



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:10 PM GMT

PDB ID : 1SDY
Title : STRUCTURE SOLUTION AND MOLECULAR DYNAMICS REFINEMENT OF THE YEAST CU,ZN ENZYME SUPEROXIDE DISMUTASE
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Deposited on : 1991-06-14
Resolution : 2.50 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

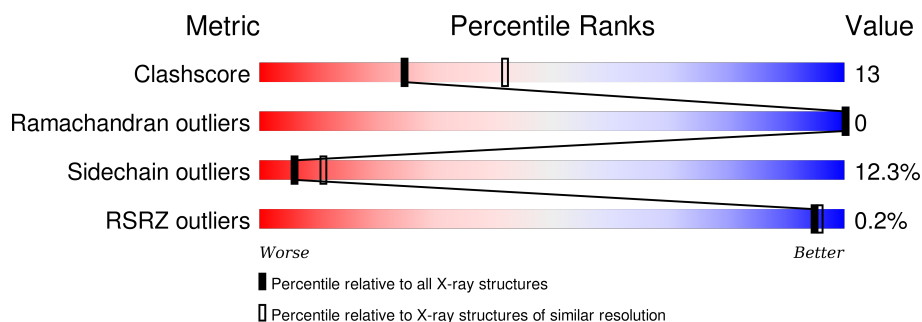
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	153	 66% 27% 5% •
1	B	153	 67% 27% 5%
1	C	153	 65% 32% •
1	D	153	 69% 24% 7% •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4945 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 COPPER,ZINC SUPEROXIDE DISMUTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	21	0	0
			1106	680	198	225	3			
1	B	153	Total	C	N	O	S	18	0	0
			1106	680	198	225	3			
1	C	153	Total	C	N	O	S	18	0	0
			1106	680	198	225	3			
1	D	153	Total	C	N	O	S	13	0	0
			1106	680	198	225	3			

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cu	0	0
			1	1		
2	A	1	Total	Cu	0	0
			1	1		
2	D	1	Total	Cu	0	0
			1	1		
2	C	1	Total	Cu	0	0
			1	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

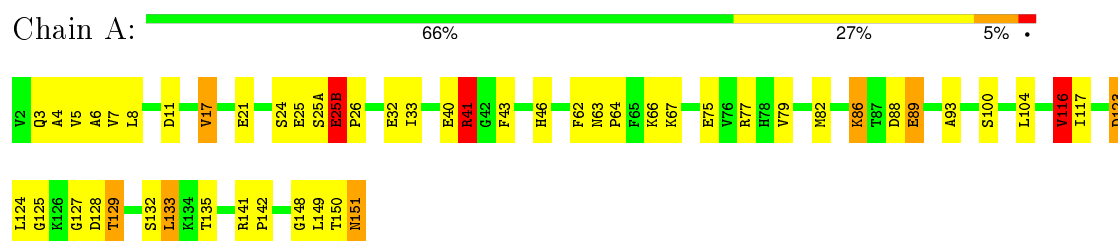
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	147	Total 147	O 147	0	0
4	B	110	Total 110	O 110	0	0
4	C	112	Total 112	O 112	0	0
4	D	144	Total 144	O 144	0	0

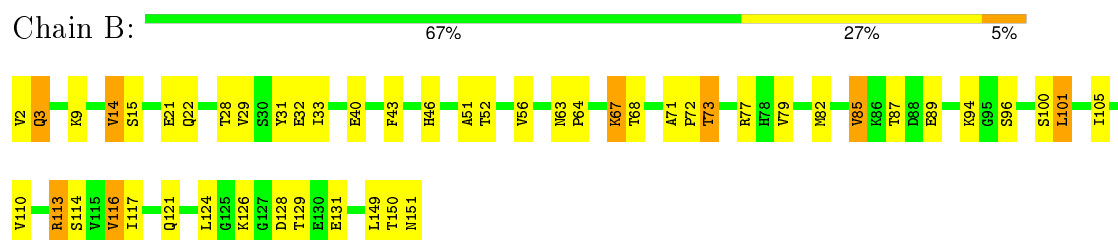
3 Residue-property plots

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 ($RSRZ > 2$). 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.

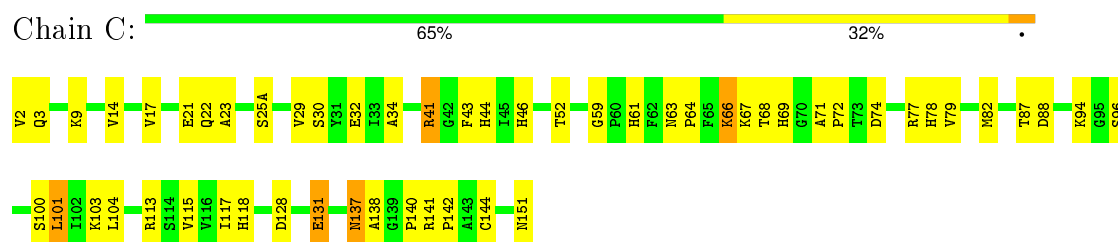
• Molecule 1: COPPER,ZINC SUPEROXIDE DISMUTASE



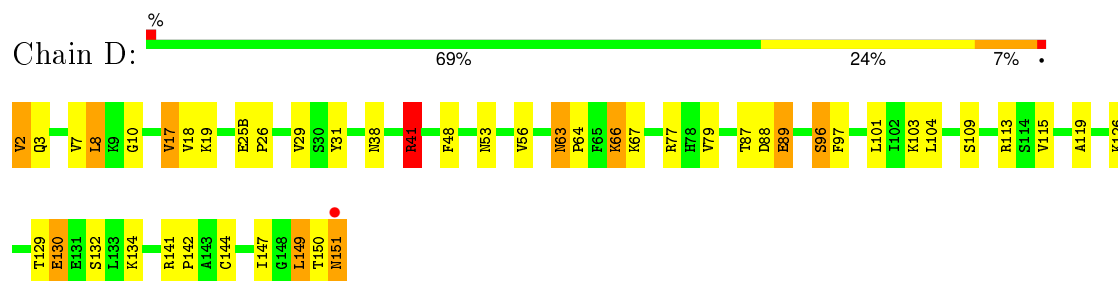
• Molecule 1: COPPER,ZINC SUPEROXIDE DISMUTASE



• Molecule 1: COPPER,ZINC SUPEROXIDE DISMUTASE



• Molecule 1: COPPER,ZINC SUPEROXIDE DISMUTASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	105.30 Å 143.00 Å 62.10 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.50 19.88 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.50) 98.1 (19.88-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.54 (at 2.50 Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.158 , (Not available) 0.165 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.156	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 93.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 32504 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4945	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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: ZN, CU

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	0.93	0/1126	1.62	17/1523 (1.1%)
1	B	0.92	0/1126	1.55	9/1523 (0.6%)
1	C	0.92	1/1126 (0.1%)	1.76	15/1523 (1.0%)
1	D	0.95	1/1126 (0.1%)	1.65	18/1523 (1.2%)
All	All	0.93	2/4504 (0.0%)	1.65	59/6092 (1.0%)

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	3	0
1	B	1	0
1	D	0	2
All	All	4	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	2	VAL	C-N	6.63	1.49	1.34
1	C	41	ARG	CZ-NH1	5.24	1.39	1.33

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	41	ARG	NE-CZ-NH2	-25.43	107.59	120.30
1	C	41	ARG	NE-CZ-NH1	21.95	131.28	120.30
1	B	113	ARG	NE-CZ-NH1	14.66	127.63	120.30
1	A	41	ARG	NE-CZ-NH1	13.99	127.30	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	41	ARG	NE-CZ-NH2	-12.92	113.84	120.30
1	D	41	ARG	NE-CZ-NH1	12.72	126.66	120.30
1	C	141	ARG	NE-CZ-NH1	12.25	126.42	120.30
1	A	141	ARG	NE-CZ-NH1	12.09	126.34	120.30
1	D	2	VAL	O-C-N	-12.05	103.42	122.70
1	A	41	ARG	NE-CZ-NH2	-11.89	114.35	120.30
1	A	141	ARG	NE-CZ-NH2	-11.22	114.69	120.30
1	B	113	ARG	NE-CZ-NH2	-10.60	115.00	120.30
1	D	141	ARG	NE-CZ-NH1	10.10	125.35	120.30
1	C	77	ARG	NE-CZ-NH2	-9.65	115.47	120.30
1	C	17	VAL	CG1-CB-CG2	-9.04	96.44	110.90
1	D	77	ARG	NE-CZ-NH2	-8.89	115.85	120.30
1	C	41	ARG	CD-NE-CZ	8.50	135.50	123.60
1	D	141	ARG	NE-CZ-NH2	-8.37	116.11	120.30
1	D	77	ARG	NE-CZ-NH1	7.75	124.17	120.30
1	D	17	VAL	CA-CB-CG1	7.34	121.91	110.90
1	C	141	ARG	NE-CZ-NH2	-7.06	116.77	120.30
1	A	77	ARG	NE-CZ-NH2	7.00	123.80	120.30
1	A	151	ASN	N-CA-CB	6.90	123.03	110.60
1	D	113	ARG	NE-CZ-NH1	-6.87	116.86	120.30
1	D	88	ASP	CB-CG-OD1	6.54	124.19	118.30
1	C	87	THR	CA-CB-CG2	-6.40	103.44	112.40
1	C	101	LEU	CB-CG-CD2	-6.38	100.15	111.00
1	B	114	SER	CB-CA-C	-6.29	98.14	110.10
1	B	28	THR	CA-CB-CG2	-6.07	103.90	112.40
1	B	113	ARG	CD-NE-CZ	6.03	132.04	123.60
1	A	25(B)	GLU	CB-CA-C	5.97	122.35	110.40
1	D	113	ARG	NE-CZ-NH2	5.87	123.24	120.30
1	C	77	ARG	NE-CZ-NH1	5.83	123.22	120.30
1	D	31	TYR	CB-CG-CD2	-5.76	117.55	121.00
1	B	73	THR	CA-CB-CG2	-5.74	104.36	112.40
1	B	87	THR	CA-CB-CG2	-5.73	104.38	112.40
1	A	123	ASP	CB-CG-OD2	-5.73	113.14	118.30
1	D	7	VAL	CG1-CB-CG2	-5.72	101.75	110.90
1	A	11	ASP	CB-CA-C	-5.69	99.02	110.40
1	C	52	THR	CA-CB-CG2	-5.64	104.50	112.40
1	B	14	VAL	CB-CA-C	5.63	122.10	111.40
1	A	7	VAL	CA-CB-CG1	-5.61	102.48	110.90
1	D	113	ARG	CD-NE-CZ	-5.60	115.76	123.60
1	D	88	ASP	CB-CG-OD2	-5.54	113.31	118.30
1	D	2	VAL	CA-C-N	5.53	129.38	117.20
1	B	77	ARG	NE-CZ-NH1	5.48	123.04	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	116	VAL	CB-CA-C	5.44	121.73	111.40
1	D	119	ALA	N-CA-CB	5.44	117.71	110.10
1	C	101	LEU	CA-CB-CG	-5.38	102.92	115.30
1	C	66	LYS	N-CA-CB	5.34	120.21	110.60
1	C	23	ALA	CB-CA-C	5.30	118.05	110.10
1	A	17	VAL	CB-CA-C	5.30	121.46	111.40
1	A	88	ASP	CB-CG-OD2	-5.29	113.54	118.30
1	A	129	THR	CA-C-N	-5.12	105.93	117.20
1	A	8	LEU	N-CA-CB	-5.12	100.17	110.40
1	D	89	GLU	CB-CA-C	-5.11	100.17	110.40
1	A	135	THR	CA-CB-CG2	-5.05	105.33	112.40
1	C	113	ARG	NE-CZ-NH2	-5.05	117.77	120.30
1	A	127	GLY	CA-C-N	-5.01	106.19	117.20

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	17	VAL	CA
1	A	116	VAL	CA
1	A	151	ASN	CA
1	B	14	VAL	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	2	VAL	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	1106	0	1068	26	0
1	B	1106	0	1067	26	0
1	C	1106	0	1067	29	0
1	D	1106	0	1067	31	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	147	0	0	2	0
4	B	110	0	0	4	0
4	C	112	0	0	4	0
4	D	144	0	0	3	0
All	All	4945	0	4269	109	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (109) 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:D:150:THR:HG23	1:D:151:ASN:N	1.81	0.95
1:B:2:VAL:HG13	1:B:22:GLN:O	1.70	0.91
1:A:46:HIS:CD2	1:A:116:VAL:HG13	2.05	0.91
1:B:63:ASN:HD21	1:B:67:LYS:H	1.22	0.86
1:B:100:SER:HB2	4:B:221:HOH:O	1.76	0.83
1:C:71:ALA:HB1	1:C:72:PRO:HD2	1.61	0.81
1:C:64:PRO:HD2	1:C:79:VAL:HG13	1.64	0.80
1:A:41:ARG:HD2	4:A:232:HOH:O	1.82	0.78
1:D:150:THR:O	1:D:151:ASN:HB2	1.82	0.77
1:B:32:GLU:HG3	1:B:94:LYS:HG3	1.69	0.75
1:C:117:ILE:O	1:C:142:PRO:HD2	1.87	0.75
1:A:64:PRO:HD2	1:A:79:VAL:HG13	1.69	0.74
1:D:63:ASN:HD21	1:D:67:LYS:H	1.33	0.74
1:D:64:PRO:HD2	1:D:79:VAL:HG13	1.68	0.74
1:A:63:ASN:HD21	1:A:67:LYS:H	1.35	0.71
1:C:2:VAL:HG13	1:C:22:GLN:O	1.90	0.71
1:B:126:LYS:HD3	4:B:246:HOH:O	1.90	0.70
1:C:63:ASN:HD21	1:C:67:LYS:H	1.41	0.68
1:A:117:ILE:O	1:A:142:PRO:HD2	1.93	0.67
1:B:2:VAL:HG11	4:D:166:HOH:O	1.97	0.65
1:C:69:HIS:HB2	1:C:78:HIS:CE1	2.32	0.65
1:D:150:THR:CG2	1:D:151:ASN:N	2.46	0.64
1:C:74:ASP:HB3	4:C:245:HOH:O	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ASN:ND2	1:A:67:LYS:H	1.97	0.62
1:D:130:GLU:HB3	4:D:244:HOH:O	2.01	0.60
1:B:43:PHE:O	1:B:82:MET:HB2	2.02	0.60
1:C:151:ASN:HB2	1:D:48:PHE:CE1	2.37	0.60
1:D:63:ASN:ND2	1:D:67:LYS:H	1.99	0.59
1:A:43:PHE:O	1:A:82:MET:HB2	2.02	0.59
1:B:63:ASN:ND2	1:B:67:LYS:H	1.96	0.59
1:D:29:VAL:O	1:D:96:SER:HA	2.04	0.58
1:B:3:GLN:HG3	1:B:21:GLU:HG3	1.86	0.58
1:C:71:ALA:HB1	1:C:72:PRO:CD	2.31	0.58
1:D:129:THR:HG23	1:D:132:SER:H	1.69	0.57
1:D:53:ASN:ND2	1:D:56:VAL:HG22	2.21	0.56
1:D:25(B):GLU:HB3	1:D:26:PRO:HD2	1.85	0.56
1:A:25:GLU:OE2	1:D:134:LYS:NZ	2.38	0.56
1:B:9:LYS:HE3	4:B:231:HOH:O	2.05	0.56
1:D:41:ARG:HD2	4:D:177:HOH:O	2.07	0.55
1:D:38:ASN:HA	1:D:87:THR:O	2.06	0.55
1:C:140:PRO:HG3	4:C:234:HOH:O	2.05	0.55
1:C:79:VAL:HB	1:C:101:LEU:HB3	1.88	0.54
1:D:64:PRO:CD	1:D:79:VAL:HG13	2.37	0.54
1:C:151:ASN:HB2	1:D:48:PHE:CZ	2.43	0.54
1:C:43:PHE:O	1:C:82:MET:HB2	2.08	0.54
1:C:29:VAL:O	1:C:96:SER:HA	2.08	0.53
1:B:40:GLU:O	1:B:121:GLN:HB2	2.07	0.53
1:B:150:THR:OG1	1:B:151:ASN:N	2.42	0.53
1:D:115:VAL:HG23	1:D:147:ILE:HD11	1.91	0.53
1:A:125:GLY:HA2	1:A:132:SER:OG	2.09	0.53
1:B:71:ALA:HB2	1:B:124:LEU:HB3	1.91	0.52
1:B:110:VAL:O	1:B:113:ARG:HG3	2.08	0.52
1:B:64:PRO:HD2	1:B:79:VAL:HG23	1.91	0.52
1:C:44:HIS:ND1	1:C:61:HIS:CE1	2.79	0.51
1:B:71:ALA:HB2	1:B:124:LEU:HD13	1.93	0.50
1:B:79:VAL:HG12	1:B:79:VAL:O	2.10	0.50
1:B:2:VAL:HG22	1:B:105:ILE:HD11	1.93	0.50
1:D:8:LEU:HD12	1:D:144:CYS:C	2.32	0.50
1:A:150:THR:OG1	1:A:151:ASN:N	2.44	0.50
1:A:123:ASP:C	1:A:123:ASP:OD1	2.51	0.49
1:A:3:GLN:HG2	1:A:151:ASN:ND2	2.27	0.49
1:D:3:GLN:HG3	1:D:151:ASN:ND2	2.27	0.49
1:C:25(A):SER:N	4:C:230:HOH:O	2.45	0.49
1:D:129:THR:OG1	1:D:130:GLU:N	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:LEU:HD12	1:A:133:LEU:HA	1.54	0.48
1:A:25(B):GLU:HB2	1:A:26:PRO:HD2	1.96	0.48
1:B:46:HIS:CD2	1:B:116:VAL:HG13	2.48	0.48
1:A:25(B):GLU:HB2	1:A:26:PRO:CD	2.45	0.47
1:D:103:LYS:O	1:D:109:SER:HA	2.15	0.46
1:D:115:VAL:O	1:D:144:CYS:HA	2.15	0.46
1:C:14:VAL:HA	1:C:34:ALA:O	2.16	0.46
1:D:53:ASN:ND2	1:D:56:VAL:CG2	2.77	0.46
1:D:66:LYS:HA	1:D:66:LYS:HD2	1.58	0.46
1:D:101:LEU:HD23	1:D:101:LEU:HA	1.44	0.45
1:A:32:GLU:C	1:A:33:ILE:HG13	2.34	0.45
1:C:118:HIS:HB3	1:C:138:ALA:O	2.16	0.45
1:A:93:ALA:HA	4:A:211:HOH:O	2.17	0.45
1:C:59:GLY:HA2	4:C:188:HOH:O	2.18	0.44
1:C:69:HIS:HB2	1:C:78:HIS:HE1	1.76	0.44
1:B:29:VAL:O	1:B:96:SER:HA	2.18	0.44
1:B:71:ALA:HB1	1:B:72:PRO:HD2	2.00	0.43
1:C:44:HIS:CE1	1:C:61:HIS:CE1	3.06	0.43
1:B:113:ARG:NH1	4:B:241:HOH:O	2.50	0.43
1:A:89:GLU:H	1:A:89:GLU:HG3	1.31	0.43
1:C:118:HIS:HA	1:C:142:PRO:HD2	2.00	0.43
1:D:97:PHE:CD1	1:D:97:PHE:C	2.92	0.42
1:C:131:GLU:HB3	1:C:137:ASN:ND2	2.34	0.42
1:C:101:LEU:HA	1:C:101:LEU:HD23	1.16	0.42
1:C:32:GLU:HG3	1:C:94:LYS:HG3	2.01	0.42
1:A:4:ALA:HA	1:A:148:GLY:O	2.19	0.42
1:D:10:GLY:HA3	1:D:142:PRO:O	2.20	0.42
1:A:43:PHE:CD2	1:A:82:MET:HE3	2.53	0.42
1:A:40:GLU:C	1:A:41:ARG:HG2	2.39	0.42
1:C:22:GLN:HE22	1:C:103:LYS:HD3	1.85	0.42
1:C:115:VAL:O	1:C:144:CYS:HA	2.20	0.42
1:B:101:LEU:HD12	1:B:101:LEU:HA	1.59	0.41
1:B:33:ILE:HD13	1:B:117:ILE:HD13	2.02	0.41
1:A:124:LEU:HD23	1:A:124:LEU:HA	1.86	0.41
1:B:31:TYR:CE2	1:B:85:VAL:HG22	2.54	0.41
1:C:88:ASP:OD1	1:C:88:ASP:C	2.57	0.41
1:D:25(B):GLU:HA	1:D:26:PRO:HD3	1.88	0.41
1:C:46:HIS:ND1	1:C:59:GLY:O	2.36	0.41
1:D:8:LEU:HD12	1:D:8:LEU:HA	1.85	0.41
1:D:149:LEU:HD12	1:D:149:LEU:HA	1.93	0.41
1:A:40:GLU:HG2	1:A:86:LYS:HG3	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:PHE:CD2	1:A:79:VAL:HG22	2.56	0.40
1:A:5:VAL:CG2	1:A:6:ALA:N	2.84	0.40
1:B:51:ALA:O	1:B:52:THR:C	2.58	0.40
1:A:5:VAL:HG22	1:A:6:ALA:N	2.36	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	151/153 (99%)	147 (97%)	4 (3%)	0	100	100
1	B	151/153 (99%)	147 (97%)	4 (3%)	0	100	100
1	C	151/153 (99%)	144 (95%)	7 (5%)	0	100	100
1	D	151/153 (99%)	147 (97%)	4 (3%)	0	100	100
All	All	604/612 (99%)	585 (97%)	19 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	118/118 (100%)	101 (86%)	17 (14%)	4	7
1	B	118/118 (100%)	103 (87%)	15 (13%)	5	10

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	118/118 (100%)	106 (90%)	12 (10%)	9	17
1	D	118/118 (100%)	104 (88%)	14 (12%)	6	12
All	All	472/472 (100%)	414 (88%)	58 (12%)	6	11

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	VAL
1	A	21	GLU
1	A	24	SER
1	A	25(A)	SER
1	A	25(B)	GLU
1	A	41	ARG
1	A	66	LYS
1	A	75	GLU
1	A	86	LYS
1	A	89	GLU
1	A	100	SER
1	A	104	LEU
1	A	116	VAL
1	A	128	ASP
1	A	129	THR
1	A	133	LEU
1	A	149	LEU
1	B	3	GLN
1	B	14	VAL
1	B	15	SER
1	B	56	VAL
1	B	67	LYS
1	B	68	THR
1	B	73	THR
1	B	85	VAL
1	B	89	GLU
1	B	101	LEU
1	B	116	VAL
1	B	128	ASP
1	B	129	THR
1	B	131	GLU
1	B	149	LEU
1	C	3	GLN
1	C	9	LYS

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Mol	Chain	Res	Type
1	C	21	GLU
1	C	30	SER
1	C	41	ARG
1	C	66	LYS
1	C	68	THR
1	C	100	SER
1	C	104	LEU
1	C	128	ASP
1	C	131	GLU
1	C	137	ASN
1	D	8	LEU
1	D	17	VAL
1	D	18	VAL
1	D	19	LYS
1	D	41	ARG
1	D	63	ASN
1	D	66	LYS
1	D	89	GLU
1	D	96	SER
1	D	104	LEU
1	D	126	LYS
1	D	130	GLU
1	D	149	LEU
1	D	151	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	B	63	ASN
1	C	63	ASN
1	C	137	ASN
1	C	151	ASN
1	D	3	GLN
1	D	63	ASN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 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	153/153 (100%)	-0.80	0 100 100	8, 21, 52, 79	8 (5%)
1	B	153/153 (100%)	-0.75	0 100 100	14, 26, 51, 65	8 (5%)
1	C	153/153 (100%)	-0.63	0 100 100	13, 23, 48, 69	8 (5%)
1	D	153/153 (100%)	-0.76	1 (0%) 89 90	12, 22, 45, 72	6 (3%)
All	All	612/612 (100%)	-0.73	1 (0%) 95 96	8, 23, 50, 79	30 (4%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	151	ASN	2.9

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(\AA^2)	Q<0.9
3	ZN	A	153	1/1	1.00	0.04	-3.31	23,23,23,23	0
3	ZN	D	153	1/1	0.99	0.05	-4.33	20,20,20,20	0
3	ZN	B	153	1/1	0.99	0.04	-6.42	23,23,23,23	0
3	ZN	C	153	1/1	0.99	0.04	-6.93	21,21,21,21	0
2	CU	A	152	1/1	1.00	0.02	-	25,25,25,25	0
2	CU	C	152	1/1	0.99	0.05	-	22,22,22,22	0
2	CU	D	152	1/1	0.99	0.05	-	25,25,25,25	0
2	CU	B	152	1/1	0.99	0.04	-	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.