:5:DT:H2''	2:Z:6:DG:C5'	2.51	0.41
1:A:71:GLN:HB2	1:B:75:VAL:HG12	2.02	0.41
1:B:66:GLU:HG3	1:B:99:MET:SD	2.61	0.41
3:X:12:DC:H2'	3:X:13:DG:C8	2.56	0.41
1:A:67:HIS:HD2	1:A:68:GLY:N	2.18	0.41
1:A:102:ILE:HA	1:A:102:ILE:HD13	1.40	0.41
1:A:110:GLU:HB3	1:A:113:LYS:HG3	2.03	0.41
1:A:138:LYS:O	1:A:142:GLU:N	2.53	0.41
1:A:190:LYS:O	1:A:193:VAL:HG13	2.21	0.41
1:B:103:LYS:HE3	1:B:103:LYS:HB3	1.75	0.41
1:B:106:VAL:O	1:B:107:VAL:HG12	2.21	0.41
1:C:131:LEU:HB2	5:C:2031:HOH:O	2.21	0.41
1:C:177:ASN:O	1:C:181:GLU:HB2	2.21	0.41
1:D:79:ASN:C	1:D:81:ARG:N	2.77	0.41
1:D:110:GLU:C	1:D:110:GLU:CD	2.88	0.41
1:D:190:LYS:CE	2:Z:9:DC:H41	2.34	0.41
1:D:213:SER:N	1:D:216:GLN:OE1	2.44	0.41
1:B:170:THR:O	1:B:170:THR:OG1	2.35	0.41
1:C:110:GLU:O	1:C:113:LYS:NZ	2.52	0.41
1:C:195:LEU:HD22	1:C:199:LYS:HD2	2.03	0.41
1:D:131:LEU:HD12	1:D:180:LEU:CD2	2.51	0.41
1:B:108:ASP:OD1	1:B:108:ASP:N	2.52	0.40
1:D:75:VAL:HG11	1:D:95:LYS:CE	2.47	0.40
1:B:69:LEU:HD12	1:B:69:LEU:C	2.45	0.40
1:B:138:LYS:HE2	1:B:169:LEU:HD12	2.03	0.40
1:D:52:PHE:N	1:D:52:PHE:HD1	2.19	0.40
1:A:184:ILE:HD11	1:A:207:ILE:CG2	2.51	0.40
1:C:52:PHE:CZ	1:C:70:LEU:HD21	2.57	0.40
1:A:195:LEU:HD12	1:A:195:LEU:HA	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:TYR:H	1:C:46:TYR:HD1	1.70	0.40
1:C:80:GLY:HA2	1:C:85:ILE:HG22	2.04	0.40
3:Y:23:DC:C4	3:Y:24:DT:O4	2.75	0.40
1:A:89:ARG:NH2	1:B:89:ARG:HH21	2.18	0.40
1:A:90:ARG:HG3	1:B:89:ARG:NH1	2.37	0.40
1:A:140:LEU:HD13	1:A:151:LEU:HD13	2.04	0.40
1:B:50:LYS:HD3	1:B:51:VAL:N	2.37	0.40
1:B:57:LEU:HD12	1:B:58:GLU:H	1.86	0.40
1:B:115:ARG:CZ	1:D:76:VAL:HG23	2.52	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	170/240 (71%)	147 (86%)	18 (11%)	5 (3%)	3	19
1	B	181/240 (75%)	152 (84%)	24 (13%)	5 (3%)	4	20
1	C	191/240 (80%)	164 (86%)	19 (10%)	8 (4%)	2	13
1	D	174/240 (72%)	151 (87%)	18 (10%)	5 (3%)	3	19
All	All	716/960 (75%)	614 (86%)	79 (11%)	23 (3%)	3	18

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	69	LEU
1	C	81	ARG
1	D	63	SER
1	D	79	ASN
1	D	99	MET
1	D	102	ILE

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Mol	Chain	Res	Type
1	A	101	THR
1	B	70	LEU
1	C	51	VAL
1	A	99	MET
1	B	142	GLU
1	C	76	VAL
1	A	68	GLY
1	B	220	LEU
1	C	50	LYS
1	C	97	ALA
1	B	221	ALA
1	C	80	GLY
1	C	85	ILE
1	C	100	PRO
1	A	216	GLN
1	A	80	GLY
1	D	80	GLY

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	153/212 (72%)	109 (71%)	44 (29%)	0	1
1	B	164/212 (77%)	113 (69%)	51 (31%)	0	0
1	C	172/212 (81%)	116 (67%)	56 (33%)	0	0
1	D	157/212 (74%)	97 (62%)	60 (38%)	0	0
All	All	646/848 (76%)	435 (67%)	211 (33%)	0	0

All (211) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	LEU
1	A	64	ILE
1	A	69	LEU
1	A	70	LEU

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Mol	Chain	Res	Type
1	A	73	VAL
1	A	74	LEU
1	A	76	VAL
1	A	88	GLU
1	A	89	ARG
1	A	90	ARG
1	A	91	LEU
1	A	94	SER
1	A	95	LYS
1	A	98	LYS
1	A	102	ILE
1	A	103	LYS
1	A	105	ILE
1	A	107	VAL
1	A	108	ASP
1	A	114	MET
1	A	115	ARG
1	A	119	LEU
1	A	124	GLN
1	A	128	LEU
1	A	132	GLU
1	A	141	LEU
1	A	145	GLN
1	A	147	THR
1	A	159	ARG
1	A	176	GLN
1	A	180	LEU
1	A	182	GLU
1	A	183	LYS
1	A	184	ILE
1	A	191	VAL
1	A	193	VAL
1	A	195	LEU
1	A	196	ASP
1	A	199	LYS
1	A	200	GLN
1	A	202	LEU
1	A	215	ARG
1	A	222	ARG
1	A	226	ILE
1	B	46	TYR
1	B	49	ARG

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Mol	Chain	Res	Type
1	B	50	LYS
1	B	54	GLU
1	B	56	SER
1	B	57	LEU
1	B	58	GLU
1	B	59	GLU
1	B	62	GLN
1	B	64	ILE
1	B	67	HIS
1	B	70	LEU
1	B	71	GLN
1	B	73	VAL
1	B	74	LEU
1	B	76	VAL
1	B	81	ARG
1	B	89	ARG
1	B	91	LEU
1	B	94	SER
1	B	96	LEU
1	B	98	LYS
1	B	99	MET
1	B	101	THR
1	B	105	ILE
1	B	106	VAL
1	B	109	ILE
1	B	113	LYS
1	B	114	MET
1	B	116	GLU
1	B	120	ILE
1	B	122	ASN
1	B	123	ILE
1	B	124	GLN
1	B	125	ARG
1	B	128	LEU
1	B	139	GLU
1	B	140	LEU
1	B	150	GLU
1	B	165	ILE
1	B	168	LEU
1	B	180	LEU
1	B	183	LYS
1	B	196	ASP

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Mol	Chain	Res	Type
1	B	208	ILE
1	B	211	LYS
1	B	215	ARG
1	B	220	LEU
1	B	225	LYS
1	B	226	ILE
1	B	227	ASN
1	C	39	ASP
1	C	46	TYR
1	C	58	GLU
1	C	59	GLU
1	C	60	LEU
1	C	62	GLN
1	C	64	ILE
1	C	66	GLU
1	C	67	HIS
1	C	69	LEU
1	C	70	LEU
1	C	71	GLN
1	C	73	VAL
1	C	74	LEU
1	C	77	SER
1	C	78	GLU
1	C	81	ARG
1	C	85	ILE
1	C	88	GLU
1	C	90	ARG
1	C	91	LEU
1	C	92	ARG
1	C	95	LYS
1	C	98	LYS
1	C	102	ILE
1	C	103	LYS
1	C	105	ILE
1	C	110	GLU
1	C	113	LYS
1	C	115	ARG
1	C	123	ILE
1	C	129	ASN
1	C	132	GLU
1	C	135	ARG
1	C	139	GLU

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Mol	Chain	Res	Type
1	C	141	LEU
1	C	143	SER
1	C	145	GLN
1	C	147	THR
1	C	149	GLU
1	C	150	GLU
1	C	151	LEU
1	C	156	LYS
1	C	157	LYS
1	C	159	ARG
1	C	165	ILE
1	C	172	SER
1	C	173	SER
1	C	179	LEU
1	C	190	LYS
1	C	191	VAL
1	C	193	VAL
1	C	201	GLU
1	C	206	SER
1	C	215	ARG
1	C	226	ILE
1	D	52	PHE
1	D	54	GLU
1	D	55	ASP
1	D	59	GLU
1	D	60	LEU
1	D	62	GLN
1	D	63	SER
1	D	64	ILE
1	D	65	LYS
1	D	66	GLU
1	D	67	HIS
1	D	70	LEU
1	D	73	VAL
1	D	74	LEU
1	D	75	VAL
1	D	76	VAL
1	D	77	SER
1	D	79	ASN
1	D	81	ARG
1	D	84	LEU
1	D	85	ILE

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Mol	Chain	Res	Type
1	D	90	ARG
1	D	91	LEU
1	D	92	ARG
1	D	95	LYS
1	D	96	LEU
1	D	99	MET
1	D	100	PRO
1	D	103	LYS
1	D	105	ILE
1	D	106	VAL
1	D	108	ASP
1	D	113	LYS
1	D	115	ARG
1	D	123	ILE
1	D	124	GLN
1	D	126	GLU
1	D	127	ASP
1	D	132	GLU
1	D	140	LEU
1	D	143	SER
1	D	144	TYR
1	D	146	MET
1	D	147	THR
1	D	148	GLN
1	D	149	GLU
1	D	150	GLU
1	D	151	LEU
1	D	155	VAL
1	D	157	LYS
1	D	166	MET
1	D	168	LEU
1	D	186	SER
1	D	190	LYS
1	D	191	VAL
1	D	196	ASP
1	D	202	LEU
1	D	215	ARG
1	D	216	GLN
1	D	226	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	GLN
1	B	47	GLN
1	B	62	GLN
1	B	79	ASN
1	B	148	GLN
1	B	161	HIS
1	D	79	ASN
1	D	148	GLN
1	D	205	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	$\# Z > 2$
4	SO4	C	1228	-	4,4,4	0.25	0	6,6,6	0.06	0
4	SO4	A	1228	-	4,4,4	0.25	0	6,6,6	0.20	0
4	SO4	A	1227	-	4,4,4	0.26	0	6,6,6	0.08	0
4	SO4	B	1228	-	4,4,4	0.26	0	6,6,6	0.08	0
4	SO4	W	1025	-	4,4,4	0.23	0	6,6,6	0.08	0
4	SO4	D	1227	-	4,4,4	0.25	0	6,6,6	0.21	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.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	172/240 (71%)	-1.08	0 100 100	23, 62, 107, 140	0
1	B	183/240 (76%)	-1.02	0 100 100	28, 71, 112, 142	0
1	C	193/240 (80%)	-0.95	0 100 100	24, 67, 126, 146	0
1	D	176/240 (73%)	-1.02	0 100 100	24, 70, 126, 137	0
2	W	24/24 (100%)	-0.60	0 100 100	27, 45, 84, 97	0
2	Z	24/24 (100%)	-0.86	0 100 100	26, 48, 109, 126	0
3	X	24/24 (100%)	-0.60	0 100 100	30, 51, 97, 102	0
3	Y	24/24 (100%)	-1.00	0 100 100	29, 59, 99, 106	0
All	All	820/1056 (77%)	-0.99	0 100 100	23, 65, 122, 146	0

There are no RSRZ outliers to report.

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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	A	1228	5/5	0.96	0.07	106,115,123,127	0
4	SO4	A	1227	5/5	0.97	0.09	121,123,124,129	0
4	SO4	B	1228	5/5	0.98	0.09	131,132,134,141	0
4	SO4	C	1228	5/5	0.98	0.06	149,149,152,156	0
4	SO4	D	1227	5/5	0.98	0.06	85,96,98,103	0
4	SO4	W	1025	5/5	0.98	0.12	124,124,129,132	0

6.5 Other polymers [i](#)

There are no such residues in this entry.