



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:32 AM GMT

PDB ID : 3ITU
Title : hPDE2A catalytic domain complexed with IBMX
Authors : Pandit, J.
Deposited on : 2009-08-28
Resolution : 1.58 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

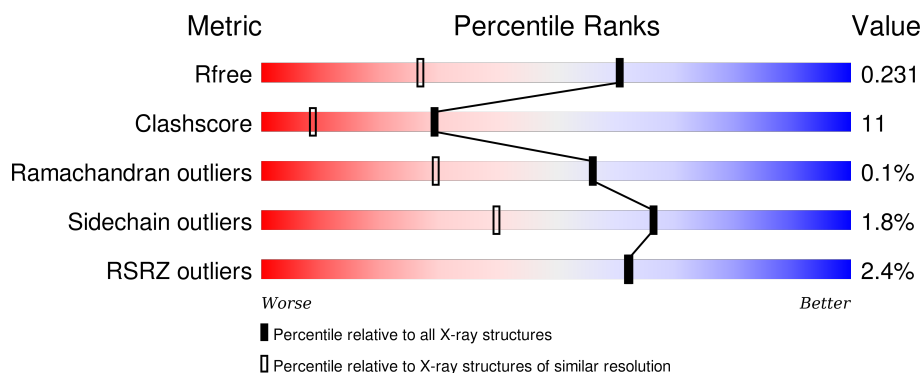
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.58 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	3815 (1.60-1.56)
Clashscore	102246	4131 (1.60-1.56)
Ramachandran outliers	100387	4021 (1.60-1.56)
Sidechain outliers	100360	4018 (1.60-1.56)
RSRZ outliers	91569	3823 (1.60-1.56)

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	345	<div> <div>2%</div> <div>82%</div> <div>14%</div> <div>• •</div> </div>
1	B	345	<div> <div>2%</div> <div>75%</div> <div>18%</div> <div>• •</div> </div>
1	C	345	<div> <div>3%</div> <div>73%</div> <div>19%</div> <div>• 5%</div> </div>
1	D	345	<div> <div>3%</div> <div>78%</div> <div>15%</div> <div>• 5%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	IBM	C	999	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12700 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 cGMP-dependent 3',5'-cyclic phosphodiesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	0	0
			2767	1763	474	505	25			
1	B	333	Total	C	N	O	S	0	0	0
			2732	1743	469	495	25			
1	C	327	Total	C	N	O	S	0	0	0
			2676	1707	460	484	25			
1	D	327	Total	C	N	O	S	0	0	0
			2676	1707	460	484	25			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	575	GLY	-	expression tag	UNP O00408
A	576	SER	-	expression tag	UNP O00408
A	577	ALA	-	expression tag	UNP O00408
A	578	MET	-	expression tag	UNP O00408
B	575	GLY	-	expression tag	UNP O00408
B	576	SER	-	expression tag	UNP O00408
B	577	ALA	-	expression tag	UNP O00408
B	578	MET	-	expression tag	UNP O00408
C	575	GLY	-	expression tag	UNP O00408
C	576	SER	-	expression tag	UNP O00408
C	577	ALA	-	expression tag	UNP O00408
C	578	MET	-	expression tag	UNP O00408
D	575	GLY	-	expression tag	UNP O00408
D	576	SER	-	expression tag	UNP O00408
D	577	ALA	-	expression tag	UNP O00408
D	578	MET	-	expression tag	UNP O00408

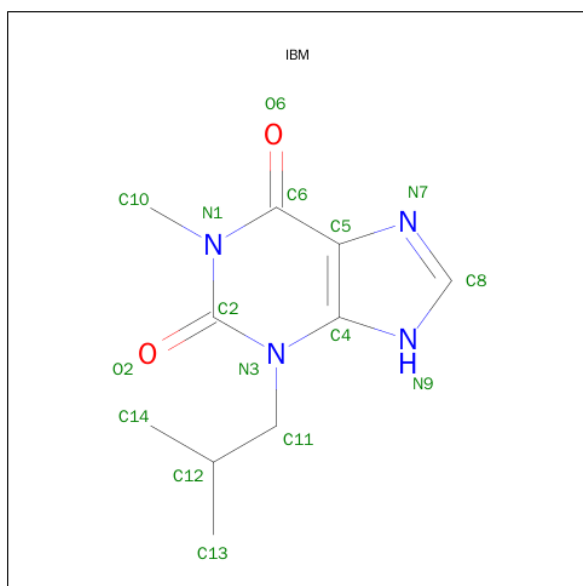
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 3-ISOBUTYL-1-METHYLNANTHINE (three-letter code: IBM) (formula: C₁₀H₁₄N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			16	10	4	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			16	10	4	2		
4	C	1	Total	C	N	O	0	0
			16	10	4	2		
4	D	1	Total	C	N	O	0	0
			16	10	4	2		

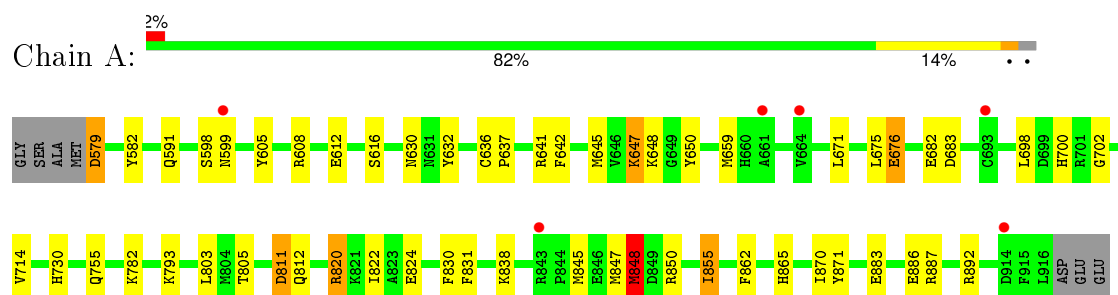
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	514	Total	O	0	0
			514	514		
5	B	447	Total	O	0	0
			447	447		
5	C	397	Total	O	0	0
			397	397		
5	D	419	Total	O	0	0
			419	419		

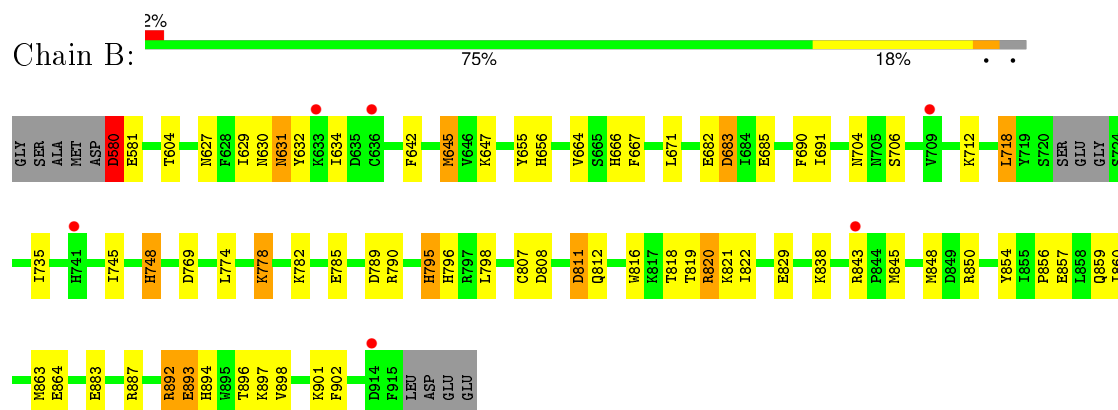
3 Residue-property plots [i](#)

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.

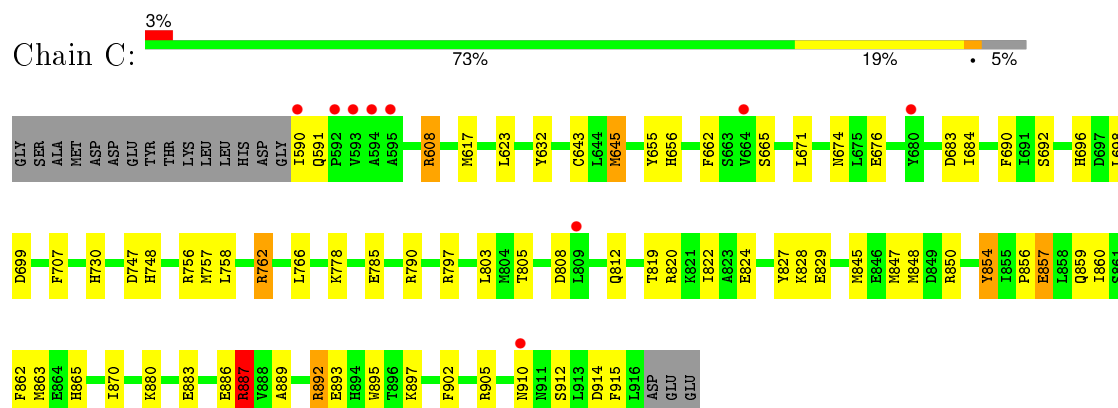
- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase



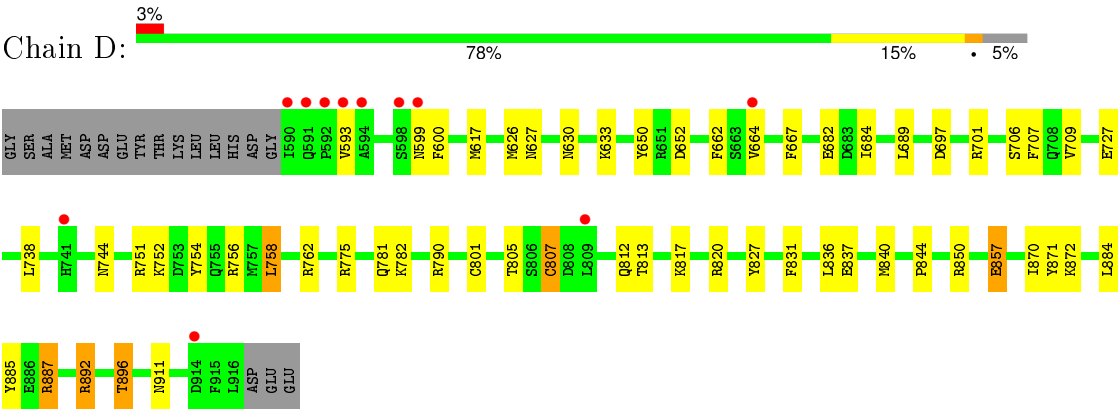
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- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase



● Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.81Å 73.29Å 91.53Å 109.30° 88.80° 89.00°	Depositor
Resolution (Å)	50.00 – 1.58 30.35 – 1.58	Depositor EDS
% Data completeness (in resolution range)	83.8 (50.00-1.58) 78.9 (30.35-1.58)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 1.58Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.174 , 0.233 0.172 , 0.231	Depositor DCC
R_{free} test set	7854 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	17.2	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.5	EDS
Estimated twinning fraction	0.015 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 157108 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12700	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.88% 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: MG, ZN, IBM

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	1.41	13/2834 (0.5%)	1.20	8/3824 (0.2%)
1	B	1.36	9/2798 (0.3%)	1.17	13/3774 (0.3%)
1	C	1.30	8/2741 (0.3%)	1.22	18/3698 (0.5%)
1	D	1.37	12/2741 (0.4%)	1.15	10/3698 (0.3%)
All	All	1.36	42/11114 (0.4%)	1.19	49/14994 (0.3%)

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	682	GLU	CB-CG	8.42	1.68	1.52
1	B	682	GLU	CG-CD	7.93	1.63	1.51
1	D	857	GLU	CG-CD	7.40	1.63	1.51
1	B	848	MET	CB-CG	7.20	1.74	1.51
1	A	650	TYR	CE2-CZ	6.69	1.47	1.38
1	C	827	TYR	CE2-CZ	6.49	1.47	1.38
1	C	662	PHE	CE2-CZ	6.43	1.49	1.37
1	B	655	TYR	CD1-CE1	6.36	1.48	1.39
1	D	727	GLU	CG-CD	6.24	1.61	1.51
1	C	854	TYR	CE2-CZ	6.14	1.46	1.38
1	A	871	TYR	CD1-CE1	6.12	1.48	1.39
1	D	754	TYR	CE2-CZ	6.12	1.46	1.38
1	A	632	TYR	CE2-CZ	6.04	1.46	1.38
1	D	885	TYR	CD1-CE1	5.87	1.48	1.39
1	C	827	TYR	CG-CD1	5.84	1.46	1.39
1	D	667	PHE	CE2-CZ	5.83	1.48	1.37
1	A	824	GLU	CG-CD	5.76	1.60	1.51
1	B	667	PHE	CG-CD1	5.72	1.47	1.38
1	A	862	PHE	CD1-CE1	5.70	1.50	1.39
1	A	582	TYR	CD2-CE2	5.61	1.47	1.39
1	D	650	TYR	CE2-CZ	5.60	1.45	1.38
1	A	871	TYR	CE1-CZ	5.58	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	676	GLU	N-CA	5.58	1.57	1.46
1	C	665	SER	CB-OG	5.57	1.49	1.42
1	C	690	PHE	CE1-CZ	5.54	1.47	1.37
1	D	697	ASP	CB-CG	5.52	1.63	1.51
1	A	632	TYR	CG-CD1	5.50	1.46	1.39
1	B	642	PHE	CD1-CE1	5.48	1.50	1.39
1	C	655	TYR	CE1-CZ	5.47	1.45	1.38
1	D	807	CYS	CA-CB	-5.42	1.42	1.53
1	A	616	SER	N-CA	5.42	1.57	1.46
1	C	690	PHE	CD2-CE2	5.36	1.50	1.39
1	D	801	CYS	CB-SG	-5.22	1.73	1.81
1	D	871	TYR	CD1-CE1	5.18	1.47	1.39
1	B	893	GLU	CG-CD	5.17	1.59	1.51
1	B	580	ASP	N-CA	5.12	1.56	1.46
1	B	690	PHE	CE1-CZ	5.10	1.47	1.37
1	D	689	LEU	N-CA	5.10	1.56	1.46
1	A	682	GLU	CD-OE1	5.10	1.31	1.25
1	A	871	TYR	CE2-CZ	5.05	1.45	1.38
1	A	642	PHE	CD2-CE2	5.05	1.49	1.39
1	D	709	VAL	CB-CG2	5.02	1.63	1.52

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	608	ARG	NE-CZ-NH2	-14.63	112.99	120.30
1	C	808	ASP	CB-CG-OD2	-11.42	108.02	118.30
1	C	608	ARG	NE-CZ-NH1	10.56	125.58	120.30
1	C	808	ASP	CB-CG-OD1	9.83	127.14	118.30
1	B	808	ASP	CB-CG-OD1	8.59	126.03	118.30
1	A	811	ASP	CB-CG-OD2	-8.27	110.85	118.30
1	A	608	ARG	NE-CZ-NH1	7.92	124.26	120.30
1	A	850	ARG	NE-CZ-NH2	-7.26	116.67	120.30
1	C	762	ARG	NE-CZ-NH2	-7.10	116.75	120.30
1	C	645	MET	CG-SD-CE	-6.73	89.43	100.20
1	C	850	ARG	NE-CZ-NH1	-6.66	116.97	120.30
1	C	766	LEU	CB-CG-CD1	-6.58	99.81	111.00
1	B	718	LEU	CA-CB-CG	6.52	130.29	115.30
1	D	775	ARG	NE-CZ-NH2	-6.50	117.05	120.30
1	C	699	ASP	CB-CG-OD2	6.47	124.12	118.30
1	B	718	LEU	CB-CG-CD1	6.19	121.52	111.00
1	A	647	LYS	CD-CE-NZ	-6.12	97.62	111.70
1	A	683	ASP	CB-CG-OD2	-6.05	112.86	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	608	ARG	CG-CD-NE	-6.01	99.17	111.80
1	C	887	ARG	NE-CZ-NH1	6.01	123.31	120.30
1	A	848	MET	CG-SD-CE	6.00	109.80	100.20
1	D	697	ASP	CB-CG-OD2	5.94	123.65	118.30
1	D	887	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	B	798	LEU	CB-CG-CD2	-5.93	100.92	111.00
1	C	905	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	B	683	ASP	CB-CG-OD2	-5.77	113.11	118.30
1	D	652	ASP	CB-CG-OD1	5.76	123.48	118.30
1	C	756	ARG	NE-CZ-NH2	-5.75	117.42	120.30
1	B	671	LEU	CB-CG-CD1	-5.74	101.24	111.00
1	B	769	ASP	CB-CG-OD1	5.70	123.43	118.30
1	B	645	MET	CG-SD-CE	-5.68	91.11	100.20
1	D	707	PHE	CB-CG-CD1	-5.68	116.82	120.80
1	B	655	TYR	CB-CG-CD1	-5.64	117.61	121.00
1	A	675	LEU	N-CA-C	5.62	126.19	111.00
1	B	671	LEU	CB-CG-CD2	5.61	120.53	111.00
1	D	626	MET	CG-SD-CE	-5.56	91.31	100.20
1	D	662	PHE	CB-CG-CD1	-5.53	116.93	120.80
1	B	811	ASP	CB-CG-OD2	-5.47	113.37	118.30
1	D	706	SER	CA-CB-OG	-5.42	96.58	111.20
1	D	758	LEU	CB-CG-CD1	-5.39	101.84	111.00
1	B	850	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	C	747	ASP	CB-CG-OD1	5.29	123.06	118.30
1	C	862	PHE	CB-CG-CD1	-5.21	117.15	120.80
1	C	707	PHE	CB-CG-CD1	-5.21	117.15	120.80
1	C	756	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	B	774	LEU	CB-CG-CD2	5.18	119.81	111.00
1	C	892	ARG	NE-CZ-NH1	5.15	122.87	120.30
1	A	676	GLU	N-CA-CB	5.07	119.72	110.60
1	D	701	ARG	NE-CZ-NH1	-5.04	117.78	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	2767	0	2709	59	0
1	B	2732	0	2679	74	0
1	C	2676	0	2630	74	0
1	D	2676	0	2630	38	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	16	0	14	3	0
4	B	16	0	14	3	0
4	C	16	0	14	1	0
4	D	16	0	14	1	0
5	A	514	0	0	36	0
5	B	447	0	0	21	2
5	C	397	0	0	30	2
5	D	419	0	0	12	0
All	All	12700	0	10704	244	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (244) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:643:CYS:HB3	5:C:1708:HOH:O	1.27	1.29
1:B:892:ARG:HH12	1:B:896:THR:CG2	1.57	1.18
1:D:892:ARG:HH12	1:D:896:THR:HG21	1.14	1.10
1:B:745:ILE:O	1:B:748:HIS:CE1	2.05	1.09
1:B:745:ILE:O	1:B:748:HIS:HE1	1.33	1.09
1:D:892:ARG:HH12	1:D:896:THR:CG2	1.66	1.08
1:A:820:ARG:HH11	1:A:820:ARG:HG2	1.24	1.02
1:A:847:MET:CE	5:A:1704:HOH:O	2.08	1.01
1:B:748:HIS:HB2	5:B:1518:HOH:O	1.62	0.98
1:B:789:ASP:H	1:B:795:HIS:HD2	1.11	0.97
1:A:676:GLU:HG3	5:A:1038:HOH:O	1.64	0.96
1:D:633:LYS:HE2	5:D:1013:HOH:O	1.66	0.95
1:B:634:ILE:HD12	1:B:748:HIS:CE1	2.01	0.95
1:B:627:ASN:HB3	5:B:1710:HOH:O	1.64	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:897:LYS:HD3	5:C:1724:HOH:O	1.68	0.93
1:B:892:ARG:NH1	1:B:896:THR:CG2	2.31	0.93
1:C:887:ARG:HH11	1:C:887:ARG:HG2	1.35	0.91
1:A:579:ASP:N	5:A:1666:HOH:O	2.05	0.89
1:C:748:HIS:HB3	5:C:1411:HOH:O	1.72	0.89
1:C:632:TYR:HE1	5:C:1712:HOH:O	1.56	0.88
1:B:894:HIS:HD2	5:B:1323:HOH:O	1.57	0.87
1:C:590:ILE:HA	1:C:617:MET:HE1	1.56	0.87
1:A:636:CYS:SG	5:A:1509:HOH:O	2.32	0.86
1:B:704:ASN:HD21	1:B:845:MET:HE1	1.38	0.86
1:D:892:ARG:NH1	1:D:896:THR:CG2	2.39	0.85
1:A:847:MET:HE3	5:A:1704:HOH:O	1.71	0.85
1:D:593:VAL:HG23	1:D:600:PHE:CD2	2.11	0.85
1:B:892:ARG:HH12	1:B:896:THR:HG23	1.39	0.85
1:A:636:CYS:HB3	5:A:1707:HOH:O	1.77	0.85
1:B:632:TYR:HB3	1:B:748:HIS:CD2	2.11	0.84
1:D:762:ARG:NH2	5:D:1738:HOH:O	2.05	0.84
1:C:914:ASP:HB3	5:C:1454:HOH:O	1.75	0.84
1:A:887:ARG:HD3	5:A:1623:HOH:O	1.78	0.83
1:A:700:HIS:HD2	1:A:702:GLY:H	1.23	0.83
1:A:847:MET:HE1	5:A:1704:HOH:O	1.72	0.83
1:B:892:ARG:NH1	1:B:896:THR:HG23	1.95	0.81
1:D:758:LEU:HB3	5:D:1738:HOH:O	1.78	0.81
1:B:812:GLN:HE21	4:B:999:IBM:H101	1.46	0.80
1:A:845:MET:HB2	1:A:848:MET:HG2	1.64	0.80
1:B:656:HIS:HD2	1:B:829:GLU:OE2	1.66	0.79
1:C:886:GLU:HG2	5:C:975:HOH:O	1.82	0.79
1:C:897:LYS:CD	5:C:1724:HOH:O	2.26	0.78
1:D:857:GLU:HG3	5:D:1669:HOH:O	1.82	0.78
1:C:656:HIS:HD2	1:C:829:GLU:OE2	1.67	0.78
1:A:883:GLU:OE1	1:A:887:ARG:NH2	2.17	0.78
1:B:685:GLU:OE1	1:B:796:HIS:HD2	1.67	0.78
1:C:617:MET:HG3	5:C:189:HOH:O	1.84	0.77
1:D:593:VAL:CG2	1:D:600:PHE:CD2	2.68	0.76
1:D:782:LYS:HE2	5:D:982:HOH:O	1.86	0.75
1:D:857:GLU:CG	5:D:1669:HOH:O	2.35	0.74
1:C:683:ASP:HB3	5:C:1712:HOH:O	1.87	0.74
1:D:627:ASN:HD21	1:D:630:ASN:HD22	1.34	0.74
1:C:845:MET:HE3	1:C:847:MET:SD	2.27	0.74
1:D:892:ARG:NH1	1:D:896:THR:HG21	1.96	0.74
1:B:789:ASP:H	1:B:795:HIS:CD2	2.02	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:704:ASN:HD21	1:B:845:MET:CE	2.01	0.73
1:B:704:ASN:ND2	1:B:845:MET:HE1	2.04	0.73
1:A:865:HIS:CE1	5:A:936:HOH:O	2.41	0.72
1:D:813:THR:O	1:D:887:ARG:HD3	1.90	0.72
1:C:845:MET:HE2	1:C:848:MET:HG3	1.72	0.72
1:C:698:LEU:O	1:C:730:HIS:HD2	1.74	0.70
1:C:887:ARG:HD3	5:C:304:HOH:O	1.91	0.70
1:B:790:ARG:O	1:B:796:HIS:HE1	1.73	0.69
5:A:1560:HOH:O	1:B:712:LYS:HB2	1.92	0.68
1:A:630:ASN:ND2	5:A:1000:HOH:O	2.25	0.68
1:C:676:GLU:HG2	5:C:1364:HOH:O	1.93	0.68
1:B:632:TYR:HB2	1:B:634:ILE:HD11	1.76	0.68
1:B:631:ASN:ND2	1:B:683:ASP:OD2	2.26	0.68
1:A:820:ARG:NH1	1:A:820:ARG:HG2	1.97	0.67
1:C:845:MET:CE	1:C:847:MET:CG	2.72	0.67
1:C:632:TYR:CE1	5:C:1712:HOH:O	2.37	0.67
1:A:700:HIS:CD2	1:A:702:GLY:H	2.09	0.67
1:A:730:HIS:HE1	5:A:5:HOH:O	1.79	0.66
1:C:591:GLN:H	1:C:617:MET:CE	2.09	0.66
1:B:812:GLN:NE2	4:B:999:IBM:H101	2.11	0.65
1:C:845:MET:HE3	1:C:847:MET:CG	2.25	0.65
1:C:730:HIS:HE1	5:C:17:HOH:O	1.79	0.65
1:D:751:ARG:HD2	5:D:1022:HOH:O	1.97	0.65
1:B:887:ARG:CD	5:B:1018:HOH:O	2.44	0.65
1:B:892:ARG:NH1	1:B:896:THR:HG21	2.12	0.64
1:C:915:PHE:HB3	5:C:1750:HOH:O	1.97	0.63
1:C:819:THR:CG2	1:C:895:TRP:HE1	2.12	0.63
1:B:897:LYS:HD2	5:B:1189:HOH:O	1.99	0.63
1:B:632:TYR:HB2	1:B:634:ILE:CD1	2.28	0.63
1:A:636:CYS:CB	5:A:1707:HOH:O	2.41	0.62
1:B:887:ARG:HG2	5:B:550:HOH:O	1.98	0.62
1:C:591:GLN:H	1:C:617:MET:HE2	1.65	0.61
1:A:648:LYS:HG3	1:D:790:ARG:HD2	1.82	0.61
1:B:887:ARG:HD3	5:B:1018:HOH:O	1.98	0.61
1:D:593:VAL:CG2	1:D:600:PHE:HD2	2.13	0.61
1:A:698:LEU:O	1:A:730:HIS:HD2	1.82	0.60
1:B:893:GLU:CG	5:B:1381:HOH:O	2.49	0.60
1:D:892:ARG:HG2	1:D:892:ARG:HH11	1.66	0.60
1:B:893:GLU:HG2	5:B:1381:HOH:O	2.01	0.60
1:C:887:ARG:CD	5:C:304:HOH:O	2.47	0.59
1:C:683:ASP:CB	5:C:1712:HOH:O	2.47	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:630:ASN:O	1:B:632:TYR:N	2.36	0.58
1:A:591:GLN:HG3	5:A:1584:HOH:O	2.01	0.58
1:A:865:HIS:HE1	5:A:936:HOH:O	1.80	0.57
1:C:865:HIS:HD2	5:C:1269:HOH:O	1.85	0.57
1:C:812:GLN:HE21	4:C:999:IBM:H101	1.68	0.57
1:D:812:GLN:HE21	4:D:999:IBM:H101	1.69	0.57
1:A:598:SER:HB2	5:A:1660:HOH:O	2.05	0.57
1:D:682:GLU:OE1	1:D:756:ARG:NH2	2.37	0.57
1:B:859:GLN:OE1	4:B:999:IBM:N7	2.37	0.57
1:B:630:ASN:O	1:B:631:ASN:C	2.43	0.56
1:A:579:ASP:OD2	5:A:1678:HOH:O	2.17	0.56
1:C:845:MET:HE2	1:C:847:MET:HG3	1.86	0.56
1:C:824:GLU:HG2	1:C:828:LYS:HE3	1.86	0.56
1:A:812:GLN:HE21	4:A:999:IBM:H101	1.69	0.56
1:D:884:LEU:HA	1:D:887:ARG:HD2	1.87	0.56
1:A:830:PHE:HE2	1:A:847:MET:CE	2.19	0.56
1:B:685:GLU:OE1	1:B:796:HIS:CD2	2.55	0.56
1:B:666:HIS:HD2	5:B:17:HOH:O	1.89	0.56
1:A:630:ASN:HB3	5:A:1499:HOH:O	2.06	0.55
1:A:820:ARG:HH11	1:A:820:ARG:CG	2.09	0.55
1:A:805:THR:HG22	1:A:870:ILE:HD13	1.88	0.55
1:C:887:ARG:NH1	1:C:887:ARG:HG2	2.07	0.55
1:C:674:ASN:O	1:C:880:LYS:CE	2.55	0.55
1:B:785:GLU:HG3	5:B:294:HOH:O	2.07	0.55
1:B:789:ASP:N	1:B:795:HIS:HD2	1.94	0.54
1:C:819:THR:HG21	1:C:895:TRP:HE1	1.71	0.54
1:B:856:PRO:HG3	1:B:902:PHE:CD2	2.43	0.54
5:A:1560:HOH:O	1:B:712:LYS:CB	2.52	0.54
1:B:631:ASN:ND2	5:B:1710:HOH:O	2.41	0.54
1:B:634:ILE:CD1	1:B:748:HIS:CE1	2.86	0.53
1:C:845:MET:HE2	1:C:847:MET:CG	2.38	0.53
1:C:887:ARG:CZ	5:C:1632:HOH:O	2.56	0.53
1:B:782:LYS:HG2	5:B:1739:HOH:O	2.08	0.53
1:C:785:GLU:HB2	5:C:1701:HOH:O	2.09	0.53
1:C:790:ARG:O	1:C:790:ARG:HG2	2.08	0.52
1:B:819:THR:HG21	1:B:863:MET:HE3	1.91	0.52
1:B:883:GLU:OE1	1:B:887:ARG:NH1	2.36	0.52
1:C:676:GLU:HA	5:C:959:HOH:O	2.10	0.52
1:B:821:LYS:HE2	5:B:1277:HOH:O	2.09	0.51
1:D:627:ASN:ND2	1:D:630:ASN:HD22	2.06	0.51
1:A:831:PHE:HZ	1:A:855:ILE:HD11	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:656:HIS:HE1	5:B:12:HOH:O	1.92	0.51
1:A:579:ASP:CA	5:A:1666:HOH:O	2.56	0.51
1:B:627:ASN:HA	5:B:1717:HOH:O	2.11	0.51
1:C:797:ARG:HG2	5:C:976:HOH:O	2.11	0.50
1:A:805:THR:CG2	1:A:870:ILE:HD13	2.42	0.50
1:A:636:CYS:HB2	1:A:637:PRO:HD3	1.94	0.50
1:A:630:ASN:CB	5:A:544:HOH:O	2.59	0.50
1:C:910:ASN:ND2	5:C:1620:HOH:O	2.44	0.50
1:A:782:LYS:HD3	5:A:972:HOH:O	2.12	0.50
1:D:752:LYS:HG3	5:D:954:HOH:O	2.11	0.49
1:A:714:VAL:HG12	5:A:1092:HOH:O	2.12	0.49
1:A:838:LYS:HD2	5:A:1718:HOH:O	2.12	0.49
1:A:630:ASN:HB3	5:A:544:HOH:O	2.10	0.49
1:D:781:GLN:HG2	5:D:1089:HOH:O	2.11	0.49
1:B:860:ILE:O	1:B:864:GLU:HG3	2.13	0.49
1:B:845:MET:HE2	5:B:1719:HOH:O	2.12	0.49
1:A:641:ARG:HD3	5:A:1743:HOH:O	2.13	0.49
1:B:580:ASP:N	5:B:1716:HOH:O	2.45	0.49
1:B:630:ASN:C	1:B:632:TYR:N	2.66	0.49
1:A:830:PHE:CE2	1:A:847:MET:CE	2.96	0.49
1:C:845:MET:CE	1:C:847:MET:HG2	2.42	0.48
1:C:797:ARG:HB3	5:C:1741:HOH:O	2.12	0.48
1:C:656:HIS:HE1	5:C:18:HOH:O	1.96	0.48
1:A:630:ASN:ND2	5:A:475:HOH:O	2.33	0.48
1:B:581:GLU:OE2	1:B:647:LYS:NZ	2.36	0.48
1:A:886:GLU:OE1	5:A:1559:HOH:O	2.20	0.48
1:B:864:GLU:OE2	1:B:892:ARG:NH2	2.44	0.48
1:C:910:ASN:OD1	1:C:912:SER:HB3	2.14	0.48
1:B:854:TYR:CD2	1:B:857:GLU:HG3	2.49	0.48
1:A:830:PHE:CE2	1:A:847:MET:HE2	2.49	0.47
1:C:819:THR:HG22	1:C:895:TRP:HE1	1.79	0.47
1:D:837:GLU:OE1	1:D:844:PRO:HB3	2.15	0.47
1:C:805:THR:HG22	1:C:870:ILE:HD13	1.96	0.47
1:D:872:LYS:HD3	5:D:967:HOH:O	2.15	0.47
1:A:793:LYS:HB2	1:A:793:LYS:HE3	1.67	0.47
1:B:898:VAL:HB	1:B:901:LYS:HD2	1.96	0.47
1:A:887:ARG:HG2	5:A:1057:HOH:O	2.14	0.47
1:D:827:TYR:HB3	1:D:831:PHE:CE2	2.50	0.47
1:D:805:THR:HG22	1:D:870:ILE:HD13	1.97	0.46
1:C:698:LEU:O	1:C:730:HIS:CD2	2.61	0.46
1:C:591:GLN:H	1:C:617:MET:HE1	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:854:TYR:HD2	1:C:857:GLU:HB3	1.80	0.46
1:A:811:ASP:HB3	1:A:822:ILE:HG13	1.97	0.46
1:A:730:HIS:CE1	5:A:5:HOH:O	2.61	0.46
1:C:608:ARG:CZ	1:C:822:ILE:HD11	2.45	0.46
1:B:778:LYS:HE3	1:B:778:LYS:HB3	1.62	0.46
1:A:599:ASN:O	1:A:605:TYR:HB2	2.15	0.46
1:C:684:ILE:HD11	1:C:757:MET:SD	2.55	0.46
1:C:860:ILE:HG23	1:C:892:ARG:HH11	1.81	0.46
1:A:782:LYS:HE2	5:A:1504:HOH:O	2.14	0.45
1:C:692:SER:O	1:C:696:HIS:HB3	2.16	0.45
1:A:847:MET:HB2	1:A:847:MET:HE2	1.70	0.45
1:D:836:LEU:O	1:D:840:MET:HG3	2.15	0.45
1:C:671:LEU:HD13	1:C:803:LEU:HD22	1.99	0.45
1:C:897:LYS:HD2	5:C:1724:HOH:O	2.05	0.45
1:B:735:ILE:HA	1:B:735:ILE:HD13	1.62	0.45
1:A:700:HIS:HE1	5:A:143:HOH:O	2.00	0.45
1:B:664:VAL:HG13	1:B:807:CYS:HB3	1.99	0.45
1:C:859:GLN:O	1:C:863:MET:HG3	2.16	0.45
1:A:647:LYS:CD	5:A:968:HOH:O	2.64	0.45
1:B:645:MET:HG2	5:B:388:HOH:O	2.16	0.44
1:A:755:GLN:CG	5:A:1560:HOH:O	2.65	0.44
1:B:854:TYR:HD2	1:B:857:GLU:HG3	1.82	0.44
1:C:889:ALA:O	1:C:893:GLU:HG2	2.17	0.44
4:A:999:IBM:H8	5:A:1704:HOH:O	2.16	0.44
1:D:664:VAL:HG13	1:D:807:CYS:HB3	2.00	0.44
1:C:645:MET:HG2	5:C:259:HOH:O	2.18	0.44
1:D:892:ARG:HG2	1:D:892:ARG:NH1	2.33	0.44
1:A:847:MET:HE1	4:A:999:IBM:H8	1.99	0.44
1:C:762:ARG:NH1	5:C:478:HOH:O	2.50	0.44
1:D:617:MET:HE3	5:D:1214:HOH:O	2.18	0.44
1:C:590:ILE:HG23	1:C:617:MET:CE	2.48	0.44
1:C:758:LEU:HD23	1:C:758:LEU:HA	1.80	0.44
1:C:856:PRO:HG3	1:C:902:PHE:CD2	2.53	0.43
1:A:645:MET:HG2	5:A:1339:HOH:O	2.17	0.43
1:B:893:GLU:HA	1:B:893:GLU:OE1	2.18	0.43
1:C:856:PRO:HG3	1:C:902:PHE:CE2	2.53	0.43
1:B:811:ASP:HB3	1:B:822:ILE:HG13	1.99	0.43
1:B:818:THR:O	1:B:822:ILE:HG12	2.18	0.43
1:A:830:PHE:HE2	1:A:847:MET:HE1	1.82	0.43
1:C:883:GLU:O	1:C:887:ARG:NH1	2.52	0.43
1:A:648:LYS:HG3	1:D:790:ARG:CD	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:816:TRP:CZ2	1:B:820:ARG:HG3	2.53	0.43
1:C:883:GLU:O	1:C:887:ARG:HG2	2.18	0.43
1:B:892:ARG:HH12	1:B:896:THR:HG22	1.65	0.42
1:D:850:ARG:HD3	1:D:850:ARG:HH11	1.69	0.42
1:B:630:ASN:ND2	5:B:235:HOH:O	2.52	0.42
1:B:604:THR:HG22	1:B:887:ARG:NH2	2.35	0.42
1:D:817:LYS:N	1:D:817:LYS:HD2	2.34	0.42
1:B:892:ARG:HH12	1:B:896:THR:HG21	1.59	0.42
1:C:859:GLN:NE2	5:C:1720:HOH:O	2.52	0.42
1:C:590:ILE:CA	1:C:617:MET:HE1	2.39	0.42
1:B:893:GLU:HG3	5:B:1381:HOH:O	2.14	0.41
1:D:756:ARG:HD2	5:D:1001:HOH:O	2.20	0.41
1:C:886:GLU:CG	5:C:975:HOH:O	2.55	0.41
1:C:845:MET:HE2	1:C:848:MET:CG	2.46	0.41
1:D:911:ASN:HA	1:D:911:ASN:HD22	1.63	0.41
1:A:612:GLU:HG3	1:A:659:MET:SD	2.60	0.41
1:C:778:LYS:HD3	5:C:979:HOH:O	2.21	0.41
1:B:629:ILE:HA	1:B:634:ILE:HG12	2.03	0.41
1:C:623:LEU:HD23	1:C:623:LEU:HA	1.91	0.41
1:B:838:LYS:HE2	1:B:843:ARG:NH2	2.36	0.40
1:C:845:MET:HE3	1:C:847:MET:HG2	2.02	0.40
1:D:738:LEU:O	1:D:744:ASN:HB2	2.21	0.40
1:A:671:LEU:HD13	1:A:803:LEU:HD22	2.04	0.40
1:B:691:ILE:HA	1:B:691:ILE:HD13	1.94	0.40
1:C:757:MET:HB3	1:C:757:MET:HE2	1.95	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:311:HOH:O	5:C:924:HOH:O[1_545]	2.00	0.20
5:B:311:HOH:O	5:C:1618:HOH:O[1_545]	2.19	0.01

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	336/345 (97%)	332 (99%)	4 (1%)	0	100	100
1	B	329/345 (95%)	318 (97%)	10 (3%)	1 (0%)	46	22
1	C	325/345 (94%)	320 (98%)	5 (2%)	0	100	100
1	D	325/345 (94%)	320 (98%)	5 (2%)	0	100	100
All	All	1315/1380 (95%)	1290 (98%)	24 (2%)	1 (0%)	56	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	631	ASN

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	305/310 (98%)	300 (98%)	5 (2%)	70	45
1	B	301/310 (97%)	293 (97%)	8 (3%)	52	22
1	C	295/310 (95%)	292 (99%)	3 (1%)	82	65
1	D	295/310 (95%)	290 (98%)	5 (2%)	68	42
All	All	1196/1240 (96%)	1175 (98%)	21 (2%)	66	40

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	579	ASP
1	A	820	ARG
1	A	848	MET
1	A	855	ILE
1	A	892	ARG
1	B	580	ASP
1	B	706	SER

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Mol	Chain	Res	Type
1	B	718	LEU
1	B	748	HIS
1	B	778	LYS
1	B	795	HIS
1	B	820	ARG
1	B	892	ARG
1	C	820	ARG
1	C	857	GLU
1	C	887	ARG
1	D	599	ASN
1	D	684	ILE
1	D	820	ARG
1	D	892	ARG
1	D	896	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	700	HIS
1	A	708	GLN
1	A	730	HIS
1	A	859	GLN
1	A	911	ASN
1	B	656	HIS
1	B	666	HIS
1	B	674	ASN
1	B	748	HIS
1	B	795	HIS
1	B	796	HIS
1	B	842	ASN
1	B	859	GLN
1	B	900	HIS
1	B	911	ASN
1	C	656	HIS
1	C	730	HIS
1	D	627	ASN
1	D	791	ASN
1	D	859	GLN
1	D	900	HIS
1	D	911	ASN

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 ⓘ

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 for Mogul analysis.

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	IBM	A	999	-	9,17,17	1.27	1 (11%)	9,25,25	2.33	1 (11%)
4	IBM	B	999	-	9,17,17	0.96	0	9,25,25	1.18	1 (11%)
4	IBM	C	999	-	9,17,17	2.07	1 (11%)	9,25,25	0.91	1 (11%)
4	IBM	D	999	-	9,17,17	1.92	3 (33%)	9,25,25	2.14	3 (33%)

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
4	IBM	A	999	-	-	0/4/4/4	0/2/2/2
4	IBM	B	999	-	-	0/4/4/4	0/2/2/2
4	IBM	C	999	-	-	0/4/4/4	0/2/2/2
4	IBM	D	999	-	-	0/4/4/4	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	999	IBM	C11-N3	-5.67	1.42	1.48
4	D	999	IBM	C4-N3	-4.34	1.34	1.39
4	A	999	IBM	C4-N3	-2.29	1.36	1.39
4	D	999	IBM	O6-C6	2.31	1.30	1.24
4	D	999	IBM	C6-N1	2.34	1.41	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	999	IBM	C6-C5-C4	-6.22	115.47	119.93
4	D	999	IBM	C4-C5-N7	-4.29	105.53	109.48
4	B	999	IBM	C4-C5-N7	-2.39	107.28	109.48
4	C	999	IBM	C6-C5-C4	-2.02	118.48	119.93
4	D	999	IBM	C10-N1-C6	2.16	124.99	118.06
4	D	999	IBM	C5-C4-N9	3.09	116.81	110.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	999	IBM	3	0
4	B	999	IBM	3	0
4	C	999	IBM	1	0
4	D	999	IBM	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	338/345 (97%)	-0.01	6 (1%) 71 72	8, 17, 31, 43	0
1	B	333/345 (96%)	0.05	6 (1%) 71 72	10, 20, 32, 44	0
1	C	327/345 (94%)	0.13	9 (2%) 56 56	11, 21, 36, 44	0
1	D	327/345 (94%)	0.16	11 (3%) 49 48	9, 20, 35, 47	0
All	All	1325/1380 (96%)	0.08	32 (2%) 62 62	8, 19, 34, 47	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	593	VAL	4.9
1	D	599	ASN	3.9
1	D	590	ILE	3.7
1	D	594	ALA	3.6
1	D	914	ASP	3.5
1	B	843	ARG	3.1
1	C	593	VAL	3.1
1	D	741	HIS	3.1
1	C	664	VAL	3.1
1	A	843	ARG	3.0
1	B	914	ASP	2.9
1	C	595	ALA	2.9
1	D	592	PRO	2.8
1	C	680	TYR	2.8
1	D	809	LEU	2.7
1	D	591	GLN	2.7
1	A	664	VAL	2.7
1	C	594	ALA	2.5
1	B	741	HIS	2.5
1	C	809	LEU	2.4
1	B	709	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	910	ASN	2.2
1	D	664	VAL	2.2
1	D	598	SER	2.2
1	A	599	ASN	2.2
1	A	693	CYS	2.2
1	C	590	ILE	2.1
1	A	661	ALA	2.1
1	B	633	LYS	2.1
1	A	914	ASP	2.1
1	C	592	PRO	2.1
1	B	636	CYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	IBM	C	999	16/16	0.93	0.16	2.09	25,28,30,31	0
4	IBM	D	999	16/16	0.95	0.14	1.10	18,22,25,29	0
3	MG	D	920	1/1	1.00	0.14	1.10	12,12,12,12	0
4	IBM	B	999	16/16	0.91	0.12	0.61	23,25,28,30	0
4	IBM	A	999	16/16	0.94	0.11	0.34	17,21,22,24	0
3	MG	C	920	1/1	0.99	0.11	-0.11	12,12,12,12	0
3	MG	A	920	1/1	1.00	0.09	-0.72	10,10,10,10	0
3	MG	B	920	1/1	1.00	0.05	-2.24	10,10,10,10	0
2	ZN	C	3	1/1	1.00	0.06	-2.85	14,14,14,14	0
2	ZN	A	1	1/1	1.00	0.06	-2.99	13,13,13,13	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ZN	D	4	1/1	1.00	0.07	-3.88	14,14,14,14	0
2	ZN	B	2	1/1	1.00	0.06	-5.08	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.