



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:06 AM GMT

PDB ID : 2PD0
Title : Protein cgd2_2020 from *Cryptosporidium parvum*
Authors : Cymborowski, M.; Chruszcz, M.; Hills, T.; Lew, J.; Melone, M.; Zhao, Y.; Artz, J.; Wernimont, A.; Edwards, A.; Sundstrom, M.; Weigelt, J.; Bochkarev, A.; Arrowsmith, C.; Hui, R.; Minor, W.; Structural Genomics Consortium (SGC)
Deposited on : 2007-03-30
Resolution : 2.30 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

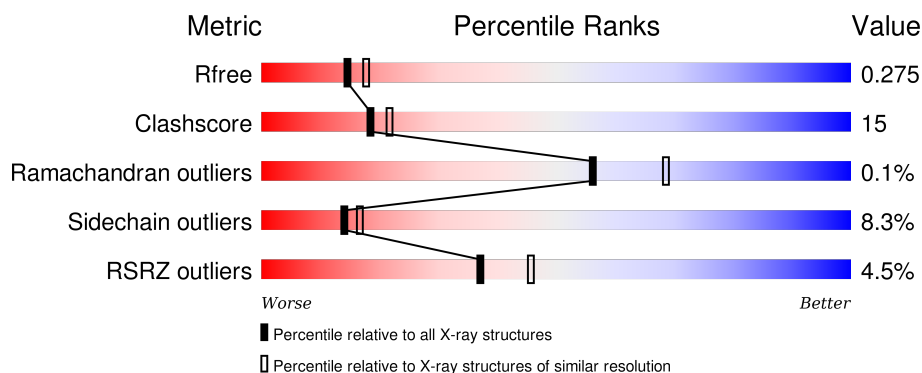
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	223	<div> <div>3%</div> <div>67% 20% 8%</div> </div>
1	B	223	<div> <div>3%</div> <div>60% 27% 8%</div> </div>
1	C	223	<div> <div>5%</div> <div>59% 27% 6% 8%</div> </div>
1	D	223	<div> <div>5%</div> <div>65% 23% 9%</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
2	MES	C	504	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7306 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 Hypothetical protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	Se	0	10	0
			1760	1126	290	338	3	3			
1	B	205	Total	C	N	O	S	Se	0	8	0
			1738	1115	288	328	4	3			
1	C	205	Total	C	N	O	S	Se	0	3	0
			1687	1083	280	319	2	3			
1	D	203	Total	C	N	O	S	Se	0	1	0
			1645	1058	271	311	2	3			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	CLONING ARTIFACT	UNP Q5CTR0
A	-17	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
A	-16	SER	-	CLONING ARTIFACT	UNP Q5CTR0
A	-15	SER	-	CLONING ARTIFACT	UNP Q5CTR0
A	-14	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
A	-13	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
A	-12	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
A	-11	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
A	-10	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
A	-9	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
A	-8	SER	-	CLONING ARTIFACT	UNP Q5CTR0
A	-7	SER	-	CLONING ARTIFACT	UNP Q5CTR0
A	-6	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
A	-5	LEU	-	CLONING ARTIFACT	UNP Q5CTR0
A	-4	VAL	-	CLONING ARTIFACT	UNP Q5CTR0
A	-3	PRO	-	CLONING ARTIFACT	UNP Q5CTR0
A	-2	ARG	-	CLONING ARTIFACT	UNP Q5CTR0
A	-1	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
A	0	SER	-	CLONING ARTIFACT	UNP Q5CTR0
A	25	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
A	131	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0

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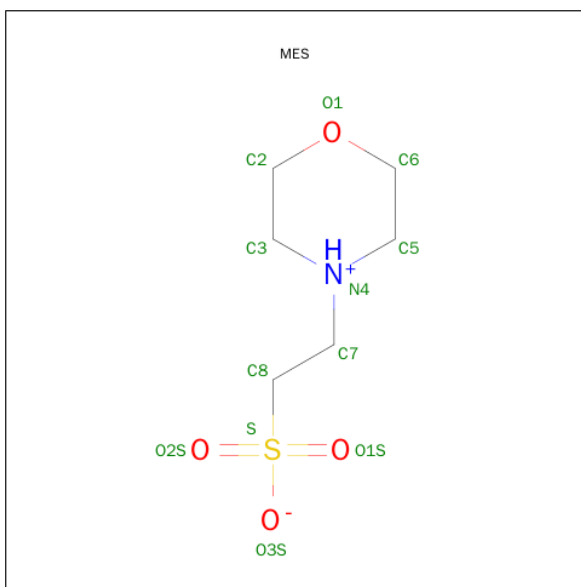
Chain	Residue	Modelled	Actual	Comment	Reference
A	164	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
B	-18	MET	-	CLONING ARTIFACT	UNP Q5CTR0
B	-17	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
B	-16	SER	-	CLONING ARTIFACT	UNP Q5CTR0
B	-15	SER	-	CLONING ARTIFACT	UNP Q5CTR0
B	-14	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
B	-13	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
B	-12	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
B	-11	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
B	-10	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
B	-9	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
B	-8	SER	-	CLONING ARTIFACT	UNP Q5CTR0
B	-7	SER	-	CLONING ARTIFACT	UNP Q5CTR0
B	-6	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
B	-5	LEU	-	CLONING ARTIFACT	UNP Q5CTR0
B	-4	VAL	-	CLONING ARTIFACT	UNP Q5CTR0
B	-3	PRO	-	CLONING ARTIFACT	UNP Q5CTR0
B	-2	ARG	-	CLONING ARTIFACT	UNP Q5CTR0
B	-1	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
B	0	SER	-	CLONING ARTIFACT	UNP Q5CTR0
B	25	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
B	131	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
B	164	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
C	-18	MET	-	CLONING ARTIFACT	UNP Q5CTR0
C	-17	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
C	-16	SER	-	CLONING ARTIFACT	UNP Q5CTR0
C	-15	SER	-	CLONING ARTIFACT	UNP Q5CTR0
C	-14	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
C	-13	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
C	-12	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
C	-11	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
C	-10	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
C	-9	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
C	-8	SER	-	CLONING ARTIFACT	UNP Q5CTR0
C	-7	SER	-	CLONING ARTIFACT	UNP Q5CTR0
C	-6	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
C	-5	LEU	-	CLONING ARTIFACT	UNP Q5CTR0
C	-4	VAL	-	CLONING ARTIFACT	UNP Q5CTR0
C	-3	PRO	-	CLONING ARTIFACT	UNP Q5CTR0
C	-2	ARG	-	CLONING ARTIFACT	UNP Q5CTR0
C	-1	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
C	0	SER	-	CLONING ARTIFACT	UNP Q5CTR0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	25	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
C	131	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
C	164	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
D	-18	MET	-	CLONING ARTIFACT	UNP Q5CTR0
D	-17	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
D	-16	SER	-	CLONING ARTIFACT	UNP Q5CTR0
D	-15	SER	-	CLONING ARTIFACT	UNP Q5CTR0
D	-14	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
D	-13	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
D	-12	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
D	-11	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
D	-10	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
D	-9	HIS	-	CLONING ARTIFACT	UNP Q5CTR0
D	-8	SER	-	CLONING ARTIFACT	UNP Q5CTR0
D	-7	SER	-	CLONING ARTIFACT	UNP Q5CTR0
D	-6	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
D	-5	LEU	-	CLONING ARTIFACT	UNP Q5CTR0
D	-4	VAL	-	CLONING ARTIFACT	UNP Q5CTR0
D	-3	PRO	-	CLONING ARTIFACT	UNP Q5CTR0
D	-2	ARG	-	CLONING ARTIFACT	UNP Q5CTR0
D	-1	GLY	-	CLONING ARTIFACT	UNP Q5CTR0
D	0	SER	-	CLONING ARTIFACT	UNP Q5CTR0
D	25	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
D	131	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0
D	164	MSE	MET	MODIFIED RESIDUE	UNP Q5CTR0

- Molecule 2 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

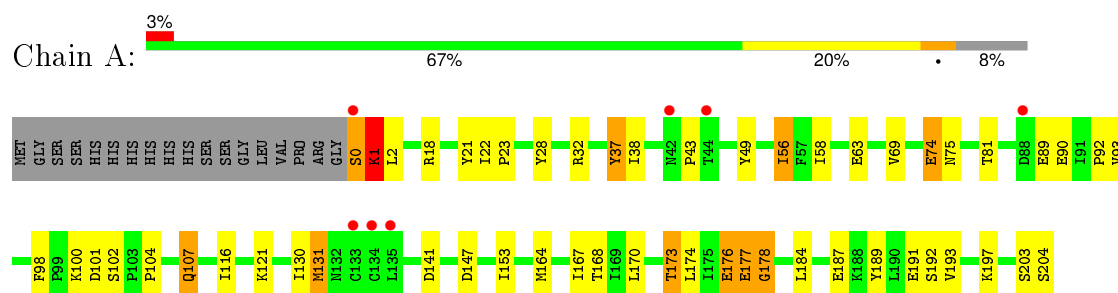
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	123	Total	O	0	0
			123	123		
3	B	130	Total	O	0	0
			130	130		
3	C	76	Total	O	0	0
			76	76		
3	D	99	Total	O	0	0
			99	99		

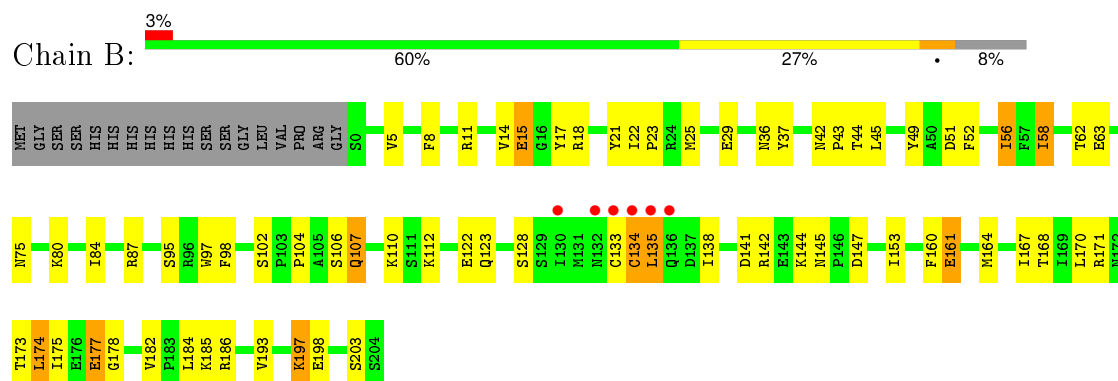
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

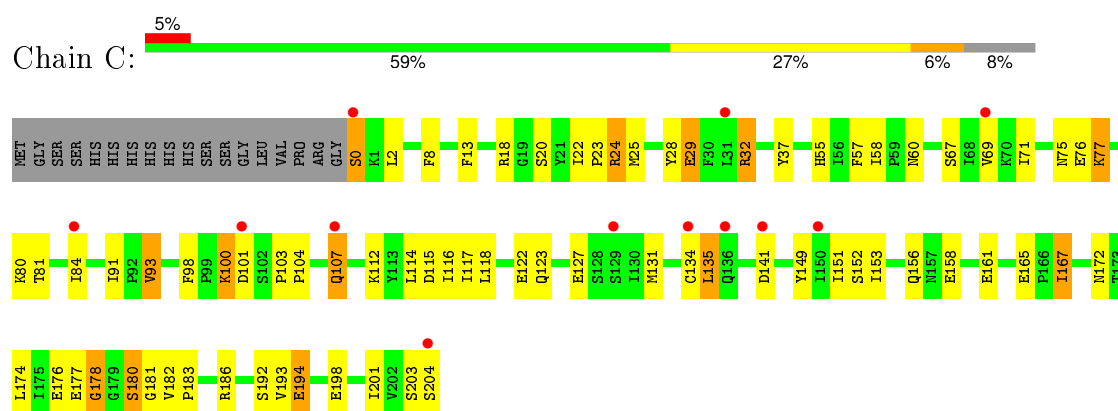
- Molecule 1: Hypothetical protein



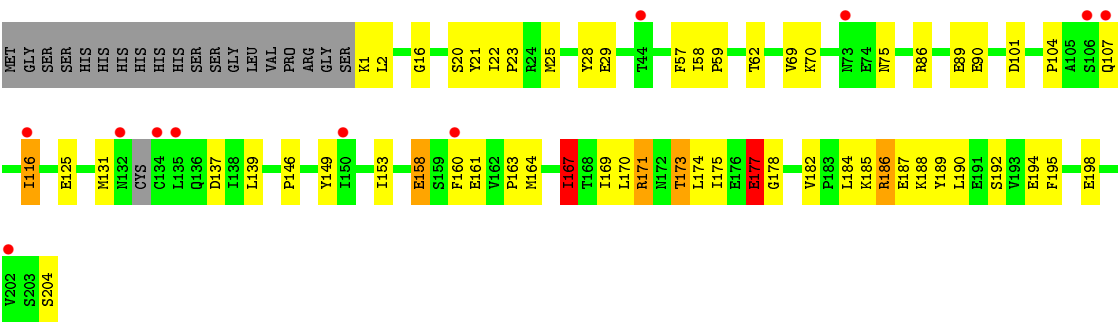
- Molecule 1: Hypothetical protein



- Molecule 1: Hypothetical protein



- Molecule 1: Hypothetical protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	49.53Å 68.52Å 79.57Å 69.29° 75.14° 93.73°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-2.30) 85.0 (20.00-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.69 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.220 , 0.280 0.224 , 0.275	Depositor DCC
R_{free} test set	2091 reflections (5.43%)	DCC
Wilson B-factor (Å ²)	28.8	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 40584 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7306	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.17% 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: MES

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	0.92	2/1796 (0.1%)	0.81	0/2420
1	B	1.04	4/1773 (0.2%)	0.91	2/2391 (0.1%)
1	C	0.87	1/1722 (0.1%)	0.82	0/2325
1	D	0.96	1/1680 (0.1%)	0.85	3/2272 (0.1%)
All	All	0.95	8/6971 (0.1%)	0.85	5/9408 (0.1%)

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	5
1	B	0	2
1	C	0	4
1	D	0	1
All	All	0	12

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	131	MSE	SE-CE	-7.25	1.52	1.95
1	B	134	CYS	CB-SG	6.68	1.93	1.82
1	B	17	TYR	CD1-CE1	-6.66	1.29	1.39
1	B	17	TYR	CD2-CE2	-5.95	1.30	1.39
1	A	37	TYR	CD1-CE1	-5.93	1.30	1.39
1	D	189	TYR	CD1-CE1	-5.81	1.30	1.39
1	B	182	VAL	CB-CG2	-5.56	1.41	1.52
1	A	37	TYR	CD2-CE2	-5.36	1.31	1.39

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	171	ARG	NE-CZ-NH2	-5.97	117.32	120.30
1	D	167	ILE	CB-CA-C	-5.80	100.00	111.60
1	B	171	ARG	NE-CZ-NH2	-5.31	117.64	120.30
1	B	147	ASP	CB-CG-OD2	-5.12	113.69	118.30
1	D	171	ARG	NE-CZ-NH1	5.11	122.85	120.30

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	0	SER	Peptide
1	A	1	LYS	Peptide
1	A	131	MSE	Peptide
1	A	177	GLU	Peptide
1	A	178	GLY	Peptide
1	B	177	GLU	Peptide
1	B	178	GLY	Peptide
1	C	0	SER	Peptide
1	C	177	GLU	Peptide
1	C	178	GLY	Peptide
1	C	180	SER	Peptide
1	D	177	GLU	Peptide

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	1760	0	1729	52	0
1	B	1738	0	1716	50	0
1	C	1687	0	1650	68	0
1	D	1645	0	1594	48	0
2	A	12	0	12	0	0
2	B	12	0	12	3	0
2	C	12	0	12	2	0
2	D	12	0	12	1	0
3	A	123	0	0	8	0
3	B	130	0	0	7	0
3	C	76	0	0	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	99	0	0	14	0
All	All	7306	0	6737	211	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 15.

All (211) 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:93:VAL:HG11	1:A:167[B]:ILE:HG12	1.28	1.15
1:A:56[A]:ILE:HD11	1:A:58:ILE:CG2	1.79	1.12
1:A:56[A]:ILE:HD11	1:A:58:ILE:HG23	1.28	1.11
1:D:1:LYS:CB	1:D:2:LEU:HA	1.82	1.10
1:C:135:LEU:HD23	1:C:135:LEU:N	1.71	1.05
1:C:84:ILE:HG23	3:C:547:HOH:O	1.63	0.95
1:C:135:LEU:CD2	1:C:135:LEU:N	2.30	0.94
1:A:93:VAL:CG1	1:A:167[B]:ILE:HG12	2.00	0.91
1:B:18:ARG:HH21	1:B:18:ARG:HG2	1.36	0.90
1:A:56[A]:ILE:CD1	1:A:58:ILE:HG23	2.04	0.87
1:C:20:SER:HA	3:C:510:HOH:O	1.78	0.82
1:A:74:GLU:HG2	1:A:75:ASN:N	1.94	0.82
1:B:161[A]:GLU:HG3	3:B:559:HOH:O	1.80	0.82
1:C:134:CYS:O	1:C:135:LEU:HD22	1.76	0.82
1:C:134:CYS:C	1:C:135:LEU:CD2	2.50	0.80
1:B:22[A]:ILE:CD1	1:B:153:ILE:HD13	2.10	0.80
1:C:167[A]:ILE:HD13	3:C:578:HOH:O	1.82	0.80
3:A:521:HOH:O	1:C:167[A]:ILE:HG12	1.82	0.79
1:C:134:CYS:C	1:C:135:LEU:HD23	2.03	0.79
1:A:193:VAL:O	1:A:197[A]:LYS:HG3	1.81	0.79
1:D:171:ARG:HD3	3:D:534:HOH:O	1.82	0.78
1:D:86:ARG:HD3	1:D:90:GLU:OE1	1.84	0.78
1:D:1:LYS:CB	1:D:2:LEU:CA	2.62	0.77
1:D:58:ILE:HD12	1:D:116:ILE:HD13	1.65	0.76
1:A:167[A]:ILE:HG12	3:A:531:HOH:O	1.86	0.75
1:B:22[A]:ILE:HD11	1:B:153:ILE:HD13	1.68	0.74
1:B:170:LEU:O	1:B:173:THR:HB	1.88	0.73
1:C:158:GLU:N	1:C:158:GLU:OE1	2.20	0.72
1:C:182:VAL:HG13	1:C:183:PRO:HD2	1.71	0.72
1:D:70:LYS:HE3	1:D:198:GLU:HG3	1.71	0.71
1:D:204:SER:HA	3:D:599:HOH:O	1.91	0.69
1:B:18:ARG:NH2	1:B:203:SER:OG	2.26	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:ARG:NE	3:B:611:HOH:O	2.27	0.68
1:B:51:ASP:HB3	3:B:619:HOH:O	1.93	0.68
1:D:86:ARG:CD	1:D:90:GLU:OE1	2.41	0.68
1:C:182:VAL:CG1	1:C:183:PRO:HD2	2.24	0.68
1:A:93:VAL:HG11	1:A:167[B]:ILE:CG1	2.16	0.67
1:C:84:ILE:HD12	3:C:547:HOH:O	1.93	0.67
1:D:57:PHE:HZ	1:D:161:GLU:HG3	1.61	0.65
1:B:42[A]:ASN:ND2	3:B:562:HOH:O	2.30	0.65
1:A:89:GLU:OE1	1:A:89:GLU:N	2.31	0.64
1:C:134:CYS:C	1:C:135:LEU:HD22	2.17	0.63
1:A:100:LYS:HB3	1:A:204:SER:HB3	1.80	0.63
1:D:58:ILE:HG22	1:D:59:PRO:O	1.97	0.63
1:D:21:TYR:HB2	3:D:599:HOH:O	1.99	0.63
1:C:100:LYS:HE3	1:C:100:LYS:HA	1.80	0.62
1:D:185:LYS:HE2	3:D:577:HOH:O	1.99	0.61
1:C:134:CYS:O	1:C:135:LEU:CD2	2.46	0.61
1:D:21:TYR:CE1	1:D:23:PRO:HB3	2.36	0.61
1:C:84:ILE:CD1	3:C:547:HOH:O	2.49	0.61
1:C:172:ASN:O	1:C:181:GLY:CA	2.49	0.61
1:B:145:ASN:ND2	3:B:627:HOH:O	2.34	0.60
1:C:123:GLN:NE2	1:C:180:SER:O	2.34	0.60
1:A:102:SER:OG	1:A:104:PRO:HD2	2.02	0.60
1:B:18:ARG:NH2	1:B:18:ARG:HG2	2.14	0.59
1:D:187:GLU:H	1:D:187:GLU:CD	2.06	0.59
1:C:165:GLU:HB3	1:C:167[A]:ILE:HG13	1.82	0.59
1:A:170:LEU:HD23	1:A:184:LEU:CD1	2.33	0.59
1:D:167:ILE:HG12	3:D:507:HOH:O	2.02	0.59
1:B:128:SER:HA	1:B:133:CYS:SG	2.43	0.59
1:A:191:GLU:HG2	1:B:160:PHE:CD1	2.38	0.58
1:A:18:ARG:NH1	3:A:620:HOH:O	2.36	0.58
1:C:167[A]:ILE:HG23	3:C:575:HOH:O	2.03	0.58
1:C:172:ASN:O	1:C:182:VAL:N	2.35	0.58
1:C:100:LYS:HD2	1:C:204:SER:HB3	1.85	0.58
1:A:92:PRO:HG2	1:C:193:VAL:HG21	1.85	0.58
1:A:164:MSE:O	1:A:192:SER:OG	2.21	0.58
1:A:197[B]:LYS:HG2	3:C:565:HOH:O	2.04	0.57
1:C:25:MSE:HG3	1:C:29:GLU:HG2	1.85	0.57
1:D:169:ILE:HG22	1:D:184:LEU:HD12	1.86	0.57
1:A:58:ILE:HD11	1:A:116:ILE:HD11	1.85	0.57
1:B:186:ARG:HD2	1:D:177:GLU:OE2	2.04	0.57
1:C:180:SER:H	1:C:181:GLY:HA2	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32[B]:ARG:NH1	1:A:141:ASP:OD1	2.39	0.55
1:D:170:LEU:O	1:D:173:THR:HB	2.06	0.55
1:C:28:TYR:HE2	1:C:32:ARG:NH1	2.04	0.54
1:C:37:TYR:HE1	3:C:573:HOH:O	1.89	0.54
1:B:43:PRO:HG2	1:B:56[A]:ILE:HG21	1.88	0.54
1:A:23:PRO:HG2	1:A:63[B]:GLU:HB2	1.88	0.54
1:B:18:ARG:HH21	1:B:18:ARG:CG	2.13	0.54
1:A:56[A]:ILE:CD1	1:A:58:ILE:CG2	2.69	0.53
1:A:49:TYR:CG	1:A:164:MSE:HE1	2.43	0.53
1:D:86:ARG:NH1	1:D:90:GLU:OE2	2.40	0.52
1:A:18:ARG:HD3	1:A:203:SER:OG	2.10	0.52
1:B:185:LYS:HE3	3:B:613:HOH:O	2.09	0.52
1:A:81:THR:HG21	1:C:81:THR:CG2	2.40	0.52
1:A:22:ILE:HD11	1:A:153:ILE:HG21	1.92	0.52
1:D:101[A]:ASP:OD1	1:D:101[A]:ASP:N	2.40	0.51
1:B:25:MSE:HE1	1:B:62:THR:HG22	1.92	0.51
1:B:173:THR:HG22	1:B:174:LEU:HG	1.93	0.51
1:B:49:TYR:CG	1:B:164:MSE:HE1	2.46	0.51
1:A:21:TYR:CE1	1:A:23:PRO:HB3	2.46	0.51
1:B:102:SER:HB2	1:B:104:PRO:HD2	1.91	0.51
1:B:36:ASN:OD1	1:B:142:ARG:NH1	2.43	0.51
1:C:127:GLU:OE1	2:C:504:MES:C7	2.58	0.50
1:C:57:PHE:HZ	1:C:161[A]:GLU:HG3	1.75	0.50
1:B:84:ILE:HG13	3:B:611:HOH:O	2.12	0.50
1:D:204:SER:C	3:D:599:HOH:O	2.48	0.50
1:D:177:GLU:H	1:D:178:GLY:HA2	1.77	0.50
1:A:187:GLU:O	1:A:191:GLU:HG3	2.11	0.49
1:A:81:THR:CG2	1:C:81:THR:HG21	2.43	0.49
1:C:103:PRO:O	1:C:107:GLN:NE2	2.43	0.49
1:D:20:SER:HA	3:D:523:HOH:O	2.13	0.49
1:D:173:THR:HG22	1:D:174:LEU:HG	1.95	0.49
1:C:69:VAL:HG13	1:C:98:PHE:CE2	2.47	0.48
1:B:56[B]:ILE:HD12	1:B:58:ILE:CD1	2.43	0.48
1:A:107:GLN:HB2	3:A:615:HOH:O	2.12	0.48
1:C:180:SER:N	1:C:181:GLY:HA2	2.28	0.48
1:D:131:MSE:SE	1:D:178:GLY:HA3	2.63	0.48
1:D:25:MSE:HG3	1:D:29:GLU:HG2	1.96	0.48
1:A:176:GLU:HG2	1:A:177:GLU:HG2	1.93	0.48
1:A:121:LYS:HD2	3:A:613:HOH:O	2.13	0.48
1:B:80:LYS:HD3	1:B:97:TRP:CZ2	2.49	0.48
1:C:23:PRO:C	1:C:24[A]:ARG:HG2	2.34	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:194:GLU:HA	1:C:194:GLU:OE1	2.14	0.47
1:D:177:GLU:N	1:D:178:GLY:HA2	2.29	0.47
1:B:87:ARG:HH21	2:B:503:MES:H31	1.80	0.47
1:A:90:GLU:HA	3:A:612:HOH:O	2.14	0.47
1:C:22:ILE:HD11	1:C:153:ILE:HD13	1.96	0.47
1:B:8:PHE:CD2	2:B:503:MES:H32	2.50	0.47
1:B:52:PHE:CD1	1:B:123:GLN:HG2	2.50	0.47
1:A:189:TYR:CE1	1:C:91:ILE:HD12	2.50	0.47
1:B:37:TYR:CD2	1:B:58:ILE:HG12	2.49	0.47
1:B:23:PRO:HG2	1:B:63:GLU:HB2	1.96	0.47
1:C:107:GLN:HG2	1:C:107:GLN:H	1.56	0.46
1:D:175:ILE:HG13	1:D:175:ILE:O	2.15	0.46
1:A:58:ILE:CD1	1:A:116:ILE:HD11	2.45	0.46
1:D:86:ARG:NH2	2:D:501:MES:O3S	2.48	0.46
1:D:22:ILE:HD12	1:D:153:ILE:HD13	1.96	0.46
1:D:182:VAL:CG1	3:D:577:HOH:O	2.64	0.46
1:B:135:LEU:CD2	1:B:135:LEU:N	2.78	0.46
1:C:75:ASN:HB2	1:C:104:PRO:HG3	1.97	0.46
1:B:141:ASP:HA	1:B:144:LYS:HE2	1.97	0.46
1:C:8:PHE:CG	2:C:504:MES:H32	2.51	0.46
1:B:175[B]:ILE:C	1:B:177:GLU:H	2.18	0.46
1:B:167:ILE:HD12	1:B:168:THR:N	2.31	0.45
1:B:15:GLU:HG3	1:B:21:TYR:CE1	2.52	0.45
1:B:25:MSE:HG2	1:B:29:GLU:HB3	1.98	0.45
1:C:180:SER:N	1:C:181:GLY:CA	2.79	0.45
1:D:16:GLY:HA2	1:D:204:SER:OG	2.17	0.45
1:B:75:ASN:HB2	1:B:104:PRO:HG3	1.98	0.45
1:C:13:PHE:CZ	1:C:22:ILE:HD13	2.52	0.45
1:B:167:ILE:HD12	1:B:168:THR:H	1.80	0.45
1:B:87:ARG:NH2	2:B:503:MES:H31	2.31	0.45
1:A:0:SER:O	1:A:147:ASP:HB2	2.17	0.45
1:C:172:ASN:O	1:C:181:GLY:HA3	2.17	0.45
1:C:93:VAL:CG1	1:C:167[A]:ILE:HG21	2.47	0.45
1:A:167[A]:ILE:HG21	3:A:611:HOH:O	2.17	0.45
1:D:86:ARG:HD2	1:D:90:GLU:OE1	2.16	0.44
1:C:167[B]:ILE:HD13	3:C:575:HOH:O	2.17	0.44
1:A:69:VAL:HG22	1:A:98:PHE:CE2	2.53	0.44
1:C:60:ASN:OD1	1:C:114:LEU:HD11	2.18	0.44
1:D:182:VAL:HG13	3:D:577:HOH:O	2.17	0.44
1:C:77:LYS:NZ	1:C:77:LYS:CB	2.81	0.44
1:B:193:VAL:O	1:B:197:LYS:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:ILE:HD12	1:C:84:ILE:HG23	1.62	0.44
1:D:57:PHE:CZ	1:D:161:GLU:HG3	2.48	0.44
1:C:167[A]:ILE:HG12	3:C:505:HOH:O	2.17	0.43
1:C:117:ILE:HB	1:C:152:SER:HB3	2.00	0.43
1:B:164:MSE:HG3	1:B:168:THR:HB	2.00	0.43
1:B:174:LEU:HD21	1:D:173:THR:HG21	1.99	0.43
1:B:112:LYS:HE2	1:B:112:LYS:HB3	1.52	0.43
1:D:163:PRO:HG2	1:D:195:PHE:CG	2.54	0.43
1:C:58:ILE:HD12	1:C:116:ILE:CD1	2.48	0.43
1:C:18:ARG:HD3	1:C:203:SER:HB2	1.99	0.43
1:B:98:PHE:CD1	1:B:104:PRO:HG2	2.53	0.43
1:C:28:TYR:CE2	1:C:32:ARG:NH1	2.85	0.43
1:B:45:LEU:HA	1:B:45:LEU:HD12	1.85	0.43
1:C:122:GLU:HG3	3:C:570:HOH:O	2.17	0.43
1:C:174:LEU:O	1:C:178:GLY:HA2	2.19	0.43
1:A:167[A]:ILE:HD12	1:A:168:THR:N	2.33	0.43
1:D:177:GLU:HA	1:D:177:GLU:OE1	2.18	0.43
1:C:55:HIS:CE1	1:C:161[A]:GLU:HG2	2.54	0.43
1:B:144:LYS:HD3	1:B:144:LYS:HA	1.84	0.43
1:D:146:PRO:HG2	1:D:149:TYR:CZ	2.53	0.43
1:B:5:VAL:HB	1:B:138:ILE:HB	2.00	0.43
1:C:93:VAL:CG2	1:C:167[B]:ILE:HG12	2.49	0.42
1:C:71:ILE:N	1:C:198:GLU:O	2.29	0.42
1:A:174:LEU:O	1:A:178:GLY:HA2	2.19	0.42
1:B:107:GLN:HG2	1:B:107:GLN:H	1.62	0.42
1:C:93:VAL:HG21	1:C:167[B]:ILE:HG12	2.02	0.42
1:A:176:GLU:HB3	3:A:560:HOH:O	2.19	0.42
1:D:204:SER:CA	3:D:599:HOH:O	2.58	0.42
1:D:164:MSE:CE	3:D:580:HOH:O	2.67	0.42
1:C:71:ILE:HG13	1:C:198:GLU:C	2.40	0.42
1:D:75:ASN:HB2	1:D:104:PRO:HG3	2.01	0.42
1:B:197:LYS:HE3	1:B:198:GLU:OE1	2.20	0.42
1:A:23:PRO:HG2	1:A:63[B]:GLU:CB	2.49	0.42
1:A:176:GLU:HG3	1:A:176:GLU:O	2.19	0.42
1:C:118:LEU:HA	1:C:149:TYR:O	2.20	0.42
1:C:55:HIS:HD1	1:C:115:ASP:CG	2.23	0.41
1:D:28:TYR:HE1	1:D:139:LEU:O	2.03	0.41
1:A:28:TYR:O	1:A:32[B]:ARG:HG2	2.20	0.41
1:A:38:ILE:HA	1:A:43:PRO:HD2	2.02	0.41
1:A:167[A]:ILE:H	1:A:167[A]:ILE:HG13	1.68	0.41
1:C:156:GLN:HB2	1:C:158:GLU:OE1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32[B]:ARG:HB3	1:A:32[B]:ARG:CZ	2.50	0.41
1:C:93:VAL:HG21	1:C:167[B]:ILE:CG1	2.51	0.41
1:D:171:ARG:NH2	3:D:526:HOH:O	2.43	0.41
1:A:1:LYS:HB2	1:A:2:LEU:H	1.79	0.41
1:A:37:TYR:CE2	1:A:58:ILE:HG22	2.55	0.41
1:D:70:LYS:HE3	1:D:198:GLU:CG	2.44	0.41
1:D:186:ARG:NH2	3:D:584:HOH:O	2.52	0.41
1:A:58:ILE:CD1	1:A:116:ILE:CD1	3.00	0.40
1:A:32[A]:ARG:CZ	1:A:32[A]:ARG:HB3	2.51	0.40
1:C:93:VAL:HG13	1:C:167[A]:ILE:HG21	2.03	0.40
1:C:151:ILE:HD12	1:C:152:SER:HB2	2.03	0.40
1:A:173:THR:HG22	1:A:174:LEU:HG	2.02	0.40
1:B:106:SER:OG	1:B:107:GLN:NE2	2.53	0.40
1:D:137:ASP:OD1	3:D:593:HOH:O	2.22	0.40
1:D:158:GLU:HB2	1:D:160:PHE:CE2	2.57	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	213/223 (96%)	205 (96%)	8 (4%)	0	100	100
1	B	211/223 (95%)	204 (97%)	7 (3%)	0	100	100
1	C	206/223 (92%)	199 (97%)	5 (2%)	2 (1%)	19	21
1	D	200/223 (90%)	197 (98%)	3 (2%)	0	100	100
All	All	830/892 (93%)	805 (97%)	23 (3%)	2 (0%)	56	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	24[A]	ARG
1	C	24[B]	ARG

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	197/203 (97%)	186 (94%)	11 (6%)	26	35
1	B	193/203 (95%)	175 (91%)	18 (9%)	11	13
1	C	185/203 (91%)	163 (88%)	22 (12%)	6	7
1	D	178/203 (88%)	163 (92%)	15 (8%)	14	16
All	All	753/812 (93%)	687 (91%)	66 (9%)	14	14

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LYS
1	A	56[A]	ILE
1	A	56[B]	ILE
1	A	74	GLU
1	A	101[A]	ASP
1	A	101[B]	ASP
1	A	107	GLN
1	A	130	ILE
1	A	131	MSE
1	A	173	THR
1	A	176	GLU
1	B	11	ARG
1	B	14	VAL
1	B	15	GLU
1	B	44	THR
1	B	56[A]	ILE
1	B	56[B]	ILE
1	B	58	ILE
1	B	95	SER
1	B	107	GLN

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Mol	Chain	Res	Type
1	B	110	LYS
1	B	122	GLU
1	B	134	CYS
1	B	135	LEU
1	B	161[A]	GLU
1	B	161[B]	GLU
1	B	174	LEU
1	B	184	LEU
1	B	197	LYS
1	C	0	SER
1	C	2	LEU
1	C	29	GLU
1	C	32	ARG
1	C	67	SER
1	C	76	GLU
1	C	77	LYS
1	C	80	LYS
1	C	93	VAL
1	C	100	LYS
1	C	101	ASP
1	C	107	GLN
1	C	112	LYS
1	C	135	LEU
1	C	141	ASP
1	C	167[A]	ILE
1	C	167[B]	ILE
1	C	176	GLU
1	C	186	ARG
1	C	192	SER
1	C	194	GLU
1	C	201	ILE
1	D	62	THR
1	D	69	VAL
1	D	89	GLU
1	D	107	GLN
1	D	116	ILE
1	D	125	GLU
1	D	158	GLU
1	D	167	ILE
1	D	173	THR
1	D	177	GLU
1	D	186	ARG

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Mol	Chain	Res	Type
1	D	188	LYS
1	D	190	LEU
1	D	192	SER
1	D	194	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	107	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

4 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	MES	A	502	-	11,12,12	0.95	0	14,16,16	2.93	6 (42%)
2	MES	B	503	-	11,12,12	0.89	0	14,16,16	2.18	3 (21%)
2	MES	C	504	-	11,12,12	0.78	0	14,16,16	2.58	6 (42%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MES	D	501	-	11,12,12	0.80	0	14,16,16	2.74	4 (28%)

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	MES	A	502	-	-	0/6/14/14	0/1/1/1
2	MES	B	503	-	-	0/6/14/14	0/1/1/1
2	MES	C	504	-	-	0/6/14/14	0/1/1/1
2	MES	D	501	-	-	0/6/14/14	0/1/1/1

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	504	MES	C6-C5-N4	-2.50	106.34	110.12
2	D	501	MES	C7-N4-C3	2.05	116.52	111.27
2	C	504	MES	C2-C3-N4	2.20	113.47	110.12
2	A	502	MES	C6-C5-N4	2.39	113.74	110.12
2	A	502	MES	C7-N4-C5	2.63	118.00	111.27
2	A	502	MES	C7-N4-C3	2.97	118.89	111.27
2	C	504	MES	C7-N4-C3	2.98	118.91	111.27
2	B	503	MES	C7-N4-C3	3.01	118.97	111.27
2	C	504	MES	C7-N4-C5	3.14	119.32	111.27
2	B	503	MES	C7-N4-C5	3.20	119.46	111.27
2	D	501	MES	C7-N4-C5	3.34	119.82	111.27
2	A	502	MES	O2S-S-C8	3.72	110.08	106.91
2	C	504	MES	O2S-S-C8	4.52	110.76	106.91
2	D	501	MES	O2S-S-C8	5.41	111.52	106.91
2	A	502	MES	O1S-S-C8	5.73	111.79	106.91
2	B	503	MES	C5-N4-C3	5.84	121.56	108.90
2	C	504	MES	C5-N4-C3	5.94	121.77	108.90
2	A	502	MES	C5-N4-C3	6.55	123.09	108.90
2	D	501	MES	C5-N4-C3	6.75	123.51	108.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	503	MES	3	0
2	C	504	MES	2	0
2	D	501	MES	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	202/223 (90%)	0.24	7 (3%) 48 56	11, 25, 41, 59	0
1	B	202/223 (90%)	0.29	6 (2%) 54 63	13, 25, 39, 53	0
1	C	202/223 (90%)	0.46	12 (5%) 26 34	15, 29, 46, 61	0
1	D	200/223 (89%)	0.47	11 (5%) 29 37	16, 30, 44, 63	0
All	All	806/892 (90%)	0.36	36 (4%) 37 46	11, 27, 44, 63	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	134	CYS	9.2
1	B	133	CYS	8.0
1	B	135	LEU	7.0
1	A	134	CYS	6.1
1	D	134	CYS	5.7
1	C	134	CYS	5.6
1	D	106	SER	5.2
1	D	132	ASN	5.1
1	B	132	ASN	4.5
1	A	42	ASN	4.5
1	C	0	SER	4.4
1	D	44	THR	3.8
1	C	84	ILE	3.8
1	A	0	SER	3.3
1	C	204	SER	3.3
1	C	150	ILE	3.3
1	D	150	ILE	3.3
1	C	129	SER	3.2
1	B	130	ILE	3.0
1	D	135	LEU	2.9
1	D	202	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	107	GLN	2.7
1	D	116	ILE	2.6
1	D	73	ASN	2.6
1	A	88	ASP	2.6
1	A	135	LEU	2.6
1	A	133	CYS	2.4
1	C	31	LEU	2.3
1	D	160	PHE	2.3
1	C	101	ASP	2.2
1	B	136	GLN	2.1
1	C	136	GLN	2.1
1	C	69	VAL	2.1
1	A	44	THR	2.1
1	C	141	ASP	2.0
1	C	107	GLN	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
2	MES	C	504	12/12	0.86	0.24	2.79	54,56,57,59	0
2	MES	B	503	12/12	0.91	0.21	1.75	39,41,43,43	0
2	MES	D	501	12/12	0.93	0.20	1.26	36,38,41,42	0
2	MES	A	502	12/12	0.93	0.16	0.51	31,37,49,50	0

6.5 Other polymers

There are no such residues in this entry.