



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:52 PM UTC

PDB ID : 3ECN / pdb_00003ecn
Title : Crystal structure of PDE8A catalytic domain in complex with IBMX
Authors : Wang, H.; Yan, Z.; Yang, S.; Cai, J.; Robinson, H.; Ke, H.
Deposited on : 2008-09-01
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

<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

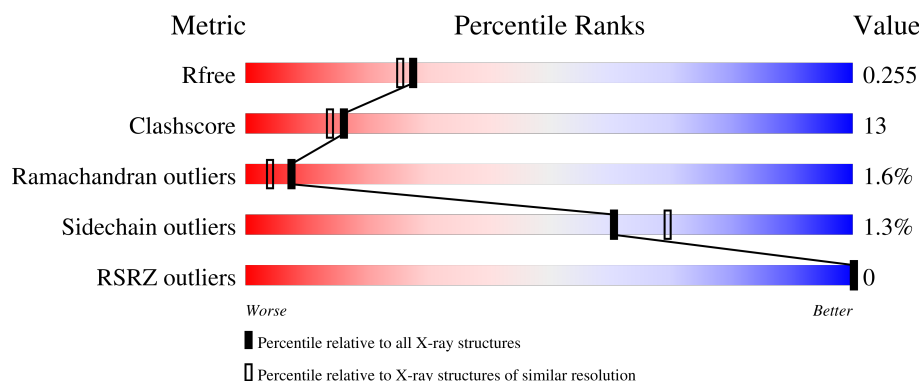
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (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 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	338	 72% 25% •
1	B	338	 73% 24% •

2 Entry composition [i](#)

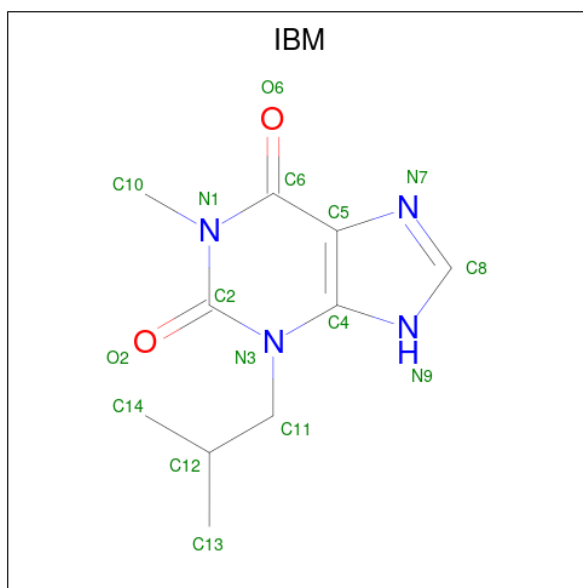
There are 5 unique types of molecules in this entry. The entry contains 5658 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 High affinity cAMP-specific and IBMX-insensitive 3',5'-cyclic phosphodiesterase 8A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	0	0
			2744	1734	464	527	19			
1	B	338	Total	C	N	O	S	0	0	0
			2744	1734	464	527	19			

- Molecule 2 is 3-ISOBUTYL-1-METHYLNANTHINE (CCD ID: IBM) (formula: C₁₀H₁₄N₄O₂).



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

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Zn 1	0	0
3	B	1	Total 1	Zn 1	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Mg 1	0	0
4	B	1	Total 1	Mg 1	0	0

- Molecule 5 is water.

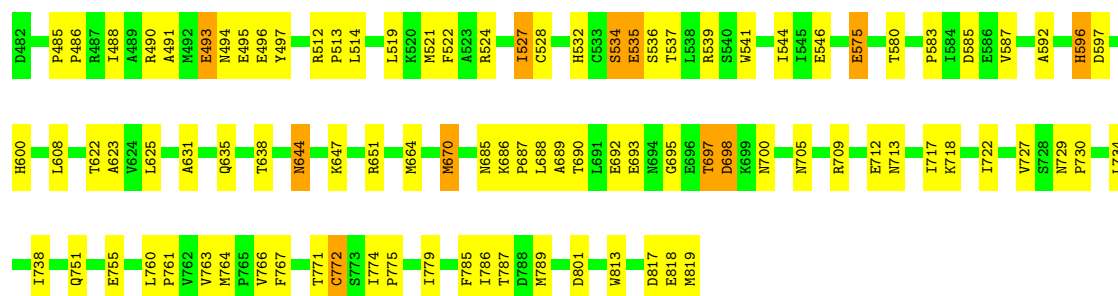
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	65	Total 65	O 65	0	0
5	B	69	Total 69	O 69	0	0

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.

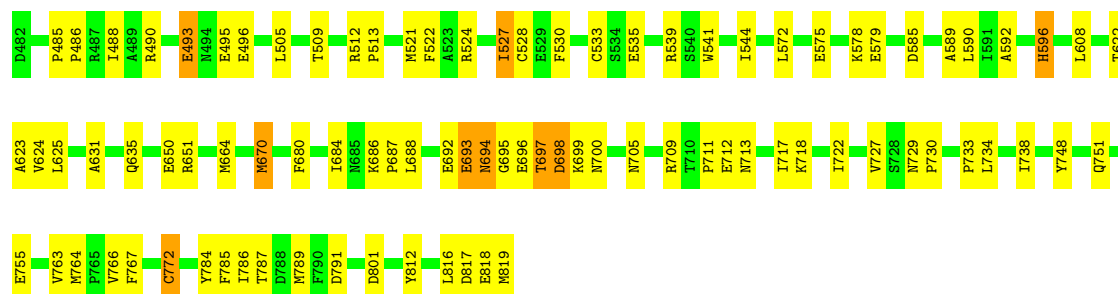
- Molecule 1: High affinity cAMP-specific and IBMX-insensitive 3',5'-cyclic phosphodiesterase 8A

Chain A: 



- Molecule 1: High affinity cAMP-specific and IBMX-insensitive 3',5'-cyclic phosphodiesterase 8A

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.33Å 101.60Å 76.57Å 90.00° 119.90° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 30.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	88.1 (30.00-2.10) 88.0 (30.00-2.10)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.10Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.221 , 0.256 0.221 , 0.255	Depositor DCC
R_{free} test set	5469 reflections (10.17%)	wwPDB-VP
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.790	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for l,k,-h-l 0.014 for -h-l,k,h 0.487 for h,-k,-h-l 0.016 for l,-k,h 0.017 for -h-l,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5658	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% 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: IBM, MG, ZN

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	0.41	0/2809	0.87	8/3811 (0.2%)
1	B	0.42	0/2809	0.88	8/3811 (0.2%)
All	All	0.41	0/5618	0.87	16/7622 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	786	ILE	N-CA-C	10.48	120.47	110.30
1	B	786	ILE	N-CA-C	10.29	120.28	110.30
1	B	624	VAL	N-CA-C	8.36	118.92	110.23
1	A	596	HIS	N-CA-C	7.11	121.58	112.34
1	B	772	CYS	N-CA-C	6.94	120.38	109.96
1	A	670	MET	N-CA-C	6.83	119.56	111.71
1	B	670	MET	N-CA-C	6.76	119.48	111.71
1	A	535	GLU	N-CA-C	-6.73	104.79	112.87
1	A	772	CYS	N-CA-C	6.31	119.79	109.76
1	A	685	ASN	N-CA-C	6.20	117.71	111.07
1	B	527	ILE	N-CA-C	5.95	116.12	110.53
1	B	693	GLU	N-CA-C	-5.80	105.30	112.90
1	B	493	GLU	N-CA-C	5.58	117.04	111.07
1	A	787	THR	N-CA-C	5.07	117.20	111.02
1	B	787	THR	N-CA-C	5.06	117.19	111.02
1	A	527	ILE	N-CA-C	5.01	115.78	110.72

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	2744	0	2631	69	0
1	B	2744	0	2631	76	0
2	A	16	0	14	3	0
2	B	16	0	14	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	65	0	0	1	0
5	B	69	0	0	1	0
All	All	5658	0	5290	142	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 13.

All (142) 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:A:493:GLU:HG3	1:A:494:ASN:H	1.25	1.02
1:A:575:GLU:H	1:A:575:GLU:CD	1.83	0.87
1:B:692:GLU:HA	1:B:696:GLU:HG2	1.54	0.85
1:B:696:GLU:HA	1:B:700:ASN:HD22	1.39	0.85
1:B:486:PRO:HB2	1:B:490:ARG:HH21	1.43	0.83
1:B:496:GLU:HG2	1:B:524:ARG:NH1	1.94	0.81
1:A:496:GLU:HG2	1:A:524:ARG:NH1	1.99	0.76
1:B:695:GLY:C	1:B:697:THR:H	1.95	0.75
1:A:763:VAL:HG23	1:A:764:MET:HG3	1.69	0.74
1:A:493:GLU:HG3	1:A:494:ASN:N	2.01	0.74
1:A:493:GLU:CG	1:A:494:ASN:H	1.98	0.72
1:B:696:GLU:HA	1:B:700:ASN:ND2	2.05	0.71
1:B:575:GLU:O	1:B:579:GLU:HG2	1.91	0.70
1:B:585:ASP:HB3	1:B:718:LYS:HD2	1.76	0.68
1:A:697:THR:HG22	1:A:698:ASP:H	1.60	0.67
1:A:585:ASP:HB3	1:A:718:LYS:HD2	1.77	0.66
1:A:535:GLU:O	1:A:539:ARG:HG3	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:688:LEU:O	1:B:692:GLU:HG3	1.97	0.64
1:B:692:GLU:CA	1:B:696:GLU:HG2	2.27	0.64
1:A:779:ILE:HD11	1:A:813:TRP:HB2	1.81	0.63
1:A:608:LEU:HD11	1:A:625:LEU:HD21	1.80	0.63
1:A:760:LEU:HD23	1:B:651:ARG:HH22	1.64	0.63
1:B:670:MET:HG3	2:B:3:IBM:H142	1.81	0.62
1:B:705:ASN:HD22	1:B:709:ARG:NH2	1.98	0.62
1:A:751:GLN:O	1:A:755:GLU:HG3	2.00	0.61
1:B:485:PRO:HG2	1:B:488:ILE:HD12	1.81	0.61
1:A:485:PRO:HG2	1:A:488:ILE:HD12	1.82	0.60
1:A:689:ALA:O	1:A:693:GLU:HG3	2.03	0.59
1:A:766:VAL:O	1:A:772:CYS:HB2	2.03	0.59
1:B:686:LYS:HB2	1:B:687:PRO:HD3	1.85	0.59
1:B:763:VAL:HG23	1:B:764:MET:HG3	1.85	0.58
1:B:521:MET:HA	1:B:521:MET:HE2	1.85	0.58
1:A:496:GLU:HG2	1:A:524:ARG:HH11	1.70	0.57
1:A:760:LEU:HD23	1:B:651:ARG:NH2	2.20	0.57
1:A:486:PRO:HB2	1:A:490:ARG:NH2	2.20	0.56
1:B:670:MET:HG3	2:B:3:IBM:C14	2.34	0.56
1:A:644:ASN:HD22	1:A:647:LYS:HB2	1.69	0.56
1:A:693:GLU:C	1:A:695:GLY:H	2.13	0.56
1:A:817:ASP:C	1:A:819:MET:H	2.13	0.56
1:B:693:GLU:C	1:B:695:GLY:H	2.14	0.56
1:B:698:ASP:C	1:B:700:ASN:H	2.13	0.56
1:B:801:ASP:HB2	5:B:35:HOH:O	2.06	0.56
1:A:534:SER:HB2	1:A:537:THR:H	1.71	0.55
1:A:521:MET:HE2	1:A:521:MET:HA	1.86	0.55
1:A:670:MET:HG3	2:A:3:IBM:H142	1.88	0.55
1:A:686:LYS:HB2	1:A:687:PRO:HD3	1.89	0.55
1:A:631:ALA:O	1:A:635:GLN:HG3	2.08	0.54
1:A:494:ASN:HB3	1:A:497:TYR:CD1	2.43	0.54
1:A:690:THR:HA	1:A:693:GLU:OE2	2.07	0.54
1:B:694:ASN:C	1:B:696:GLU:H	2.17	0.53
1:B:535:GLU:O	1:B:539:ARG:HG3	2.09	0.53
1:A:528:CYS:O	1:A:532:HIS:N	2.41	0.52
1:A:705:ASN:HD22	1:A:709:ARG:NH2	2.07	0.52
1:A:534:SER:C	1:A:536:SER:N	2.67	0.52
1:B:512:ARG:N	1:B:513:PRO:CD	2.73	0.52
1:A:764:MET:HE2	1:A:767:PHE:CD1	2.46	0.51
1:B:522:PHE:CG	1:B:527:ILE:HD12	2.46	0.51
1:B:729:ASN:HB2	1:B:730:PRO:HD3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:785:PHE:CE1	2:A:3:IBM:H131	2.45	0.51
1:A:764:MET:HB3	1:A:767:PHE:HD1	1.76	0.51
1:A:801:ASP:HB2	5:A:835:HOH:O	2.10	0.51
1:B:496:GLU:HG2	1:B:524:ARG:HH11	1.73	0.51
1:A:519:LEU:HD21	1:A:539:ARG:HG2	1.94	0.50
1:B:631:ALA:O	1:B:635:GLN:HG3	2.12	0.50
1:A:491:ALA:C	1:A:493:GLU:H	2.19	0.49
1:A:688:LEU:O	1:A:692:GLU:HG3	2.12	0.49
1:B:817:ASP:O	1:B:819:MET:HG3	2.12	0.49
1:A:512:ARG:N	1:A:513:PRO:CD	2.76	0.49
1:B:817:ASP:C	1:B:819:MET:H	2.20	0.49
1:B:521:MET:HE1	1:B:524:ARG:NH1	2.27	0.49
1:B:764:MET:HE2	1:B:767:PHE:CD1	2.48	0.48
1:A:522:PHE:CG	1:A:527:ILE:HD12	2.48	0.48
1:B:698:ASP:C	1:B:700:ASN:N	2.71	0.48
1:A:729:ASN:HB2	1:A:730:PRO:HD3	1.95	0.48
1:B:680:PHE:CD1	1:B:684:ILE:HD12	2.48	0.48
1:B:784:TYR:HD1	1:B:785:PHE:CE1	2.32	0.48
1:A:755:GLU:OE2	1:A:763:VAL:HG22	2.14	0.48
1:B:541:TRP:O	1:B:544:ILE:HG22	2.13	0.48
1:B:751:GLN:O	1:B:755:GLU:HG3	2.14	0.48
1:B:528:CYS:C	1:B:530:PHE:H	2.22	0.47
1:A:638:THR:HG22	1:A:644:ASN:OD1	2.14	0.47
1:A:734:LEU:O	1:A:738:ILE:HG13	2.14	0.47
1:A:486:PRO:O	1:A:490:ARG:HG3	2.14	0.47
1:A:695:GLY:O	1:A:700:ASN:ND2	2.48	0.47
1:B:734:LEU:O	1:B:738:ILE:HG13	2.15	0.47
1:A:521:MET:HE1	1:A:524:ARG:CZ	2.45	0.47
1:A:485:PRO:CG	1:A:488:ILE:HD12	2.44	0.46
1:A:727:VAL:CG2	1:A:789:MET:HE1	2.45	0.46
1:A:534:SER:C	1:A:536:SER:H	2.23	0.46
1:B:766:VAL:O	1:B:772:CYS:HB2	2.15	0.46
1:A:761:PRO:HD3	1:B:651:ARG:NH1	2.30	0.46
1:B:698:ASP:O	1:B:700:ASN:N	2.48	0.46
1:B:727:VAL:CG2	1:B:789:MET:HE1	2.46	0.45
1:B:695:GLY:C	1:B:697:THR:N	2.66	0.45
1:B:713:ASN:O	1:B:717:ILE:HG13	2.17	0.45
1:B:589:ALA:CB	1:B:664:MET:HE1	2.47	0.45
1:A:712:GLU:H	1:A:712:GLU:CD	2.24	0.45
1:A:541:TRP:O	1:A:544:ILE:HG22	2.16	0.45
1:A:760:LEU:HA	1:B:651:ARG:HH12	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:490:ARG:HA	1:B:493:GLU:HG3	1.99	0.45
1:B:664:MET:HE2	1:B:722:ILE:HB	1.99	0.45
1:B:692:GLU:HA	1:B:696:GLU:CG	2.37	0.44
1:B:696:GLU:O	1:B:700:ASN:HB3	2.16	0.44
1:B:711:PRO:HG2	1:B:712:GLU:OE2	2.17	0.44
1:A:592:ALA:O	1:A:596:HIS:HB3	2.18	0.44
1:A:817:ASP:O	1:A:819:MET:N	2.50	0.44
1:B:697:THR:HG22	1:B:698:ASP:N	2.33	0.44
1:B:697:THR:O	1:B:698:ASP:O	2.36	0.44
1:B:592:ALA:O	1:B:596:HIS:HB3	2.18	0.44
1:B:817:ASP:C	1:B:819:MET:N	2.76	0.44
1:B:485:PRO:CG	1:B:488:ILE:HD12	2.48	0.43
1:B:695:GLY:C	1:B:696:GLU:HG3	2.43	0.43
1:B:505:LEU:O	1:B:509:THR:HG23	2.19	0.43
1:B:528:CYS:HB3	1:B:533:CYS:O	2.18	0.43
1:A:583:PRO:O	1:A:587:VAL:HG23	2.18	0.43
1:B:650:GLU:O	1:B:651:ARG:C	2.61	0.43
1:A:597:ASP:O	1:A:600:HIS:HB2	2.19	0.43
1:B:817:ASP:O	1:B:819:MET:N	2.52	0.43
1:A:580:THR:HG21	1:A:717:ILE:CD1	2.49	0.42
1:A:638:THR:HA	1:A:644:ASN:OD1	2.19	0.42
2:A:3:IBM:O2	2:A:3:IBM:H133	2.20	0.42
1:A:693:GLU:C	1:A:695:GLY:N	2.77	0.42
1:B:693:GLU:C	1:B:695:GLY:N	2.78	0.42
1:B:785:PHE:CE1	2:B:3:IBM:H131	2.54	0.42
1:B:622:THR:O	1:B:623:ALA:C	2.62	0.42
1:A:817:ASP:C	1:A:819:MET:N	2.77	0.42
1:B:585:ASP:CB	1:B:718:LYS:HD2	2.49	0.41
1:A:575:GLU:CD	1:A:575:GLU:N	2.62	0.41
1:A:713:ASN:O	1:A:717:ILE:HG13	2.20	0.41
1:B:816:LEU:O	1:B:817:ASP:C	2.63	0.41
1:A:664:MET:HE2	1:A:722:ILE:HB	2.02	0.41
1:A:774:ILE:N	1:A:775:PRO:HD2	2.36	0.41
1:B:528:CYS:C	1:B:530:PHE:N	2.79	0.41
1:B:527:ILE:HD11	1:B:590:LEU:HD13	2.03	0.40
1:B:572:LEU:O	1:B:578:LYS:CD	2.68	0.40
1:B:748:TYR:CE2	1:B:764:MET:HE1	2.56	0.40
1:A:622:THR:O	1:A:623:ALA:C	2.64	0.40
1:B:608:LEU:HD11	1:B:625:LEU:HD21	2.04	0.40
1:B:712:GLU:CD	1:B:712:GLU:H	2.29	0.40
1:B:812:TYR:CE2	1:B:816:LEU:HD11	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:LEU:HB3	1:A:546:GLU:CD	2.45	0.40
1:B:694:ASN:C	1:B:696:GLU:N	2.79	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	336/338 (99%)	314 (94%)	17 (5%)	5 (2%)	8	4
1	B	336/338 (99%)	306 (91%)	24 (7%)	6 (2%)	6	3
All	All	672/676 (99%)	620 (92%)	41 (6%)	11 (2%)	7	4

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	493	GLU
1	A	698	ASP
1	B	697	THR
1	B	698	ASP
1	A	644	ASN
1	B	699	LYS
1	B	818	GLU
1	A	818	GLU
1	A	651	ARG
1	B	596	HIS
1	B	694	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	303/303 (100%)	298 (98%)	5 (2%)	53	62
1	B	303/303 (100%)	300 (99%)	3 (1%)	68	76
All	All	606/606 (100%)	598 (99%)	8 (1%)	61	69

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	495	GLU
1	A	534	SER
1	A	575	GLU
1	A	697	THR
1	A	771	THR
1	B	495	GLU
1	B	733	PRO
1	B	791	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	543	GLN
1	A	644	ASN
1	A	676	HIS
1	A	705	ASN
1	A	770	ASN
1	A	778	GLN
1	A	804	GLN
1	A	809	ASN
1	B	685	ASN
1	B	700	ASN
1	B	705	ASN
1	B	758	GLN
1	B	770	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](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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
2	IBM	A	3	-	17,17,17	0.99	1 (5%)	19,25,25	1.80	5 (26%)
2	IBM	B	3	-	17,17,17	0.99	1 (5%)	19,25,25	1.81	5 (26%)

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	IBM	A	3	-	-	2/4/4/4	0/2/2/2
2	IBM	B	3	-	-	2/4/4/4	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3	IBM	C5-C6	-2.53	1.39	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	3	IBM	C5-C6	-2.46	1.39	1.45

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3	IBM	N3-C2-N1	4.28	121.99	116.72
2	A	3	IBM	N3-C2-N1	4.27	121.99	116.72
2	B	3	IBM	C6-N1-C2	-3.20	120.45	125.66
2	A	3	IBM	C6-N1-C2	-3.15	120.54	125.66
2	A	3	IBM	C11-N3-C2	2.58	122.92	117.67
2	B	3	IBM	C11-N3-C2	2.55	122.85	117.67
2	B	3	IBM	C5-C6-N1	2.30	119.26	115.02
2	A	3	IBM	C5-C6-N1	2.23	119.13	115.02
2	B	3	IBM	O2-C2-N3	-2.22	118.99	121.98
2	A	3	IBM	O2-C2-N3	-2.17	119.05	121.98

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	3	IBM	N3-C11-C12-C13
2	A	3	IBM	N3-C11-C12-C14
2	B	3	IBM	N3-C11-C12-C13
2	B	3	IBM	N3-C11-C12-C14

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	3	IBM	3	0
2	B	3	IBM	3	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 [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	338/338 (100%)	-0.94	0 100 100	31, 46, 84, 111	0
1	B	338/338 (100%)	-0.94	0 100 100	31, 45, 85, 113	0
All	All	676/676 (100%)	-0.94	0 100 100	31, 46, 85, 113	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(Å ²)	Q<0.9
2	IBM	B	3	16/16	0.97	0.08	72,74,79,79	0
2	IBM	A	3	16/16	0.98	0.07	71,75,79,79	0
4	MG	A	2	1/1	0.99	0.03	37,37,37,37	0
4	MG	B	2	1/1	0.99	0.04	38,38,38,38	0
3	ZN	A	1	1/1	1.00	0.02	40,40,40,40	0
3	ZN	B	1	1/1	1.00	0.01	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.