



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:22 PM UTC

PDB ID : 2DEW / pdb_00002dew
Title : Crystal structure of human peptidylarginine deiminase 4 in complex with histone H3 N-terminal tail including Arg8
Authors : Arita, K.; Shimizu, T.; Hashimoto, H.; Hidaka, Y.; Yamada, M.; Sato, M.
Deposited on : 2006-02-18
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

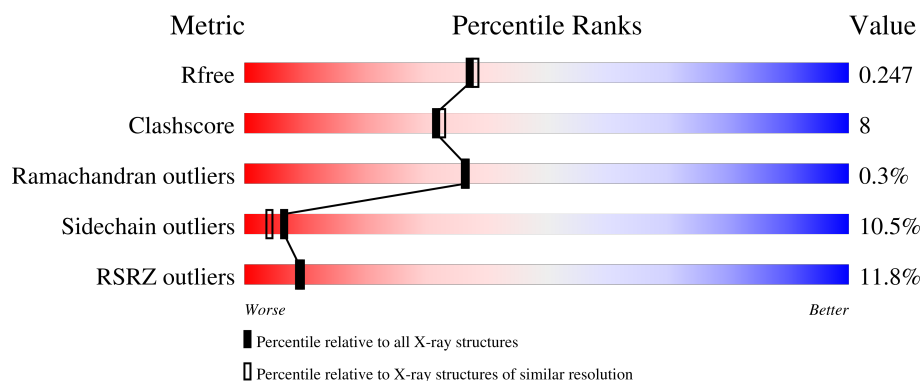
1 Overall quality at a glance

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	X	671	
2	A	10	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5234 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 Protein-arginine deiminase type IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	629	Total	C	N	O	S	0	0	0
			4952	3161	831	925	35			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	-7	GLY	-	cloning artifact	UNP Q9UM07
X	-6	PRO	-	cloning artifact	UNP Q9UM07
X	-5	LEU	-	cloning artifact	UNP Q9UM07
X	-4	GLY	-	cloning artifact	UNP Q9UM07
X	-3	SER	-	cloning artifact	UNP Q9UM07
X	-2	PRO	-	cloning artifact	UNP Q9UM07
X	-1	GLU	-	cloning artifact	UNP Q9UM07
X	0	PHE	-	cloning artifact	UNP Q9UM07
X	645	ALA	CYS	engineered mutation	UNP Q9UM07

- Molecule 2 is a protein called 10-mer peptide from histone H3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	A	5	Total	C	N	O	0	0	0
			38	22	9	7			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	X	5	Total	Ca	0	0
			5	5		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	X	1	Total	O	S	0	0
			5	4	1		
4	X	1	Total	O	S	0	0
			5	4	1		
4	X	1	Total	O	S	0	0
			5	4	1		

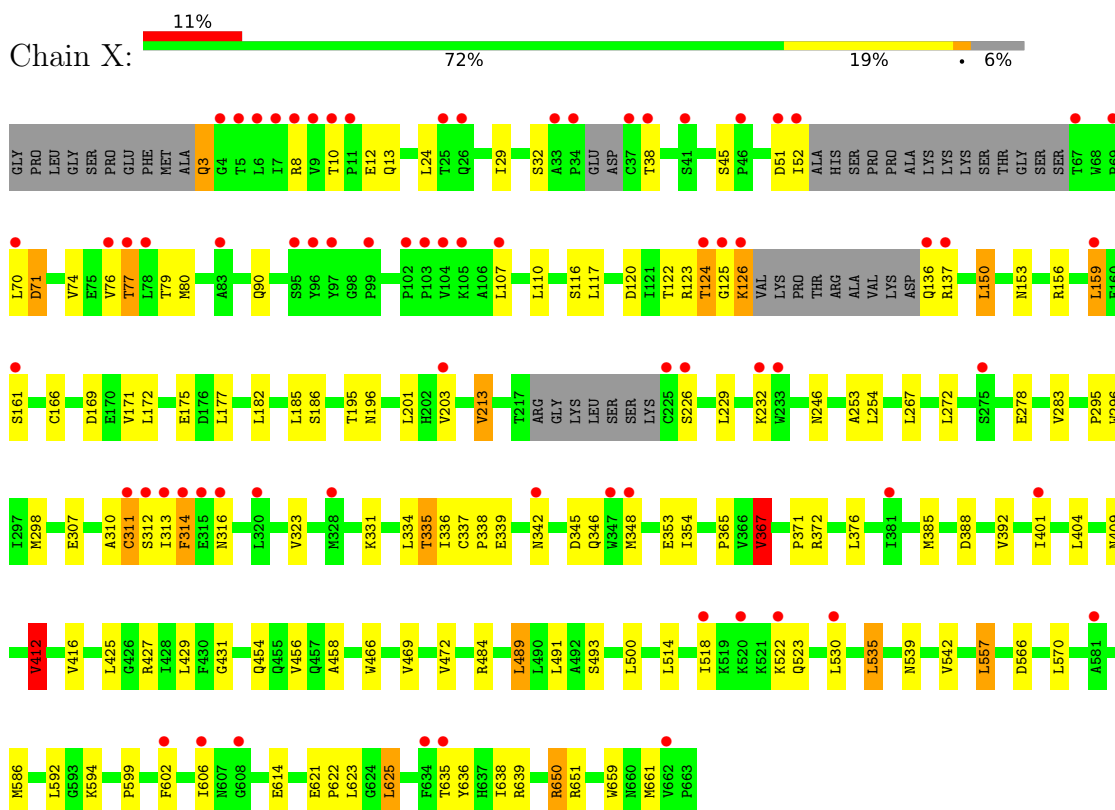
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	X	222	Total	O	0	0
			222	222		
5	A	2	Total	O	0	0
			2	2		

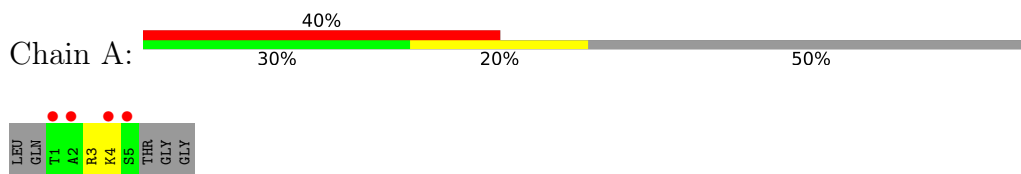
3 Residue-property plots [i](#)

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: Protein-arginine deiminase type IV



- Molecule 2: 10-mer peptide from histone H3



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	146.27Å 60.77Å 115.15Å 90.00° 124.26° 90.00°	Depositor
Resolution (Å)	38.34 – 2.10 38.34 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.9 (38.34-2.10) 97.9 (38.34-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.14 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.202 , 0.241 0.209 , 0.247	Depositor DCC
R_{free} test set	4878 reflections (10.16%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5234	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, SO4

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	X	0.72	0/5071	0.88	5/6881 (0.1%)
2	A	0.95	0/37	1.35	1/47 (2.1%)
All	All	0.72	0/5108	0.88	6/6928 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	367	VAL	CB-CA-C	-9.71	99.21	110.96
1	X	316	ASN	N-CA-C	8.22	122.31	108.90
1	X	412	VAL	CB-CA-C	-6.21	99.60	110.71
2	A	4	LYS	N-CA-C	6.19	117.82	111.14
1	X	314	PHE	N-CA-C	-5.64	102.00	108.49
1	X	311	CYS	N-CA-C	-5.37	99.62	108.49

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	X	4952	0	4913	81	0
2	A	38	0	45	1	0
3	X	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	X	15	0	0	0	0
5	A	2	0	0	0	0
5	X	222	0	0	4	1
All	All	5234	0	4958	81	1

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 8.

All (81) 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:X:651:ARG:NH1	5:X:1096:HOH:O	2.05	0.89
1:X:416:VAL:HG21	1:X:557:LEU:O	1.78	0.84
1:X:313:ILE:HB	1:X:338:PRO:HA	1.63	0.80
1:X:120:ASP:HB2	1:X:126:LYS:HD3	1.64	0.78
1:X:650:ARG:HH11	1:X:650:ARG:HG3	1.55	0.71
1:X:311:CYS:HB3	1:X:348:MET:HE2	1.79	0.64
1:X:203:VAL:HG12	1:X:267:LEU:HD23	1.80	0.64
1:X:310:ALA:O	1:X:336:ILE:HA	1.97	0.64
1:X:354:ILE:HD13	1:X:367:VAL:HG13	1.79	0.64
1:X:296:TRP:H	1:X:454:GLN:HE21	1.47	0.62
1:X:416:VAL:CG2	1:X:557:LEU:O	2.46	0.62
1:X:307:GLU:OE2	1:X:335:THR:CG2	2.48	0.61
1:X:195:THR:HG22	1:X:196:ASN:ND2	2.18	0.59
1:X:313:ILE:CB	1:X:338:PRO:HA	2.32	0.59
1:X:409:ASN:HD22	1:X:472:VAL:H	1.51	0.59
1:X:307:GLU:OE2	1:X:335:THR:HG21	2.05	0.57
1:X:586:MET:HA	1:X:599:PRO:HG2	1.86	0.57
1:X:334:LEU:C	1:X:334:LEU:HD13	2.30	0.56
1:X:203:VAL:CG1	1:X:267:LEU:CD2	2.84	0.56
1:X:123:ARG:HD3	1:X:659:TRP:CD1	2.41	0.55
1:X:334:LEU:HD11	1:X:336:ILE:CD1	2.36	0.55
1:X:466:TRP:CZ3	1:X:542:VAL:HG13	2.42	0.54
1:X:311:CYS:SG	1:X:311:CYS:O	2.66	0.53
1:X:10:THR:HG22	1:X:32:SER:CB	2.38	0.53
1:X:153:ASN:HB3	1:X:166:CYS:HB3	1.91	0.53
1:X:296:TRP:H	1:X:454:GLN:NE2	2.06	0.53
1:X:71:ASP:C	1:X:71:ASP:OD1	2.51	0.53
1:X:409:ASN:HD21	1:X:431:GLY:HA3	1.74	0.52
1:X:51:ASP:HB2	1:X:77:THR:HG23	1.90	0.52
1:X:52:ILE:HD12	1:X:76:VAL:HG22	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:365:PRO:HD2	1:X:388:ASP:O	2.10	0.51
1:X:45:SER:OG	1:X:90:GLN:NE2	2.44	0.51
1:X:232:LYS:O	1:X:232:LYS:HD3	2.11	0.50
1:X:650:ARG:HH11	1:X:650:ARG:CG	2.22	0.50
1:X:311:CYS:CB	1:X:337:CYS:HB3	2.42	0.50
1:X:425:LEU:HD12	1:X:456:VAL:HG22	1.94	0.50
1:X:298:MET:HE3	1:X:353:GLU:CD	2.37	0.50
1:X:213:VAL:HG13	1:X:229:LEU:HB2	1.93	0.49
1:X:311:CYS:HB2	1:X:342:ASN:OD1	2.13	0.49
1:X:203:VAL:HG12	1:X:267:LEU:CD2	2.43	0.49
1:X:29:ILE:CD1	1:X:76:VAL:HG21	2.43	0.48
1:X:313:ILE:HG12	1:X:336:ILE:CG2	2.43	0.48
1:X:298:MET:HE2	1:X:412:VAL:HG22	1.95	0.48
1:X:635:THR:HG23	1:X:636:TYR:HD1	1.77	0.48
1:X:166:CYS:HB2	1:X:254:LEU:HD22	1.95	0.48
1:X:416:VAL:HG23	5:X:1107:HOH:O	2.14	0.48
1:X:518:ILE:HD11	1:X:602:PHE:CD1	2.48	0.48
1:X:226:SER:HB2	5:X:1025:HOH:O	2.13	0.48
1:X:29:ILE:HD12	1:X:76:VAL:HG21	1.96	0.47
1:X:124:THR:OG1	1:X:125:GLY:N	2.45	0.47
1:X:150:LEU:HD13	1:X:253:ALA:HB2	1.97	0.47
1:X:518:ILE:HD11	1:X:602:PHE:CG	2.49	0.46
1:X:311:CYS:HB3	1:X:337:CYS:HB3	1.98	0.46
1:X:345:ASP:OD2	1:X:372:ARG:NH2	2.50	0.45
1:X:623:LEU:HB2	1:X:625:LEU:CD2	2.46	0.45
1:X:10:THR:HG22	1:X:32:SER:HB2	1.99	0.45
1:X:296:TRP:N	1:X:454:GLN:HE21	2.11	0.45
1:X:456:VAL:HG23	1:X:661:MET:SD	2.56	0.44
1:X:650:ARG:HD3	5:X:1100:HOH:O	2.17	0.44
1:X:469:VAL:HG11	2:A:3:ARG:HD3	1.98	0.44
1:X:535:LEU:HD22	1:X:539:ASN:ND2	2.33	0.44
1:X:74:VAL:HG13	1:X:74:VAL:O	2.18	0.43
1:X:295:PRO:HA	1:X:454:GLN:NE2	2.34	0.43
1:X:156:ARG:HD2	1:X:161:SER:O	2.19	0.43
1:X:3:GLN:HE21	1:X:24:LEU:H	1.67	0.43
1:X:171:VAL:HG11	1:X:177:LEU:HD21	2.01	0.43
1:X:126:LYS:HE3	1:X:182:LEU:HB3	2.01	0.42
1:X:354:ILE:CD1	1:X:367:VAL:HG13	2.48	0.42
1:X:371:PRO:O	1:X:372:ARG:C	2.60	0.42
1:X:489:LEU:HD13	1:X:491:LEU:HG	2.01	0.42
1:X:122:THR:OG1	1:X:124:THR:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:186:SER:OG	1:X:246:ASN:ND2	2.53	0.41
1:X:296:TRP:CZ3	1:X:298:MET:HG3	2.54	0.41
1:X:493:SER:CB	1:X:566:ASP:HB3	2.50	0.41
1:X:621:GLU:N	1:X:622:PRO:CD	2.83	0.41
1:X:71:ASP:O	1:X:74:VAL:HG12	2.20	0.41
1:X:456:VAL:CG2	1:X:661:MET:SD	3.09	0.41
1:X:80:MET:HE3	1:X:90:GLN:HG3	2.03	0.41
1:X:339:GLU:HA	1:X:342:ASN:HB2	2.03	0.41
1:X:272:LEU:CD2	1:X:283:VAL:HG22	2.51	0.40
1:X:296:TRP:CH2	1:X:298:MET:CG	3.05	0.40

All (1) 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:X:931:HOH:O	5:X:993:HOH:O[2_555]	2.17	0.03

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	X	619/671 (92%)	587 (95%)	30 (5%)	2 (0%)	36	36
2	A	3/10 (30%)	3 (100%)	0	0	100	100
All	All	622/681 (91%)	590 (95%)	30 (5%)	2 (0%)	36	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	159	LEU
1	X	458	ALA

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	X	559/593 (94%)	500 (89%)	59 (11%)	6	4
2	A	4/7 (57%)	4 (100%)	0	100	100
All	All	563/600 (94%)	504 (90%)	59 (10%)	6	4

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	3	GLN
1	X	8	ARG
1	X	12	GLU
1	X	13	GLN
1	X	38	THR
1	X	70	LEU
1	X	71	ASP
1	X	77	THR
1	X	79	THR
1	X	107	LEU
1	X	110	LEU
1	X	116	SER
1	X	117	LEU
1	X	124	THR
1	X	126	LYS
1	X	136	GLN
1	X	137	ARG
1	X	150	LEU
1	X	159	LEU
1	X	169	ASP
1	X	172	LEU
1	X	175	GLU
1	X	185	LEU
1	X	201	LEU
1	X	213	VAL
1	X	278	GLU
1	X	312	SER

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Mol	Chain	Res	Type
1	X	314	PHE
1	X	323	VAL
1	X	331	LYS
1	X	335	THR
1	X	346	GLN
1	X	367	VAL
1	X	376	LEU
1	X	385	MET
1	X	392	VAL
1	X	401	ILE
1	X	404	LEU
1	X	412	VAL
1	X	427	ARG
1	X	429	LEU
1	X	484	ARG
1	X	489	LEU
1	X	500	LEU
1	X	514	LEU
1	X	522	LYS
1	X	523	GLN
1	X	530	LEU
1	X	535	LEU
1	X	557	LEU
1	X	570	LEU
1	X	592	LEU
1	X	594	LYS
1	X	606	ILE
1	X	614	GLU
1	X	625	LEU
1	X	638	ILE
1	X	639	ARG
1	X	650	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	3	GLN
1	X	13	GLN
1	X	90	GLN
1	X	178	GLN
1	X	196	ASN
1	X	246	ASN

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Mol	Chain	Res	Type
1	X	409	ASN
1	X	438	ASN
1	X	448	GLN
1	X	454	GLN
1	X	538	HIS
1	X	640	HIS

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 8 ligands modelled in this entry, 5 are monoatomic - leaving 3 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	SO4	X	905	-	4,4,4	0.23	0	6,6,6	0.21	0
4	SO4	X	906	-	4,4,4	0.20	0	6,6,6	0.15	0
4	SO4	X	907	-	4,4,4	0.26	0	6,6,6	0.13	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	X	629/671 (93%)	0.92	71 (11%) 10 10	40, 50, 67, 80	0
2	A	5/10 (50%)	2.37	4 (80%) 0 0	58, 61, 70, 71	0
All	All	634/681 (93%)	0.93	75 (11%) 9 9	40, 50, 67, 80	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	312	SER	5.7
1	X	313	ILE	5.3
1	X	314	PHE	4.3
1	X	518	ILE	4.1
1	X	6	LEU	4.0
1	X	159	LEU	3.9
1	X	275	SER	3.9
1	X	76	VAL	3.8
1	X	126	LYS	3.8
1	X	52	ILE	3.7
1	X	315	GLU	3.7
1	X	70	LEU	3.6
1	X	34	PRO	3.5
1	X	11	PRO	3.4
1	X	320	LEU	3.4
2	A	5	SER	3.3
1	X	9	VAL	3.3
2	A	4	LYS	3.2
1	X	316	ASN	3.1
1	X	342	ASN	3.1
1	X	33	ALA	3.1
1	X	41	SER	3.1
2	A	2	ALA	3.0
1	X	83	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
1	X	99	PRO	3.0
1	X	124	THR	2.9
1	X	125	GLY	2.8
1	X	8	ARG	2.7
1	X	233	TRP	2.8
1	X	635	THR	2.7
1	X	137	ARG	2.7
1	X	5	THR	2.7
1	X	107	LEU	2.7
1	X	161	SER	2.7
1	X	38	THR	2.6
1	X	4	GLY	2.6
1	X	69	PRO	2.6
1	X	608	GLY	2.6
1	X	78	LEU	2.6
1	X	530	LEU	2.6
1	X	37	CYS	2.6
1	X	662	VAL	2.5
1	X	97	TYR	2.5
1	X	232	LYS	2.5
1	X	225	CYS	2.5
1	X	25	THR	2.5
2	A	1	THR	2.5
1	X	10	THR	2.4
1	X	67	THR	2.4
1	X	602	PHE	2.4
1	X	381	ILE	2.4
1	X	328	MET	2.4
1	X	520	LYS	2.4
1	X	311	CYS	2.4
1	X	7	ILE	2.3
1	X	77	THR	2.3
1	X	95	SER	2.2
1	X	105	LYS	2.2
1	X	136	GLN	2.2
1	X	522	LYS	2.2
1	X	203	VAL	2.2
1	X	103	PRO	2.1
1	X	634	PHE	2.1
1	X	104	VAL	2.1
1	X	348	MET	2.1
1	X	51	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	X	96	TYR	2.1
1	X	581	ALA	2.1
1	X	606	ILE	2.1
1	X	26	GLN	2.0
1	X	102	PRO	2.0
1	X	401	ILE	2.0
1	X	226	SER	2.0
1	X	347	TRP	2.0
1	X	46	PRO	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 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
4	SO4	X	907	5/5	0.67	0.14	75,75,75,76	0
4	SO4	X	905	5/5	0.78	0.15	90,90,91,91	0
4	SO4	X	906	5/5	0.83	0.10	86,87,87,87	0
3	CA	X	904	1/1	0.96	0.17	44,44,44,44	0
3	CA	X	902	1/1	0.97	0.13	53,53,53,53	0
3	CA	X	901	1/1	0.98	0.15	47,47,47,47	0
3	CA	X	903	1/1	0.98	0.14	50,50,50,50	0
3	CA	X	900	1/1	0.99	0.12	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.