



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:42 AM UTC

PDB ID : 5T0W / pdb_00005t0w
Title : Crystal structure of the ancestral amino acid-binding protein AncCDT-1, a precursor of cyclohexadienyl dehydratase
Authors : Clifton, B.E.; Carr, P.D.; Jackson, C.J.
Deposited on : 2016-08-16
Resolution : 2.59 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

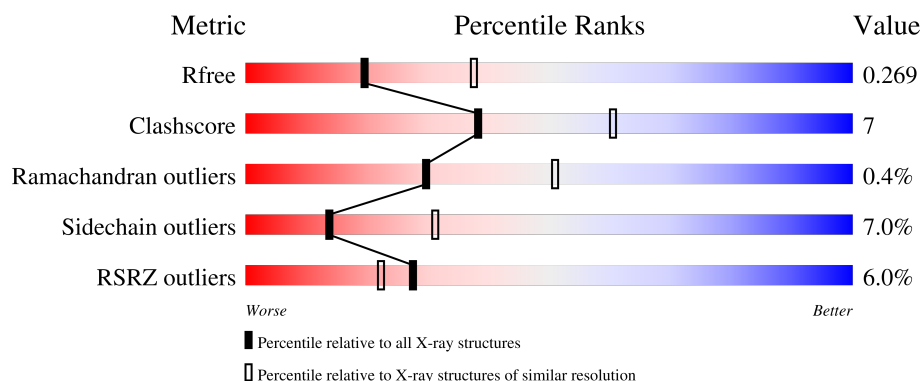
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	247	
1	B	247	
1	C	247	
1	D	247	

2 Entry composition [i](#)

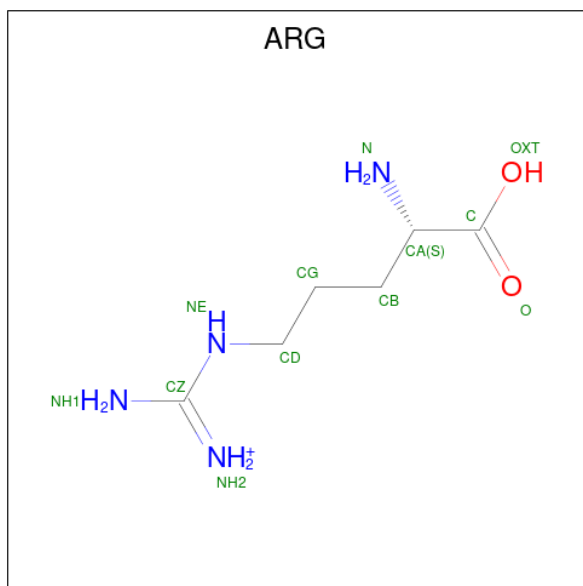
There are 3 unique types of molecules in this entry. The entry contains 7011 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 AncCDT-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	0	0
			1769	1136	287	340	6			
1	B	227	Total	C	N	O	S	0	0	0
			1734	1118	281	329	6			
1	C	229	Total	C	N	O	S	0	0	0
			1718	1106	276	330	6			
1	D	225	Total	C	N	O	S	0	1	0
			1721	1103	282	330	6			

- Molecule 2 is ARGinine (CCD ID: ARG) (formula: C₆H₁₅N₄O₂).



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

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

- Molecule 3 is water.

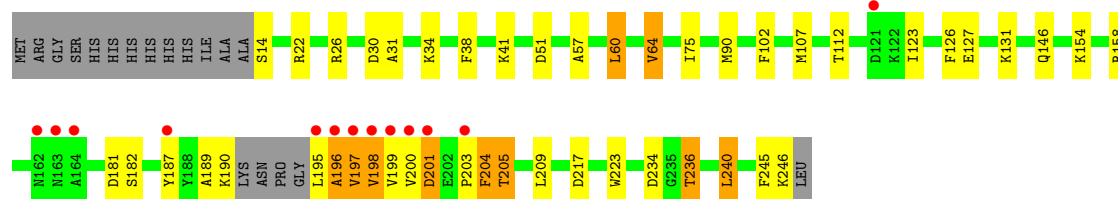
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	3	Total	O	0	0
			3	3		
3	D	7	Total	O	0	0
			7	7		

3 Residue-property plots [i](#)

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.

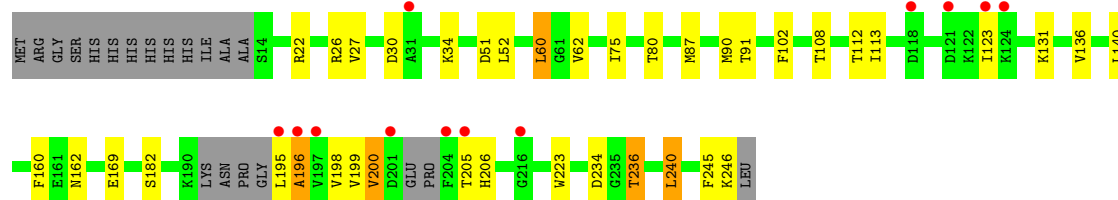
• Molecule 1: AncCDT-1

Chain A: 



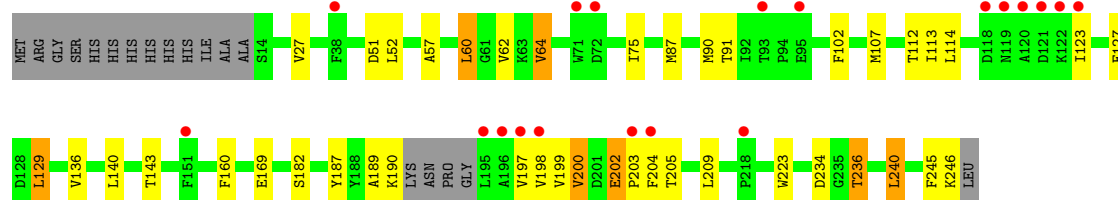
• Molecule 1: AncCDT-1

Chain B: 



• Molecule 1: AncCDT-1

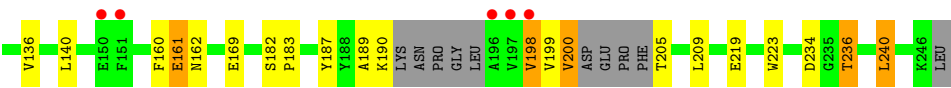
Chain C: 



• Molecule 1: AncCDT-1

Chain D: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.03Å 68.88Å 318.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.19 – 2.59 33.19 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.9 (33.19-2.59) 99.0 (33.19-2.59)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.57Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.253 , 0.287 (Not available) , 0.269	Depositor DCC
R_{free} test set	1686 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	40.2	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	7011	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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 [i](#)

5.1 Standard geometry [i](#)

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	1/1807 (0.1%)	0.95	6/2453 (0.2%)
1	B	0.73	0/1770	0.89	3/2402 (0.1%)
1	C	0.67	0/1756	0.86	2/2395 (0.1%)
1	D	0.76	0/1758	0.88	1/2386 (0.0%)
All	All	0.76	1/7091 (0.0%)	0.89	12/9636 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	197	VAL	N-CA	5.56	1.53	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	198	VAL	N-CA-C	8.02	120.87	108.44
1	A	198	VAL	N-CA-C	7.56	119.64	108.45
1	D	198	VAL	N-CA-C	6.64	118.36	108.46
1	A	123	ILE	N-CA-C	6.63	116.43	106.42
1	A	200	VAL	CB-CA-C	-6.43	100.94	110.33
1	C	204	PHE	N-CA-C	6.41	119.65	110.10
1	A	196	ALA	N-CA-C	5.96	123.48	110.80
1	A	64	VAL	CB-CA-C	-5.95	101.76	110.81
1	A	205	THR	N-CA-C	5.89	123.36	110.80
1	B	196	ALA	N-CA-C	5.62	122.77	110.80
1	B	123	ILE	N-CA-C	5.40	114.58	106.42
1	C	64	VAL	CB-CA-C	-5.37	102.65	110.81

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	1769	0	1717	26	0
1	B	1734	0	1680	21	0
1	C	1718	0	1618	26	0
1	D	1721	0	1668	25	0
2	A	12	0	12	1	0
2	B	12	0	12	2	0
2	C	12	0	12	2	0
2	D	12	0	12	1	0
3	A	11	0	0	1	0
3	B	3	0	0	0	0
3	D	7	0	0	0	0
All	All	7011	0	6731	95	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 7.

All (95) 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:31:ALA:O	1:A:38:PHE:HA	1.88	0.74
1:B:160:PHE:CZ	1:B:169:GLU:HG3	2.23	0.73
1:A:203:PRO:O	1:A:204:PHE:CD2	2.43	0.72
1:D:160:PHE:CZ	1:D:169:GLU:HG3	2.26	0.71
1:B:234:ASP:OD1	1:B:236:THR:OG1	2.10	0.70
1:D:234:ASP:OD1	1:D:236:THR:OG1	2.10	0.69
1:C:160:PHE:CZ	1:C:169:GLU:HG3	2.27	0.69
1:C:234:ASP:OD1	1:C:236:THR:OG1	2.09	0.68
1:A:234:ASP:OD1	1:A:236:THR:OG1	2.11	0.68
1:B:34:LYS:HZ1	1:D:219:GLU:CD	2.03	0.66
1:A:38:PHE:HB2	3:A:406:HOH:O	1.96	0.65
1:A:203:PRO:O	1:A:204:PHE:CG	2.50	0.64
1:B:245:PHE:O	1:B:246:LYS:C	2.41	0.63
1:C:114:LEU:HG	1:C:197:VAL:CG1	2.28	0.63
1:C:91:THR:HG1	2:C:301:ARG:N	1.97	0.63
1:A:245:PHE:O	1:A:246:LYS:C	2.43	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:202:GLU:CB	1:C:203:PRO:CD	2.79	0.60
1:C:51:ASP:CB	1:C:240:LEU:HD21	2.31	0.59
1:B:51:ASP:CB	1:B:240:LEU:HD21	2.33	0.59
1:C:75:ILE:HD11	1:C:90:MET:HE1	1.85	0.58
1:A:34:LYS:HD2	1:A:187:TYR:CE1	2.39	0.58
1:B:75:ILE:HD11	1:B:90:MET:HE1	1.85	0.57
1:B:108:THR:HG22	1:B:206:HIS:ND1	2.19	0.57
1:D:75:ILE:HD11	1:D:90:MET:HE1	1.87	0.57
1:D:60:LEU:HD13	1:D:223:TRP:NE1	2.20	0.56
1:A:60:LEU:HD13	1:A:223:TRP:NE1	2.20	0.56
1:B:60:LEU:HD13	1:B:223:TRP:NE1	2.20	0.56
1:A:75:ILE:HD11	1:A:90:MET:HE1	1.88	0.56
1:C:60:LEU:HD13	1:C:223:TRP:NE1	2.22	0.55
1:C:57:ALA:CB	1:C:64:VAL:HG22	2.36	0.55
1:D:34:LYS:HD3	1:D:187:TYR:OH	2.07	0.54
1:B:51:ASP:HB2	1:B:240:LEU:HD21	1.90	0.54
1:C:123:ILE:HD12	1:C:129:LEU:CD1	2.38	0.53
1:B:27:VAL:HG22	1:B:87:MET:HE3	1.91	0.53
1:D:189:ALA:O	1:D:190:LYS:CB	2.57	0.52
1:A:57:ALA:CB	1:A:64:VAL:HG22	2.39	0.52
1:D:117:LYS:HA	1:D:198:VAL:HG11	1.91	0.52
1:D:27:VAL:HG22	1:D:87:MET:HE3	1.91	0.52
1:D:51:ASP:HB3	1:D:240:LEU:HD21	1.91	0.52
1:C:27:VAL:HG22	1:C:87:MET:HE3	1.91	0.52
1:C:51:ASP:HB2	1:C:240:LEU:HD21	1.90	0.52
1:B:91:THR:HG1	2:B:301:ARG:N	2.09	0.51
1:A:34:LYS:HG2	1:A:38:PHE:CG	2.44	0.51
1:B:34:LYS:NZ	1:D:219:GLU:CD	2.68	0.51
1:D:182:SER:OG	1:D:183:PRO:HD3	2.10	0.51
1:A:154:LYS:O	1:D:161:GLU:OE2	2.28	0.50
1:D:91:THR:HG1	2:D:301:ARG:N	2.10	0.50
1:C:245:PHE:O	1:C:246:LYS:C	2.55	0.49
1:C:90:MET:HE2	1:C:102:PHE:HZ	1.76	0.49
1:A:90:MET:HE2	1:A:102:PHE:HZ	1.78	0.49
1:A:204:PHE:O	1:A:205:THR:OG1	2.30	0.49
1:A:51:ASP:HB3	1:A:240:LEU:HD21	1.95	0.48
1:D:90:MET:HE2	1:D:102:PHE:HZ	1.78	0.48
1:A:107:MET:HG2	1:A:209:LEU:HD12	1.94	0.48
1:B:112:THR:CG2	1:B:199:VAL:HG13	2.44	0.48
1:A:34:LYS:HG2	1:A:38:PHE:CD1	2.49	0.48
1:B:90:MET:HE2	1:B:102:PHE:HZ	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ALA:O	1:A:190:LYS:CB	2.61	0.47
1:D:107:MET:HG2	1:D:209:LEU:HD12	1.96	0.47
1:D:16:LEU:C	1:D:16:LEU:HD13	2.40	0.47
1:D:112:THR:CG2	1:D:199:VAL:HG13	2.45	0.47
1:A:22:ARG:NH1	1:A:26:ARG:NH1	2.63	0.47
1:D:39:LYS:HB3	1:D:45:TYR:CE1	2.51	0.46
1:A:112:THR:CG2	1:A:199:VAL:HG13	2.46	0.46
1:C:202:GLU:CB	1:C:203:PRO:HD3	2.45	0.46
1:C:107:MET:HG2	1:C:209:LEU:HD12	1.98	0.46
1:C:112:THR:CG2	1:C:199:VAL:HG13	2.46	0.46
1:C:189:ALA:O	1:C:190:LYS:CB	2.64	0.45
1:B:108:THR:HG23	1:B:206:HIS:HB3	1.98	0.45
1:D:117:LYS:HA	1:D:198:VAL:CG1	2.46	0.45
1:D:113:ILE:HB	1:D:200:VAL:HG13	1.99	0.44
1:D:60:LEU:CD1	1:D:223:TRP:NE1	2.81	0.44
1:A:34:LYS:CG	1:A:38:PHE:CG	3.01	0.44
1:B:22:ARG:NH1	1:B:26:ARG:NH1	2.66	0.43
1:C:113:ILE:HB	1:C:200:VAL:HG13	1.99	0.43
1:B:113:ILE:HB	1:B:200:VAL:HG13	1.99	0.43
1:B:30:ASP:HB2	2:B:301:ARG:HH22	1.83	0.42
1:A:60:LEU:CD1	1:A:223:TRP:NE1	2.82	0.42
1:D:60:LEU:HB3	1:D:62:VAL:HG22	2.02	0.42
1:D:22:ARG:NH1	1:D:24:THR:O	2.53	0.42
1:A:181:ASP:OD2	2:A:301:ARG:N	2.53	0.42
1:A:90:MET:HE2	1:A:102:PHE:CZ	2.55	0.41
1:C:90:MET:HE2	1:C:102:PHE:CZ	2.54	0.41
1:C:187:TYR:CD1	1:C:187:TYR:C	2.98	0.41
1:B:162:ASN:OD1	1:B:162:ASN:C	2.64	0.41
1:C:114:LEU:HG	1:C:197:VAL:HG11	2.00	0.41
1:A:14:SER:OG	1:A:217:ASP:OD1	2.35	0.41
1:A:201:ASP:O	1:A:203:PRO:HD3	2.21	0.41
1:B:90:MET:HE2	1:B:102:PHE:CZ	2.56	0.41
1:C:75:ILE:HD11	1:C:90:MET:CE	2.51	0.41
1:C:234:ASP:OD1	1:C:234:ASP:C	2.64	0.41
1:C:60:LEU:HB3	1:C:62:VAL:HG22	2.02	0.41
1:B:60:LEU:HB3	1:B:62:VAL:HG22	2.03	0.40
1:C:143:THR:HG23	2:C:301:ARG:O	2.22	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	225/247 (91%)	211 (94%)	12 (5%)	2 (1%)	14	30
1	B	221/247 (90%)	211 (96%)	9 (4%)	1 (0%)	24	46
1	C	225/247 (91%)	213 (95%)	11 (5%)	1 (0%)	30	51
1	D	220/247 (89%)	211 (96%)	9 (4%)	0	100	100
All	All	891/988 (90%)	846 (95%)	41 (5%)	4 (0%)	30	51

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	ALA
1	B	196	ALA
1	C	202	GLU
1	A	204	PHE

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	184/209 (88%)	169 (92%)	15 (8%)	10	24
1	B	177/209 (85%)	165 (93%)	12 (7%)	14	32
1	C	171/209 (82%)	159 (93%)	12 (7%)	14	31
1	D	176/209 (84%)	165 (94%)	11 (6%)	16	36
All	All	708/836 (85%)	658 (93%)	50 (7%)	14	31

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ASP
1	A	41	LYS
1	A	60	LEU
1	A	126	PHE
1	A	127	GLU
1	A	131	LYS
1	A	146	GLN
1	A	158	ARG
1	A	182	SER
1	A	195	LEU
1	A	197	VAL
1	A	198	VAL
1	A	201	ASP
1	A	236	THR
1	A	240	LEU
1	B	52	LEU
1	B	60	LEU
1	B	80	THR
1	B	131	LYS
1	B	136	VAL
1	B	140	LEU
1	B	182	SER
1	B	195	LEU
1	B	200	VAL
1	B	205	THR
1	B	236	THR
1	B	240	LEU
1	C	52	LEU
1	C	60	LEU
1	C	127	GLU
1	C	129	LEU
1	C	136	VAL
1	C	140	LEU
1	C	182	SER
1	C	198	VAL
1	C	200	VAL
1	C	205	THR
1	C	236	THR
1	C	240	LEU
1	D	25	LEU
1	D	52	LEU
1	D	60	LEU

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Mol	Chain	Res	Type
1	D	127	GLU
1	D	136	VAL
1	D	140	LEU
1	D	161	GLU
1	D	200	VAL
1	D	205	THR
1	D	236	THR
1	D	240	LEU

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

Mol	Chain	Res	Type
1	C	146	GLN

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](#)

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	ARG	A	301	-	10,11,11	0.74	0	9,13,13	0.57	0
2	ARG	D	301	-	10,11,11	0.70	0	9,13,13	0.66	0
2	ARG	B	301	-	10,11,11	0.72	0	9,13,13	0.64	0
2	ARG	C	301	-	10,11,11	0.67	0	9,13,13	0.84	1 (11%)

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	ARG	A	301	-	-	1/11/11/11	-
2	ARG	D	301	-	-	0/11/11/11	-
2	ARG	B	301	-	-	0/11/11/11	-
2	ARG	C	301	-	-	1/11/11/11	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	ARG	OXT-C-O	-2.28	118.90	124.08

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	ARG	NE-CD-CG-CB
2	C	301	ARG	OXT-C-CA-N

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	ARG	1	0
2	D	301	ARG	1	0
2	B	301	ARG	2	0
2	C	301	ARG	2	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	229/247 (92%)	0.13	13 (5%) 29 24	18, 30, 58, 74	0
1	B	227/247 (91%)	0.41	12 (5%) 32 26	23, 42, 77, 111	0
1	C	229/247 (92%)	0.75	19 (8%) 17 13	34, 55, 90, 144	0
1	D	225/247 (91%)	0.39	11 (4%) 35 29	16, 40, 84, 102	1 (0%)
All	All	910/988 (92%)	0.42	55 (6%) 27 22	16, 42, 81, 144	1 (0%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	196	ALA	5.7
1	B	204	PHE	5.7
1	B	196	ALA	5.4
1	A	196	ALA	5.1
1	C	121	ASP	5.0
1	C	195	LEU	4.4
1	B	195	LEU	4.3
1	C	119	ASN	4.0
1	A	197	VAL	3.9
1	C	198	VAL	3.7
1	A	195	LEU	3.5
1	C	72	ASP	3.4
1	C	204	PHE	3.4
1	D	197	VAL	3.4
1	D	126	PHE	3.1
1	C	203	PRO	3.0
1	D	121	ASP	3.0
1	C	197	VAL	3.0
1	D	198	VAL	3.0
1	A	203	PRO	3.0
1	A	198	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	120	ALA	2.8
1	D	119	ASN	2.8
1	B	201	ASP	2.6
1	C	196	ALA	2.6
1	D	120	ALA	2.6
1	C	123	ILE	2.5
1	A	162	ASN	2.5
1	B	121	ASP	2.4
1	C	118	ASP	2.4
1	A	199	VAL	2.4
1	A	201	ASP	2.4
1	C	38	PHE	2.4
1	D	132	PRO	2.3
1	C	71	TRP	2.3
1	B	216	GLY	2.3
1	B	123	ILE	2.3
1	B	124	LYS	2.3
1	A	187	TYR	2.3
1	D	118	ASP	2.2
1	C	95	GLU	2.2
1	D	150	GLU	2.2
1	C	151	PHE	2.2
1	B	197	VAL	2.1
1	B	31	ALA	2.1
1	C	122	LYS	2.1
1	C	93	THR	2.1
1	C	218	PRO	2.1
1	A	200	VAL	2.0
1	A	164	ALA	2.0
1	A	121	ASP	2.0
1	B	118	ASP	2.0
1	D	151	PHE	2.0
1	A	163	ASN	2.0
1	B	205	THR	2.0

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ARG	C	301	12/12	0.92	0.10	41,43,45,47	0
2	ARG	B	301	12/12	0.95	0.07	20,23,26,29	0
2	ARG	D	301	12/12	0.97	0.06	17,17,18,19	0
2	ARG	A	301	12/12	0.99	0.04	7,7,9,9	0

6.5 Other polymers [i](#)

There are no such residues in this entry.