



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:09 AM GMT

PDB ID : 2PO3
Title : Crystal Structure Analysis of DesI in the presence of its TDP-sugar product
Authors : Burgie, E.S.; Holden, H.M.
Deposited on : 2007-04-25
Resolution : 2.10 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

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

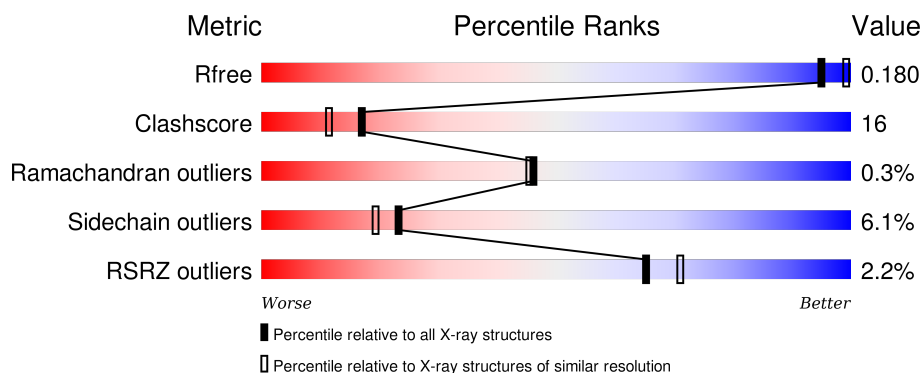
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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 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	424	 62% 25% 5% 7%
1	B	424	 4% 54% 29% 9% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6635 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.

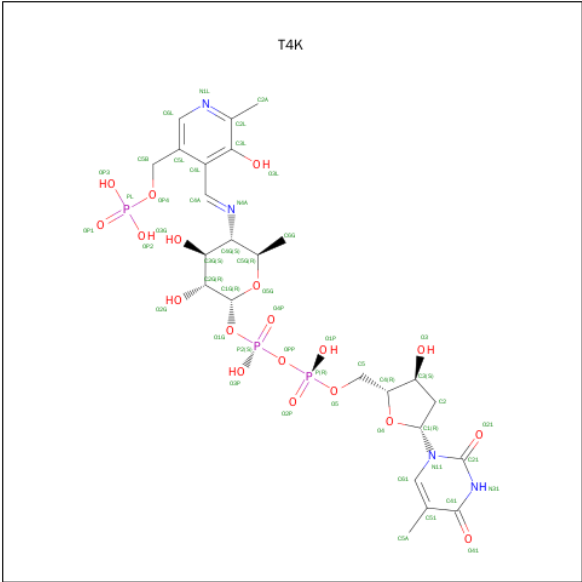
- Molecule 1 is a protein called 4-dehydrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	1	0
			2969	1847	561	551	10			
1	B	392	Total	C	N	O	S	0	3	0
			2973	1850	561	552	10			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1416	GLY	-	EXPRESSION TAG	UNP Q9ZGH0
A	1417	LEU	-	EXPRESSION TAG	UNP Q9ZGH0
A	1418	GLU	-	EXPRESSION TAG	UNP Q9ZGH0
A	1419	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
A	1420	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
A	1421	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
A	1422	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
A	1423	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
A	1424	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
B	2416	GLY	-	EXPRESSION TAG	UNP Q9ZGH0
B	2417	LEU	-	EXPRESSION TAG	UNP Q9ZGH0
B	2418	GLU	-	EXPRESSION TAG	UNP Q9ZGH0
B	2419	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
B	2420	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
B	2421	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
B	2422	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
B	2423	HIS	-	EXPRESSION TAG	UNP Q9ZGH0
B	2424	HIS	-	EXPRESSION TAG	UNP Q9ZGH0

- Molecule 2 is (2R,3R,4S,5S,6R)-3,4-DIHYDROXY-5-[(3-HYDROXY-2-METHYL-5-[(P HOSPHONOOXY)METHYL]PYRIDIN-4-YL}METHYL)IMINO]-6-METHYLTETRAHYDRO-2H-PYRAN-2-YL [(2R,3S,5R)-3-HYDROXY-5-(5-METHYL-2,4-DIOXO-3,4-DIHYDROPYRIMIDIN-1(2H)-YL)TETRAHYDROFURAN-2-YL]METHYL DIHYDROGEN DIPHOSPHATE (three-letter code: T4K) (formula: C₂₄H₃₅N₄O₁₉P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			50	24	4	19	3		
2	B	1	Total	C	N	O	P	0	0
			50	24	4	19	3		

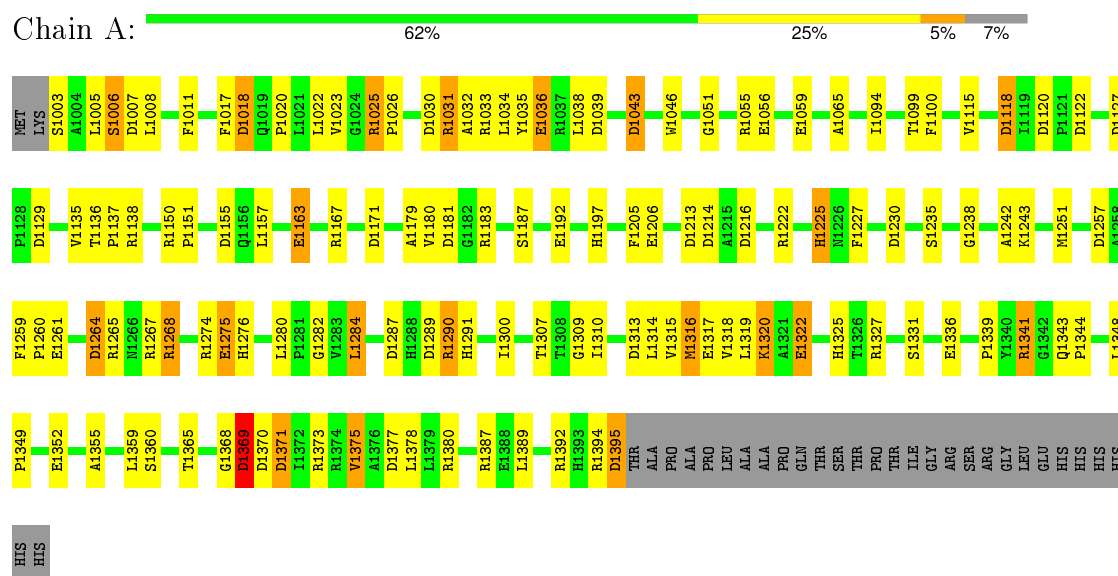
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	342	Total	O	0	0
			342	342		
3	B	251	Total	O	0	0
			251	251		

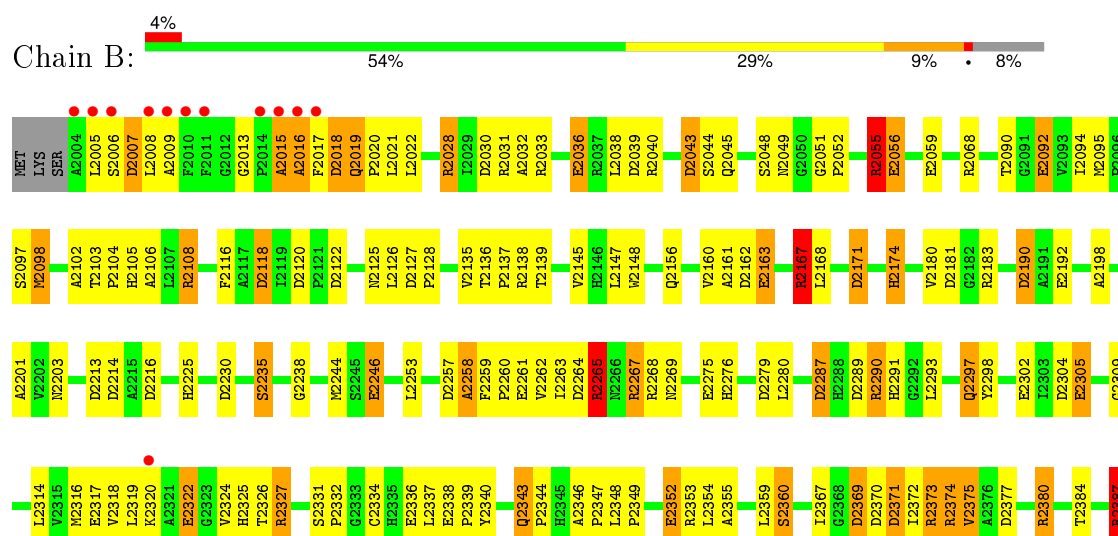
3 Residue-property plots

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: 4-dehydrase



- Molecule 1: 4-dehydrase



E2388	
L2389	
T2390	
A2391	
R2392	
H2393	
R2394	
D2395	
THR	
ALA	
PRO	
ALA	
PRO	
LEU	
ALA	
ALA	
PRO	
GLN	
THR	
SER	
THR	
PRO	
THR	
ILE	
GLY	
ARG	
SER	
ARG	
GLY	
LEU	
GLU	
HIS	
HIS	
HIS	
HIS	
HIS	

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.50 Å 66.80 Å 242.00 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.84 – 2.10 44.84 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.2 (44.84-2.10) 97.3 (44.84-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.63 (at 2.10 Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.180 , 0.248 0.177 , 0.180	Depositor DCC
R_{free} test set	5595 reflections (11.15%)	DCC
Wilson B-factor (Å ²)	25.2	Xtriage
Anisotropy	0.634	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 132.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 55830 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6635	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: T4K

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	0.87	10/3041 (0.3%)	1.41	58/4140 (1.4%)
1	B	0.91	18/3053 (0.6%)	1.51	73/4156 (1.8%)
All	All	0.89	28/6094 (0.5%)	1.46	131/8296 (1.6%)

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	2056	GLU	CD-OE2	7.08	1.33	1.25
1	B	2163	GLU	CD-OE2	6.65	1.32	1.25
1	A	1261	GLU	CD-OE2	6.41	1.32	1.25
1	A	1059	GLU	CD-OE2	6.30	1.32	1.25
1	A	1275	GLU	CD-OE2	6.26	1.32	1.25
1	A	1163	GLU	CD-OE2	6.26	1.32	1.25
1	B	2275	GLU	CD-OE2	6.24	1.32	1.25
1	A	1056	GLU	CD-OE2	6.14	1.32	1.25
1	B	2036	GLU	CD-OE2	6.08	1.32	1.25
1	B	2322	GLU	CD-OE2	6.05	1.32	1.25
1	B	2302	GLU	CD-OE2	5.87	1.32	1.25
1	B	2261	GLU	CD-OE2	5.83	1.32	1.25
1	B	2338	GLU	CD-OE2	5.78	1.32	1.25
1	B	2092	GLU	CD-OE2	5.66	1.31	1.25
1	A	1317	GLU	CD-OE2	5.63	1.31	1.25
1	B	2059	GLU	CD-OE2	5.59	1.31	1.25
1	B	2192	GLU	CD-OE2	5.56	1.31	1.25
1	A	1192	GLU	CD-OE2	5.49	1.31	1.25
1	B	2317	GLU	CD-OE2	5.41	1.31	1.25
1	B	2336	GLU	CD-OE2	5.37	1.31	1.25
1	A	1206	GLU	CD-OE2	5.34	1.31	1.25
1	B	2352	GLU	CD-OE2	5.32	1.31	1.25
1	B	2305	GLU	CD-OE2	5.29	1.31	1.25
1	B	2388[A]	GLU	CD-OE2	5.22	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	2388[B]	GLU	CD-OE2	5.22	1.31	1.25
1	B	2246	GLU	CD-OE2	5.10	1.31	1.25
1	A	1322	GLU	CD-OE2	5.06	1.31	1.25
1	A	1036	GLU	CD-OE2	5.01	1.31	1.25

All (131) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2394	ARG	NE-CZ-NH1	9.44	125.02	120.30
1	B	2214	ASP	CB-CG-OD2	-9.14	110.08	118.30
1	A	1043	ASP	CB-CG-OD2	-8.91	110.28	118.30
1	B	2214	ASP	CB-CG-OD1	8.84	126.26	118.30
1	B	2127	ASP	CB-CG-OD2	-8.64	110.52	118.30
1	A	1377	ASP	CB-CG-OD2	-8.44	110.70	118.30
1	B	2033	ARG	NE-CZ-NH1	8.38	124.49	120.30
1	B	2369	ASP	CB-CG-OD2	-8.36	110.78	118.30
1	B	2373	ARG	NE-CZ-NH1	8.02	124.31	120.30
1	B	2127	ASP	CB-CG-OD1	7.93	125.44	118.30
1	B	2387	ARG	NE-CZ-NH1	7.83	124.22	120.30
1	B	2327	ARG	NE-CZ-NH1	7.75	124.17	120.30
1	A	1213	ASP	CB-CG-OD2	-7.73	111.34	118.30
1	B	2181	ASP	CB-CG-OD2	-7.62	111.44	118.30
1	B	2171	ASP	CB-CG-OD1	7.46	125.02	118.30
1	A	1395	ASP	CB-CG-OD2	-7.33	111.70	118.30
1	A	1394	ARG	NE-CZ-NH1	7.28	123.94	120.30
1	A	1371	ASP	CB-CG-OD2	-7.25	111.78	118.30
1	A	1043	ASP	CB-CG-OD1	7.18	124.76	118.30
1	B	2122	ASP	CB-CG-OD2	-7.16	111.86	118.30
1	B	2395	ASP	CB-CG-OD2	-7.10	111.91	118.30
1	B	2120	ASP	CB-CG-OD2	-7.00	112.00	118.30
1	B	2039	ASP	CB-CG-OD2	-6.97	112.03	118.30
1	B	2290	ARG	NE-CZ-NH1	6.95	123.78	120.30
1	B	2162	ASP	CB-CG-OD2	-6.88	112.11	118.30
1	B	2304	ASP	CB-CG-OD1	6.87	124.48	118.30
1	B	2287	ASP	CB-CG-OD2	-6.81	112.17	118.30
1	B	2007	ASP	CB-CG-OD2	-6.80	112.18	118.30
1	B	2353	ARG	NE-CZ-NH1	6.79	123.70	120.30
1	B	2373	ARG	NE-CZ-NH2	-6.79	116.91	120.30
1	A	1120	ASP	CB-CG-OD1	6.77	124.39	118.30
1	A	1213	ASP	CB-CG-OD1	6.74	124.36	118.30
1	B	2118	ASP	CB-CG-OD2	-6.74	112.24	118.30
1	B	2015	ALA	O-C-N	6.70	133.43	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2289	ASP	CB-CG-OD2	-6.63	112.33	118.30
1	B	2171	ASP	CB-CG-OD2	-6.60	112.36	118.30
1	A	1118	ASP	CB-CG-OD2	-6.54	112.41	118.30
1	A	1313	ASP	CB-CG-OD2	-6.54	112.41	118.30
1	A	1222	ARG	NE-CZ-NH1	6.53	123.57	120.30
1	B	2122	ASP	CB-CG-OD1	6.52	124.17	118.30
1	A	1289	ASP	CB-CG-OD1	6.51	124.16	118.30
1	B	2377	ASP	CB-CG-OD1	6.48	124.14	118.30
1	A	1214	ASP	CB-CG-OD2	-6.48	112.47	118.30
1	A	1287	ASP	CB-CG-OD2	-6.42	112.52	118.30
1	B	2028	ARG	NE-CZ-NH1	6.33	123.47	120.30
1	A	1287	ASP	CB-CG-OD1	6.33	124.00	118.30
1	B	2264	ASP	CB-CG-OD2	-6.29	112.64	118.30
1	B	2395	ASP	CB-CG-OD1	6.29	123.96	118.30
1	B	2068	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	B	2371	ASP	CB-CG-OD2	-6.28	112.65	118.30
1	A	1181	ASP	CB-CG-OD1	6.28	123.95	118.30
1	A	1007	ASP	CB-CG-OD2	-6.27	112.66	118.30
1	B	2007	ASP	CB-CG-OD1	6.24	123.91	118.30
1	A	1018	ASP	CB-CG-OD2	-6.23	112.69	118.30
1	A	1264	ASP	CB-CG-OD2	-6.23	112.69	118.30
1	A	1341	ARG	NE-CZ-NH1	6.23	123.42	120.30
1	B	2120	ASP	CB-CG-OD1	6.23	123.91	118.30
1	A	1290	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	B	2216	ASP	CB-CG-OD1	6.18	123.86	118.30
1	A	1370	ASP	CB-CG-OD2	-6.16	112.76	118.30
1	B	2267	ARG	NE-CZ-NH1	6.16	123.38	120.30
1	B	2370	ASP	CB-CG-OD2	-6.02	112.89	118.30
1	B	2043	ASP	CB-CG-OD2	-6.01	112.89	118.30
1	A	1127	ASP	CB-CG-OD1	5.98	123.68	118.30
1	A	1289	ASP	CB-CG-OD2	-5.95	112.95	118.30
1	B	2258	ALA	N-CA-CB	-5.94	101.78	110.10
1	B	2183	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	B	2268	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	B	2392	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	B	2216	ASP	CB-CG-OD2	-5.86	113.02	118.30
1	B	2380	ARG	NE-CZ-NH1	5.85	123.22	120.30
1	A	1030	ASP	CB-CG-OD2	-5.82	113.07	118.30
1	A	1274	ARG	NE-CZ-NH1	5.81	123.21	120.30
1	B	2030	ASP	CB-CG-OD2	-5.78	113.10	118.30
1	A	1214	ASP	CB-CG-OD1	5.76	123.49	118.30
1	B	2340	TYR	CB-CG-CD2	-5.75	117.55	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1377	ASP	CB-CG-OD1	5.75	123.47	118.30
1	A	1181	ASP	CB-CG-OD2	-5.74	113.14	118.30
1	A	1394	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	B	2055	ARG	NE-CZ-NH2	-5.71	117.44	120.30
1	A	1369	ASP	CB-CG-OD2	-5.71	113.17	118.30
1	B	2257	ASP	CB-CG-OD1	5.68	123.42	118.30
1	A	1395	ASP	CB-CG-OD1	5.68	123.41	118.30
1	A	1122	ASP	CB-CG-OD1	5.64	123.37	118.30
1	B	2370	ASP	CB-CG-OD1	5.63	123.37	118.30
1	A	1167	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	A	1257	ASP	CB-CG-OD1	5.59	123.33	118.30
1	A	1171	ASP	CB-CG-OD1	5.58	123.32	118.30
1	B	2304	ASP	CB-CG-OD2	-5.56	113.30	118.30
1	A	1127	ASP	CB-CG-OD2	-5.54	113.31	118.30
1	B	2265	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	2343	GLN	CB-CA-C	5.48	121.36	110.40
1	B	2387	ARG	NE-CZ-NH2	-5.43	117.58	120.30
1	A	1230	ASP	CB-CG-OD2	-5.43	113.41	118.30
1	B	2183	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	B	2213	ASP	CB-CG-OD2	-5.43	113.42	118.30
1	B	2028	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	1039	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	A	1055	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	B	2190	ASP	CB-CG-OD2	-5.36	113.48	118.30
1	B	2230	ASP	CB-CG-OD2	-5.34	113.49	118.30
1	A	1129	ASP	CB-CG-OD2	-5.33	113.50	118.30
1	B	2369	ASP	CB-CG-OD1	5.30	123.07	118.30
1	A	1122	ASP	CB-CG-OD2	-5.29	113.54	118.30
1	B	2016	ALA	N-CA-CB	5.29	117.51	110.10
1	B	2268	ARG	NE-CZ-NH2	-5.29	117.66	120.30
1	A	1365	THR	CA-CB-CG2	-5.28	105.00	112.40
1	B	2291	HIS	CA-CB-CG	-5.28	104.62	113.60
1	A	1230	ASP	CB-CG-OD1	5.24	123.02	118.30
1	A	1155	ASP	CB-CG-OD1	5.22	123.00	118.30
1	B	2264	ASP	CB-CG-OD1	5.22	123.00	118.30
1	A	1025	ARG	NE-CZ-NH2	-5.19	117.71	120.30
1	A	1120	ASP	CB-CG-OD2	-5.19	113.63	118.30
1	B	2380	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	A	1118	ASP	CB-CG-OD1	5.18	122.96	118.30
1	A	1007	ASP	CB-CG-OD1	5.17	122.96	118.30
1	B	2377	ASP	CB-CG-OD2	-5.17	113.65	118.30
1	B	2257	ASP	CB-CG-OD2	-5.16	113.66	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1268	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	B	2167	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	A	1138[A]	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	A	1138[B]	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	B	2343	GLN	N-CA-CB	5.12	119.81	110.60
1	B	2279	ASP	CB-CG-OD2	-5.10	113.71	118.30
1	B	2043	ASP	CB-CG-OD1	5.07	122.86	118.30
1	A	1216	ASP	CB-CG-OD2	-5.05	113.75	118.30
1	A	1030	ASP	CB-CG-OD1	5.05	122.84	118.30
1	A	1369	ASP	CB-CG-OD1	5.04	122.84	118.30
1	A	1375	VAL	CA-CB-CG2	5.02	118.42	110.90
1	A	1216	ASP	CB-CG-OD1	5.01	122.81	118.30
1	B	2068	ARG	NE-CZ-NH2	-5.00	117.80	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2969	0	2896	71	0
1	B	2973	0	2901	126	0
2	A	50	0	31	6	0
2	B	50	0	31	1	0
3	A	342	0	0	11	0
3	B	251	0	0	13	0
All	All	6635	0	5859	194	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 16.

All (194) 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:B:2201:ALA:HB2	1:B:2297[B]:GLN:HG3	1.41	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2009:ALA:HA	1:B:2013:GLY:O	1.67	0.94
1:B:2017:PHE:HB2	1:B:2374:ARG:CZ	1.99	0.93
1:A:1022:LEU:HD23	1:A:1325:HIS:HB2	1.55	0.89
1:B:2167:ARG:HG3	1:B:2167:ARG:HH11	1.42	0.82
1:A:1348:LEU:O	1:A:1352:GLU:HG3	1.82	0.79
1:B:2032:ALA:O	1:B:2036:GLU:HG3	1.84	0.78
1:A:1320:LYS:HE3	1:A:1325:HIS:HD2	1.49	0.78
1:B:2343:GLN:HB2	1:B:2344:PRO:HD2	1.65	0.77
1:B:2320:LYS:HG3	3:B:3346:HOH:O	1.85	0.77
1:B:2348:LEU:O	1:B:2352:GLU:HG3	1.85	0.75
1:A:1264:ASP:HB3	3:A:3574:HOH:O	1.86	0.75
1:B:2156:GLN:O	1:B:2160:VAL:HG23	1.87	0.75
1:A:1290:ARG:HD3	3:A:4667:HOH:O	1.87	0.74
1:A:1369:ASP:HB3	3:A:3577:HOH:O	1.87	0.74
1:B:2201:ALA:CB	1:B:2297[B]:GLN:HG3	2.17	0.73
1:B:2174:HIS:HB3	1:B:2298:TYR:HE1	1.52	0.72
1:A:1023:VAL:HG11	1:A:1327:ARG:HG2	1.69	0.72
1:B:2201:ALA:HB2	1:B:2297[B]:GLN:CG	2.18	0.72
1:B:2136:THR:HB	1:B:2137:PRO:HD2	1.71	0.72
1:B:2118:ASP:OD2	1:B:2349:PRO:HD2	1.91	0.71
1:B:2314:LEU:O	1:B:2318:VAL:HG23	1.91	0.71
1:B:2016:ALA:HB3	1:B:2322:GLU:HG3	1.73	0.70
1:B:2018:ASP:O	1:B:2020:PRO:HD3	1.91	0.69
1:B:2167:ARG:NH1	1:B:2167:ARG:HG3	2.04	0.67
1:B:2017:PHE:HB2	1:B:2374:ARG:NH1	2.10	0.67
1:B:2040[A]:ARG:NH2	1:B:2052:PRO:HG2	2.10	0.67
1:A:1320:LYS:CE	1:A:1325:HIS:HD2	2.08	0.66
1:B:2016:ALA:CB	1:B:2322:GLU:HG3	2.26	0.66
1:A:1023:VAL:CG1	1:A:1327:ARG:HG2	2.26	0.66
1:A:1205:PHE:CE1	1:A:1251:MET:HB2	2.30	0.66
1:A:1315:VAL:O	1:A:1319:LEU:HD13	1.96	0.66
1:B:2160:VAL:O	1:B:2163:GLU:HG2	1.96	0.65
1:B:2280:LEU:HD21	1:B:2380:ARG:HA	1.78	0.65
1:B:2009:ALA:HB2	1:B:2015:ALA:HA	1.79	0.65
1:B:2007:ASP:O	1:B:2013:GLY:HA3	1.96	0.65
1:A:1320:LYS:HE3	1:A:1325:HIS:CD2	2.31	0.64
1:B:2056:GLU:HG2	1:B:2253:LEU:HD21	1.79	0.64
1:B:2343:GLN:HB2	1:B:2344:PRO:CD	2.27	0.63
1:A:1327:ARG:NH1	2:A:1500:T4K:O4P	2.29	0.63
1:A:1018:ASP:O	1:A:1020:PRO:HD3	1.98	0.63
1:B:2346:ALA:HB1	1:B:2347:PRO:HD2	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2016:ALA:CB	1:B:2322:GLU:CG	2.77	0.62
1:B:2028:ARG:HH12	1:B:2031:ARG:HH21	1.47	0.62
1:B:2259:PHE:HB3	1:B:2260:PRO:HD3	1.82	0.61
1:B:2022:LEU:HD23	1:B:2325:HIS:HB2	1.82	0.61
1:B:2016:ALA:HB3	1:B:2322:GLU:CG	2.29	0.61
1:A:1022:LEU:HD23	1:A:1325:HIS:CB	2.29	0.61
1:A:1373:ARG:NH1	3:A:3083:HOH:O	2.34	0.61
1:A:1227:PHE:CE2	1:B:2102:ALA:HB2	2.37	0.60
1:B:2380:ARG:O	1:B:2384:THR:HG23	2.01	0.60
1:B:2051:GLY:HA3	1:B:2246:GLU:CD	2.22	0.60
1:B:2028:ARG:NH1	1:B:2031:ARG:HH21	2.00	0.59
1:B:2287:ASP:O	1:B:2290:ARG:HG2	2.02	0.59
1:B:2393:HIS:O	1:B:2395:ASP:N	2.34	0.58
1:B:2371:ASP:O	1:B:2375:VAL:HG12	2.04	0.58
1:B:2244:MET:HG3	3:B:6230:HOH:O	2.03	0.58
1:B:2016:ALA:O	1:B:2017:PHE:HD1	1.86	0.58
1:B:2334:CYS:O	1:B:2337:LEU:HB2	2.03	0.58
1:B:2258:ALA:O	1:B:2262:VAL:HG23	2.04	0.57
1:B:2267:ARG:HH12	1:B:2293:LEU:HD12	1.69	0.57
1:B:2390:THR:HG22	1:B:2394:ARG:HD2	1.86	0.57
1:B:2017:PHE:CD2	1:B:2374:ARG:HD2	2.40	0.57
1:B:2108:ARG:NH1	1:B:2108:ARG:HG3	2.19	0.57
1:A:1322:GLU:HG2	3:A:5801:HOH:O	2.03	0.57
1:A:1043:ASP:OD1	1:B:2031:ARG:NH2	2.35	0.56
1:B:2108:ARG:HG3	1:B:2108:ARG:HH11	1.70	0.56
1:A:1011:PHE:CE2	1:A:1392:ARG:HG2	2.40	0.56
1:B:2056:GLU:HG2	1:B:2253:LEU:CD2	2.37	0.55
1:A:1280:LEU:HD21	1:A:1380:ARG:HA	1.88	0.55
1:B:2290:ARG:NH2	3:B:5604:HOH:O	2.39	0.55
1:A:1118:ASP:OD2	1:A:1349:PRO:HD2	2.06	0.55
1:B:2267:ARG:NH1	1:B:2293:LEU:HD12	2.23	0.54
1:A:1320:LYS:NZ	1:A:1325:HIS:HD2	2.06	0.54
1:B:2174:HIS:CB	1:B:2298:TYR:HE1	2.20	0.53
1:A:1344:PRO:HD3	3:A:4676:HOH:O	2.06	0.53
1:B:2138:ARG:NH1	3:B:3852:HOH:O	2.40	0.53
1:B:2319:LEU:HB3	1:B:2324:VAL:HB	1.91	0.53
1:B:2387:ARG:HD3	3:B:5610:HOH:O	2.08	0.53
1:A:1265:ARG:HD3	1:A:1268:ARG:CZ	2.39	0.53
1:B:2016:ALA:C	1:B:2017:PHE:HD1	2.13	0.52
1:B:2136:THR:HB	1:B:2137:PRO:CD	2.38	0.52
1:B:2094:ILE:HD12	1:B:2139:THR:HG21	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2135:VAL:HG12	1:B:2136:THR:N	2.24	0.52
1:A:1031:ARG:NH1	1:B:2043:ASP:OD1	2.39	0.52
1:B:2168:LEU:O	1:B:2190:ASP:HB2	2.10	0.51
1:A:1316:MET:HG2	1:A:1359:LEU:HD11	1.92	0.51
1:B:2309:GLY:HA3	1:B:2387:ARG:HD2	1.93	0.51
1:B:2343:GLN:HG3	3:B:4833:HOH:O	2.10	0.51
1:B:2098:MET:HE1	1:B:2355:ALA:HA	1.92	0.51
1:A:1331:SER:HB2	1:A:1355:ALA:O	2.10	0.50
1:A:1282:GLY:N	1:A:1307:THR:HG21	2.25	0.50
1:B:2097:SER:OG	1:B:2334:CYS:N	2.45	0.50
1:B:2005:LEU:HD22	1:B:2008:LEU:CD1	2.40	0.50
1:B:2393:HIS:C	1:B:2395:ASP:H	2.14	0.50
1:B:2326:THR:O	1:B:2327:ARG:HG2	2.12	0.50
1:B:2174:HIS:CG	1:B:2298:TYR:HE1	2.29	0.50
1:B:2016:ALA:H	1:B:2322:GLU:HG3	1.77	0.50
1:B:2235:SER:HB2	3:B:3256:HOH:O	2.11	0.50
1:A:1005:LEU:O	1:A:1008:LEU:HB2	2.11	0.50
1:A:1046:TRP:CE2	1:A:1051:GLY:HA2	2.47	0.49
1:A:1314:LEU:O	1:A:1318:VAL:HG23	2.13	0.49
1:B:2009:ALA:HB2	1:B:2015:ALA:CA	2.41	0.49
1:B:2265:ARG:HG2	1:B:2265:ARG:NH1	2.28	0.49
2:B:2500:T4K:O3G	2:B:2500:T4K:H4A1	2.13	0.48
1:B:2125:ASN:O	1:B:2126:LEU:C	2.51	0.48
1:A:1238:GLY:HA2	1:B:2339:PRO:HD3	1.96	0.48
1:A:1322:GLU:HG3	1:A:1378:LEU:HD11	1.95	0.48
1:A:1368:GLY:O	1:A:1371:ASP:HB2	2.13	0.48
1:B:2297[B]:GLN:NE2	3:B:3801:HOH:O	2.42	0.48
1:B:2103:THR:HB	1:B:2104:PRO:HD3	1.95	0.48
1:A:1284:LEU:HD21	3:A:3041:HOH:O	2.13	0.48
1:B:2005:LEU:HD22	1:B:2008:LEU:HD11	1.95	0.47
1:A:1197:HIS:HE1	2:A:1500:T4K:H5A3	1.79	0.47
1:B:2265:ARG:HD2	1:B:2269:ASN:OD1	2.15	0.47
1:B:2108:ARG:CG	1:B:2108:ARG:HH11	2.26	0.47
1:B:2316:MET:HG2	1:B:2359:LEU:HD13	1.95	0.47
1:B:2009:ALA:CB	1:B:2015:ALA:HA	2.42	0.47
1:B:2095:MET:O	1:B:2116:PHE:HA	2.15	0.47
1:B:2017:PHE:HB2	1:B:2374:ARG:NH2	2.29	0.47
1:B:2043:ASP:HB2	3:B:3809:HOH:O	2.14	0.46
1:A:1275:GLU:HG3	1:A:1276:HIS:N	2.29	0.46
1:B:2009:ALA:HA	1:B:2013:GLY:C	2.33	0.46
1:A:1259:PHE:HB3	1:A:1260:PRO:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2040[A]:ARG:HH22	1:B:2052:PRO:HG2	1.80	0.46
1:A:1136:THR:HB	1:A:1137:PRO:CD	2.45	0.46
1:B:2346:ALA:HB1	1:B:2347:PRO:CD	2.44	0.45
1:B:2038:LEU:HA	1:B:2038:LEU:HD23	1.67	0.45
1:A:1197:HIS:CE1	2:A:1500:T4K:H5A3	2.51	0.45
1:B:2055:ARG:HA	1:B:2055:ARG:HD2	1.64	0.45
1:B:2174:HIS:HB3	1:B:2298:TYR:CE1	2.42	0.45
1:B:2316:MET:HG2	1:B:2359:LEU:CD1	2.46	0.45
1:B:2147:LEU:O	1:B:2148:TRP:HB2	2.16	0.45
1:B:2160:VAL:O	1:B:2161:ALA:C	2.54	0.45
1:B:2263:ILE:O	1:B:2267:ARG:HG3	2.16	0.45
1:B:2044:SER:O	1:B:2045:GLN:HB2	2.17	0.45
1:B:2019:GLN:HE21	1:B:2019:GLN:HB3	1.67	0.45
1:A:1310:ILE:HA	3:A:3044:HOH:O	2.17	0.44
1:A:1017:PHE:HB3	3:A:4313:HOH:O	2.17	0.44
1:B:2090:THR:HG23	3:B:4621:HOH:O	2.16	0.44
1:A:1336:GLU:O	1:A:1341:ARG:NH1	2.48	0.44
1:B:2387:ARG:HB2	3:B:5610:HOH:O	2.16	0.44
1:B:2048:SER:OG	1:B:2049:ASN:N	2.51	0.44
1:B:2180:VAL:HG13	1:B:2180:VAL:O	2.17	0.44
1:A:1151:PRO:HG2	1:A:1180:VAL:HB	2.00	0.44
1:A:1135:VAL:HG12	1:A:1136:THR:N	2.33	0.43
1:A:1023:VAL:HG12	1:A:1325:HIS:O	2.18	0.43
1:B:2145:VAL:HG22	1:B:2171:ASP:HB3	2.00	0.43
1:A:1150:ARG:HB3	1:A:1291:HIS:HB2	2.00	0.43
1:B:2156:GLN:HG3	3:B:3500:HOH:O	2.18	0.43
1:B:2021:LEU:HD22	1:B:2367:ILE:CD1	2.49	0.43
1:B:2276:HIS:HD2	1:B:2372:ILE:HG22	1.83	0.43
1:A:1094:ILE:HA	1:A:1115:VAL:O	2.19	0.43
1:B:2201:ALA:CA	1:B:2297[B]:GLN:HG3	2.49	0.43
1:B:2135:VAL:CG1	1:B:2136:THR:N	2.82	0.43
1:A:1135:VAL:O	1:A:1136:THR:HG22	2.19	0.43
1:A:1327:ARG:NH1	2:A:1500:T4K:H52	2.34	0.42
1:B:2174:HIS:CG	1:B:2298:TYR:CE1	3.06	0.42
1:B:2326:THR:HB	1:B:2360:SER:O	2.20	0.42
2:A:1500:T4K:H4A1	2:A:1500:T4K:HO3G	1.85	0.42
1:A:1179:ALA:HA	1:A:1183:ARG:O	2.20	0.42
1:B:2009:ALA:HB2	1:B:2015:ALA:N	2.35	0.42
1:B:2098:MET:HE2	1:B:2355:ALA:HB2	2.02	0.42
1:A:1343:GLN:NE2	3:A:4388:HOH:O	2.51	0.42
1:A:1280:LEU:HA	1:A:1280:LEU:HD12	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1046:TRP:CZ2	1:A:1051:GLY:HA2	2.55	0.41
1:B:2316:MET:HE3	1:B:2316:MET:HB3	1.88	0.41
1:B:2016:ALA:HB2	1:B:2322:GLU:CG	2.48	0.41
1:A:1157:LEU:HA	1:A:1157:LEU:HD23	1.93	0.41
1:A:1032:ALA:O	1:A:1036:GLU:HG3	2.20	0.41
1:B:2305:GLU:O	1:B:2309:GLY:N	2.52	0.41
1:B:2092:GLU:HB2	1:B:2139:THR:HA	2.01	0.41
1:B:2316:MET:HE2	3:B:3902:HOH:O	2.19	0.41
1:A:1316:MET:HG2	1:A:1359:LEU:CD1	2.50	0.41
1:B:2016:ALA:O	1:B:2017:PHE:CD1	2.72	0.41
1:B:2103:THR:HB	1:B:2104:PRO:CD	2.51	0.41
1:B:2354:LEU:O	1:B:2354:LEU:HD12	2.20	0.41
1:A:1006:SER:HB3	1:A:1392:ARG:HH22	1.86	0.41
1:A:1035:TYR:HD1	1:A:1038:LEU:HD12	1.85	0.41
1:A:1267:ARG:NH1	3:A:3023:HOH:O	2.45	0.41
1:B:2016:ALA:CB	1:B:2322:GLU:HG2	2.49	0.41
1:A:1034:LEU:HD13	1:A:1251:MET:HA	2.03	0.41
1:A:1389:LEU:O	1:A:1389:LEU:HD23	2.21	0.41
1:A:1309:GLY:HA3	1:A:1387:ARG:HG3	2.03	0.41
1:A:1225:HIS:HB2	1:A:1242:ALA:HB3	2.03	0.41
1:A:1339:PRO:HD3	1:B:2238:GLY:HA2	2.02	0.40
1:A:1025:ARG:HG2	1:A:1026:PRO:HD2	2.03	0.40
1:A:1316:MET:HE3	1:A:1316:MET:HB3	1.89	0.40
1:B:2331:SER:HA	1:B:2332:PRO:HA	1.66	0.40
1:A:1099:THR:OG1	1:A:1100:PHE:N	2.54	0.40
1:A:1327:ARG:HH11	2:A:1500:T4K:H52	1.87	0.40
1:B:2265:ARG:CG	1:B:2265:ARG:HH11	2.35	0.40
1:A:1065:ALA:O	1:A:1187:SER:HB3	2.21	0.40
1:B:2105:HIS:O	1:B:2106:ALA:C	2.59	0.40

There are no symmetry-related clashes.

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	392/424 (92%)	379 (97%)	13 (3%)	0	100	100
1	B	393/424 (93%)	372 (95%)	19 (5%)	2 (0%)	34	30
All	All	785/848 (93%)	751 (96%)	32 (4%)	2 (0%)	46	45

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2394	ARG
1	B	2198	ALA

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	297/321 (92%)	281 (95%)	16 (5%)	27	24
1	B	298/321 (93%)	277 (93%)	21 (7%)	19	15
All	All	595/642 (93%)	558 (94%)	37 (6%)	23	19

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1003	SER
1	A	1006	SER
1	A	1031	ARG
1	A	1033	ARG
1	A	1163	GLU
1	A	1225	HIS
1	A	1235	SER
1	A	1243	LYS
1	A	1284	LEU
1	A	1300	ILE
1	A	1316	MET
1	A	1320	LYS
1	A	1360	SER
1	A	1369	ASP

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Mol	Chain	Res	Type
1	A	1375	VAL
1	A	1395	ASP
1	B	2006	SER
1	B	2018	ASP
1	B	2019	GLN
1	B	2055	ARG
1	B	2098	MET
1	B	2108	ARG
1	B	2128	PRO
1	B	2167	ARG
1	B	2174	HIS
1	B	2203	ASN
1	B	2225	HIS
1	B	2235	SER
1	B	2265	ARG
1	B	2297[A]	GLN
1	B	2297[B]	GLN
1	B	2360	SER
1	B	2369	ASP
1	B	2373	ARG
1	B	2374	ARG
1	B	2375	VAL
1	B	2387	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1069	HIS
1	A	1325	HIS
1	A	1343	GLN
1	B	2019	GLN
1	B	2130	GLN
1	B	2156	GLN
1	B	2276	HIS

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 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	T4K	A	1500	-	45,53,53	1.86	8 (17%)	60,81,81	1.70	7 (11%)
2	T4K	B	2500	-	45,53,53	1.94	7 (15%)	60,81,81	1.84	9 (15%)

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	T4K	A	1500	-	-	1/28/64/64	0/4/4/4
2	T4K	B	2500	-	-	1/28/64/64	0/4/4/4

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2500	T4K	C4G-N4A	-8.72	1.39	1.46
2	A	1500	T4K	C4G-N4A	-7.03	1.41	1.46
2	B	2500	T4K	C4L-C4A	-3.01	1.41	1.46
2	B	2500	T4K	PL-OP3	2.32	1.63	1.54
2	A	1500	T4K	C5G-C4G	2.43	1.56	1.52
2	B	2500	T4K	PL-OP1	2.58	1.59	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1500	T4K	P-O2P	2.59	1.60	1.51
2	A	1500	T4K	C61-N11	2.81	1.39	1.35
2	B	2500	T4K	P2-O4P	3.05	1.62	1.51
2	B	2500	T4K	P-O2P	3.09	1.62	1.51
2	A	1500	T4K	P2-O4P	3.19	1.62	1.51
2	B	2500	T4K	C4A-N4A	3.43	1.33	1.27
2	A	1500	T4K	PL-OP1	3.50	1.62	1.51
2	A	1500	T4K	C4L-C3L	3.55	1.45	1.40
2	A	1500	T4K	C4A-N4A	3.70	1.34	1.27

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2500	T4K	P-OPP-P2	-7.63	111.30	132.73
2	A	1500	T4K	P-OPP-P2	-5.79	116.46	132.73
2	B	2500	T4K	C51-C41-N31	-5.00	119.57	125.14
2	A	1500	T4K	C51-C41-N31	-4.99	119.59	125.14
2	A	1500	T4K	C1G-O5G-C5G	-3.26	108.08	113.64
2	B	2500	T4K	C1G-O5G-C5G	-3.08	108.38	113.64
2	A	1500	T4K	OP4-PL-OP1	2.06	112.38	107.14
2	B	2500	T4K	O1G-C1G-C2G	2.10	112.31	108.39
2	B	2500	T4K	O5G-C5G-C6G	2.15	111.34	106.64
2	A	1500	T4K	OPP-P-O5	2.17	108.69	102.94
2	B	2500	T4K	OP3-PL-OP4	2.41	113.52	106.56
2	B	2500	T4K	OPP-P-O5	2.49	109.55	102.94
2	B	2500	T4K	C4G-N4A-C4A	3.19	123.07	118.70
2	A	1500	T4K	C4G-N4A-C4A	4.56	124.95	118.70
2	A	1500	T4K	C41-N31-C21	6.39	120.77	115.25
2	B	2500	T4K	C41-N31-C21	6.46	120.83	115.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1500	T4K	C3G-C4G-N4A-C4A
2	B	2500	T4K	C3G-C4G-N4A-C4A

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1500	T4K	6	0
2	B	2500	T4K	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	393/424 (92%)	-0.65	0 100 100	13, 27, 59, 93	0
1	B	392/424 (92%)	-0.26	17 (4%) 39 48	12, 32, 69, 99	0
All	All	785/848 (92%)	-0.45	17 (2%) 65 71	12, 30, 63, 99	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2011	PHE	5.0
1	B	2005	LEU	4.8
1	B	2017	PHE	3.8
1	B	2008	LEU	3.7
1	B	2010	PHE	3.0
1	B	2004	ALA	3.0
1	B	2391	ALA	2.9
1	B	2015	ALA	2.7
1	B	2393	HIS	2.7
1	B	2016	ALA	2.6
1	B	2394	ARG	2.6
1	B	2320	LYS	2.5
1	B	2395	ASP	2.5
1	B	2009	ALA	2.5
1	B	2389	LEU	2.4
1	B	2006	SER	2.4
1	B	2014	PRO	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, 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
2	T4K	B	2500	50/50	0.95	0.12	0.52	14,54,100,100	0
2	T4K	A	1500	50/50	0.97	0.09	-0.27	13,28,67,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.