



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 09:18 PM UTC

PDB ID : 3DW8 / pdb_00003dw8
Title : Structure of a Protein Phosphatase 2A Holoenzyme with B55 subunit
Authors : Xu, Y.; Chen, Y.; Zhang, P.; Jeffrey, P.D.; Shi, Y.
Deposited on : 2008-07-21
Resolution : 2.85 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

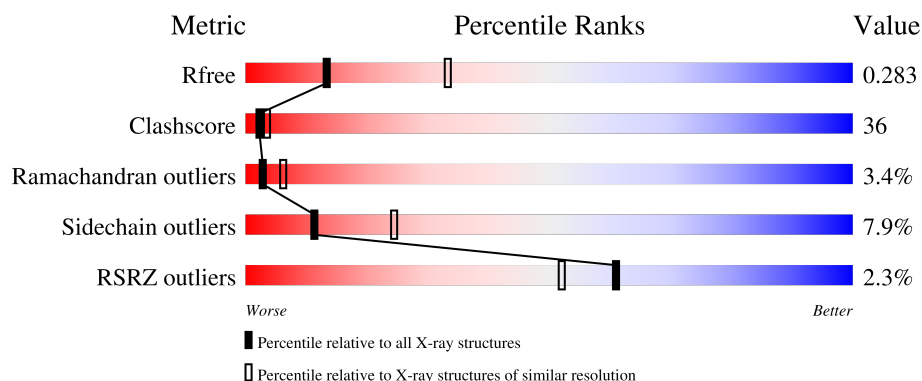
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

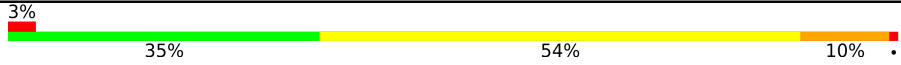
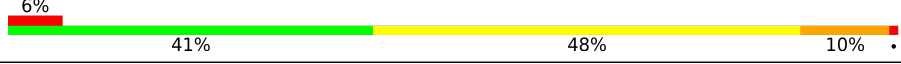
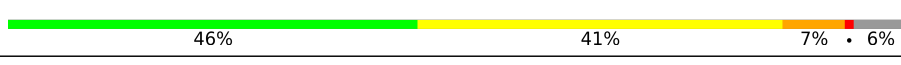
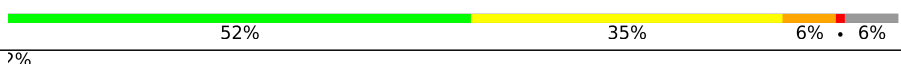
The reported resolution of this entry is 2.85 Å.

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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	582	
1	D	582	
2	B	447	
2	E	447	
3	C	309	

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Mol	Chain	Length	Quality of chain
3	F	309	% 36%50%7%7%
4	G	7	43%14%43%
4	H	7	43%29%14%14%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 20717 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 Serine/threonine-protein phosphatase 2A 65 kDa regulatory subunit A alpha isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	582	Total	C	N	O	S	0	0	0
			4535	2882	764	861	28			
1	D	582	Total	C	N	O	S	0	0	0
			4535	2882	764	861	28			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	MET	-	expression tag	UNP P30153
D	8	MET	-	expression tag	UNP P30153

- Molecule 2 is a protein called Serine/threonine-protein phosphatase 2A 55 kDa regulatory subunit B alpha isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	421	Total	C	N	O	S	0	0	0
			3429	2158	595	658	18			
2	E	421	Total	C	N	O	S	0	0	0
			3428	2157	595	658	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	310	VAL	ILE	engineered mutation	UNP P63151
E	310	VAL	ILE	engineered mutation	UNP P63151

- Molecule 3 is a protein called Serine/threonine-protein phosphatase 2A catalytic subunit alpha isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	288	Total	C	N	O	S	0	0	0
			2322	1471	396	440	15			
3	F	288	Total	C	N	O	S	0	0	0
			2322	1471	396	440	15			

- Molecule 4 is a protein called microcystin LR.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	7	Total	C	N	O	0	0	0
			71	49	10	12			
4	H	7	Total	C	N	O	0	0	0
			71	49	10	12			

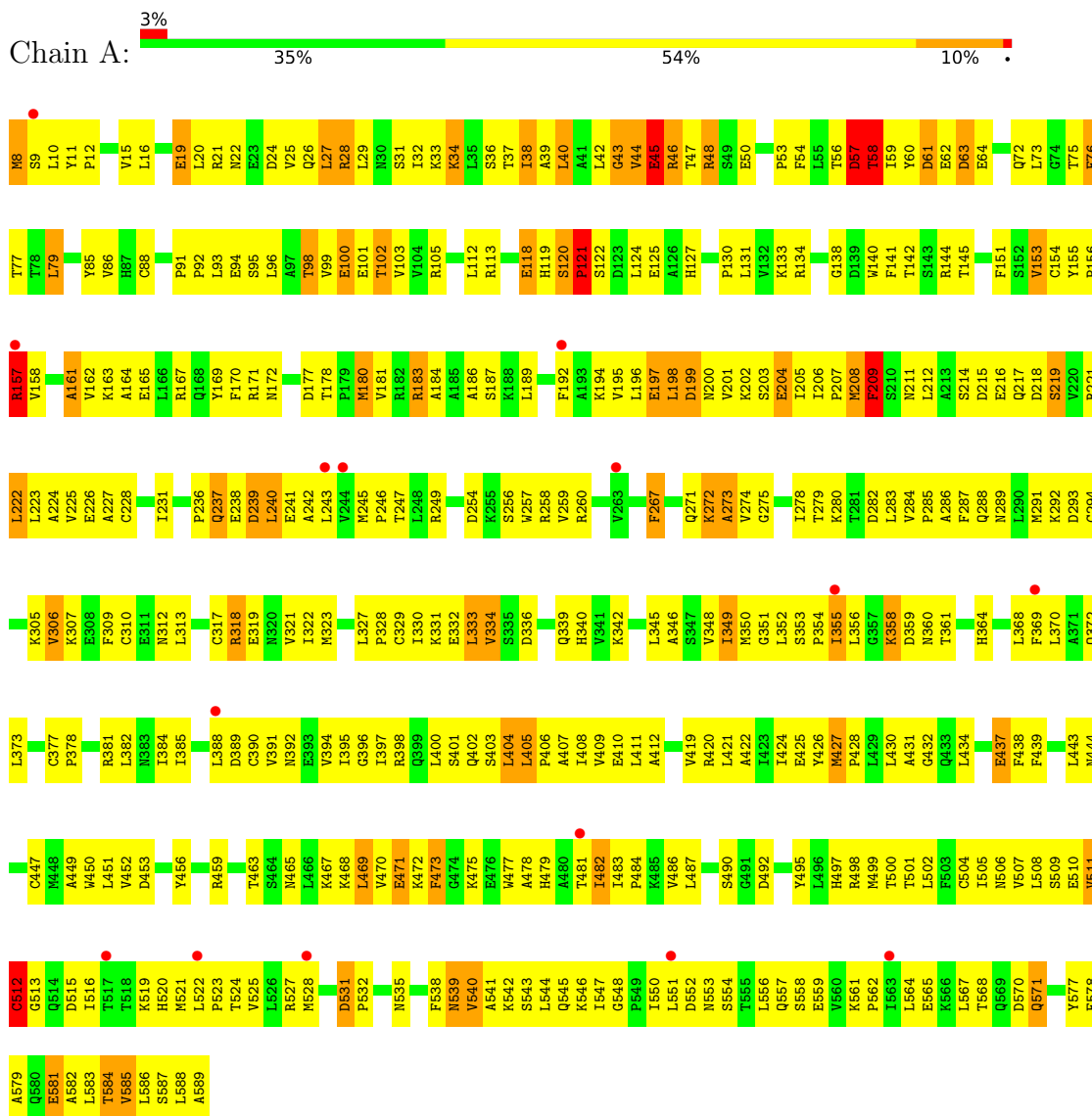
- Molecule 5 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	2	Total	Mn	0	0
			2	2		
5	F	2	Total	Mn	0	0
			2	2		

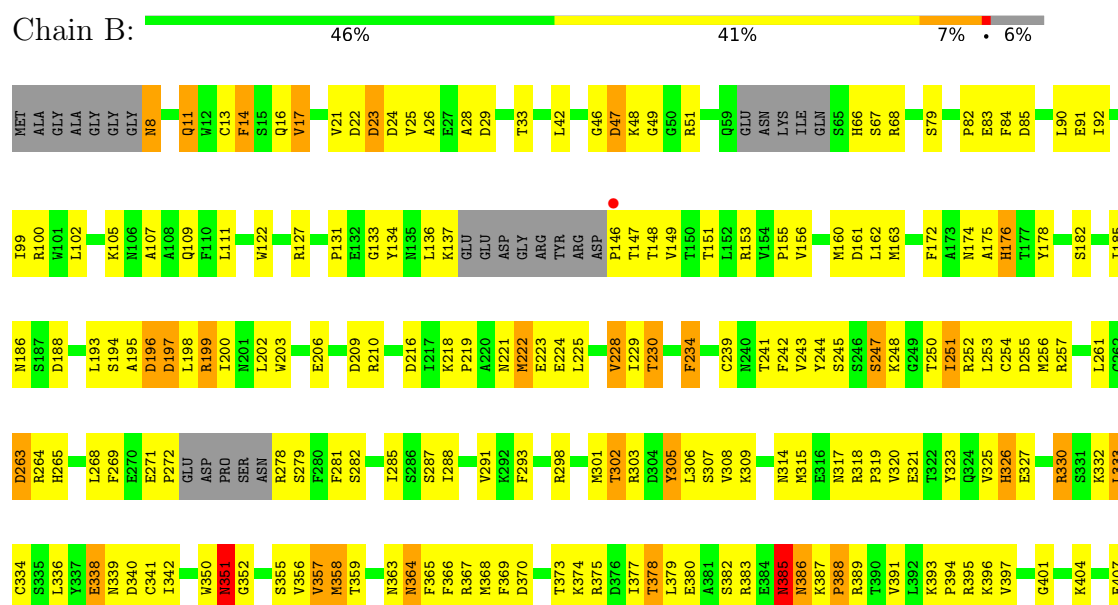
3 Residue-property plots

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.

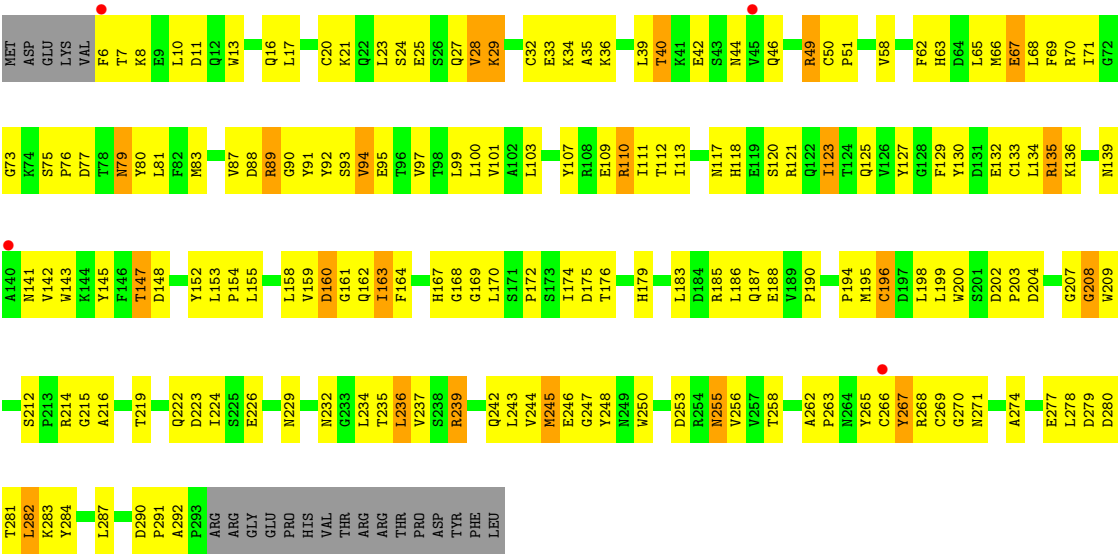
- Molecule 1: Serine/threonine-protein phosphatase 2A 65 kDa regulatory subunit A alpha isoform



- Molecule 1: Serine/threonine-protein phosphatase 2A 65 kDa regulatory subunit A alpha isoform



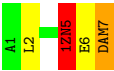




● Molecule 4: microcystin LR



● Molecule 4: microcystin LR



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	247.34Å 121.39Å 172.48Å 90.00° 132.60° 90.00°	Depositor
Resolution (Å)	49.63 – 2.85 49.63 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.63-2.85) 99.2 (49.63-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.40 (at 2.91Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.228 , 0.285 0.227 , 0.283	Depositor DCC
R_{free} test set	4180 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	78.6	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 74.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.010 for -h-2*1,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	20717	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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, 0.333 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: ACB, DAM, FGA, MN, 1ZN, DAL

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	0.54	1/4609 (0.0%)	1.03	23/6256 (0.4%)
1	D	0.56	1/4609 (0.0%)	1.04	35/6256 (0.6%)
2	B	0.65	0/3501	1.08	22/4730 (0.5%)
2	E	0.69	0/3500	1.11	21/4728 (0.4%)
3	C	0.49	0/2379	0.98	8/3227 (0.2%)
3	F	0.45	0/2379	0.98	9/3227 (0.3%)
4	G	0.44	0/17	0.57	0/19
4	H	0.41	0/17	0.64	0/19
All	All	0.58	2/21011 (0.0%)	1.04	118/28462 (0.4%)

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
4	G	0	2
4	H	0	2
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	208	MET	SD-CE	8.85	2.01	1.79
1	D	180	MET	SD-CE	6.35	1.95	1.79

The worst 5 of 118 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	48	LYS	N-CA-C	-10.59	99.75	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	48	LYS	N-CA-C	-9.81	100.72	112.89
1	A	47	THR	N-CA-C	-9.75	100.61	111.14
1	D	540	VAL	N-CA-C	-9.30	102.84	111.67
1	D	45	GLU	N-CA-C	-8.06	103.04	113.12

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	G	5	1ZN	Mainchain,Peptide
4	H	5	1ZN	Mainchain,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	4535	0	4642	436	0
1	D	4535	0	4642	395	0
2	B	3429	0	3323	220	0
2	E	3428	0	3322	190	0
3	C	2322	0	2223	116	0
3	F	2322	0	2223	172	0
4	G	71	0	61	4	0
4	H	71	0	61	2	0
5	C	2	0	0	0	0
5	F	2	0	0	0	0
All	All	20717	0	20497	1486	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 36.

The worst 5 of 1486 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:208:MET:SD	1:A:208:MET:CE	2.01	1.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:38:ILE:HG12	3:C:108:ARG:HH21	1.10	1.11
1:A:478:ALA:HA	1:A:482:ILE:HD12	1.30	1.10
1:D:22:ASN:ND2	1:D:27:LEU:HD12	1.69	1.08
1:A:373:LEU:HD21	1:A:404:LEU:HB3	1.39	1.04

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	580/582 (100%)	441 (76%)	110 (19%)	29 (5%)	1	2
1	D	580/582 (100%)	465 (80%)	88 (15%)	27 (5%)	2	3
2	B	413/447 (92%)	368 (89%)	38 (9%)	7 (2%)	7	16
2	E	413/447 (92%)	364 (88%)	41 (10%)	8 (2%)	6	14
3	C	286/309 (93%)	246 (86%)	35 (12%)	5 (2%)	7	16
3	F	286/309 (93%)	234 (82%)	40 (14%)	12 (4%)	2	3
4	G	1/7 (14%)	1 (100%)	0	0	100	100
4	H	1/7 (14%)	1 (100%)	0	0	100	100
All	All	2560/2690 (95%)	2120 (83%)	352 (14%)	88 (3%)	3	6

5 of 88 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	58	THR
1	A	157	ARG
1	A	241	GLU
1	A	318	ARG
1	A	558	SER

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	509/509 (100%)	470 (92%)	39 (8%)	12	25
1	D	509/509 (100%)	468 (92%)	41 (8%)	11	23
2	B	386/408 (95%)	350 (91%)	36 (9%)	8	18
2	E	386/408 (95%)	355 (92%)	31 (8%)	11	24
3	C	254/274 (93%)	235 (92%)	19 (8%)	12	26
3	F	254/274 (93%)	238 (94%)	16 (6%)	16	34
4	G	2/2 (100%)	2 (100%)	0	100	100
4	H	2/2 (100%)	2 (100%)	0	100	100
All	All	2302/2386 (96%)	2120 (92%)	182 (8%)	11	24

5 of 182 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	204	GLU
2	E	169	ARG
1	D	222	LEU
1	D	476	GLU
2	E	303	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
2	E	438	ASN
2	E	443	GLN
3	F	125	GLN
2	B	438	ASN
2	B	351	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

10 non-standard protein/DNA/RNA residues 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
4	1ZN	H	5	4	21,23,24	1.46	3 (14%)	25,29,31	1.10	1 (4%)
4	FGA	G	6	4	6,8,9	2.60	2 (33%)	6,9,11	0.50	0
4	ACB	G	3	4	7,8,9	1.12	0	8,10,12	0.75	0
4	DAM	H	7	3,4	4,5,6	2.50	2 (50%)	3,5,7	3.89	3 (100%)
4	1ZN	G	5	4	21,23,24	1.38	3 (14%)	25,29,31	1.10	2 (8%)
4	FGA	H	6	4	6,8,9	2.57	2 (33%)	6,9,11	0.54	0
4	DAM	G	7	3,4	4,5,6	2.02	1 (25%)	3,5,7	3.85	3 (100%)
4	ACB	H	3	4	7,8,9	1.10	0	8,10,12	1.01	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	1ZN	H	5	4	-	5/23/25/27	0/1/1/1
4	FGA	G	6	4	-	4/8/8/9	-
4	ACB	G	3	4	-	2/10/10/12	-
4	DAM	H	7	3,4	-	0/0/4/6	-
4	1ZN	G	5	4	-	5/23/25/27	0/1/1/1
4	FGA	H	6	4	-	4/8/8/9	-
4	DAM	G	7	3,4	-	0/0/4/6	-
4	ACB	H	3	4	-	2/10/10/12	-

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	6	FGA	O-C	5.35	1.38	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	6	FGA	O-C	5.31	1.37	1.22
4	H	7	DAM	C-CA	4.46	1.52	1.45
4	H	5	1ZN	C3-C2	3.77	1.57	1.52
4	G	7	DAM	C-CA	3.45	1.50	1.45

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	7	DAM	CM-N-CA	-4.78	116.72	123.98
4	G	7	DAM	CM-N-CA	-4.69	116.86	123.98
4	H	7	DAM	O-C-CA	-4.22	119.84	125.33
4	G	7	DAM	O-C-CA	-4.16	119.92	125.33
4	G	5	1ZN	CA-C18-C	-3.34	106.43	110.67

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	3	ACB	C4-CB-CG-OD1
4	H	3	ACB	C4-CB-CG-OD1
4	G	5	1ZN	C12-C13-C15-C16
4	H	5	1ZN	C12-C13-C15-C16
4	H	5	1ZN	C14-C13-C15-C16

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	5	1ZN	1	0
4	G	6	FGA	1	0
4	H	7	DAM	1	0
4	G	7	DAM	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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 [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	582/582 (100%)	0.41	15 (2%) 57 48	53, 104, 164, 173	0
1	D	582/582 (100%)	0.42	33 (5%) 29 22	48, 100, 173, 188	0
2	B	421/447 (94%)	-0.08	1 (0%) 91 91	38, 67, 115, 131	0
2	E	421/447 (94%)	-0.21	2 (0%) 87 85	34, 60, 111, 133	0
3	C	288/309 (93%)	0.14	5 (1%) 69 63	60, 88, 119, 154	0
3	F	288/309 (93%)	0.26	4 (1%) 73 67	75, 99, 129, 157	0
4	G	2/7 (28%)	0.79	0 100 100	95, 95, 95, 117	0
4	H	2/7 (28%)	1.10	0 100 100	108, 108, 108, 117	0
All	All	2586/2690 (96%)	0.19	60 (2%) 61 52	34, 88, 161, 188	0

The worst 5 of 60 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	400	LEU	4.3
3	C	6	PHE	3.7
1	D	368	LEU	3.7
2	B	146	PRO	3.4
1	D	404	LEU	3.4

6.2 Non-standard residues in protein, DNA, RNA chains [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	DAL	H	1	5/6	0.81	0.16	108,108,110,112	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	FGA	H	6	9/10	0.86	0.17	102,111,118,118	0
4	DAM	G	7	6/7	0.86	0.15	99,101,102,104	0
4	ACB	H	3	9/10	0.87	0.15	108,111,113,113	0
4	FGA	G	6	9/10	0.89	0.18	100,101,106,107	0
4	1ZN	H	5	23/24	0.91	0.21	106,109,113,115	0
4	DAL	G	1	5/6	0.92	0.12	92,94,96,98	0
4	ACB	G	3	9/10	0.93	0.12	93,95,102,104	0
4	1ZN	G	5	23/24	0.94	0.19	99,102,106,107	0
4	DAM	H	7	6/7	0.94	0.11	112,113,116,116	0

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
5	MN	C	501	1/1	0.98	0.03	69,69,69,69	0
5	MN	C	502	1/1	0.98	0.03	69,69,69,69	0
5	MN	F	511	1/1	0.98	0.05	82,82,82,82	0
5	MN	F	512	1/1	0.98	0.04	91,91,91,91	0

6.5 Other polymers [i](#)

There are no such residues in this entry.