



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:54 PM UTC

PDB ID : 1RHS / pdb_00001rhs
Title : SULFUR-SUBSTITUTED RHODANESE
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Deposited on : 1997-07-16
Resolution : 1.36 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

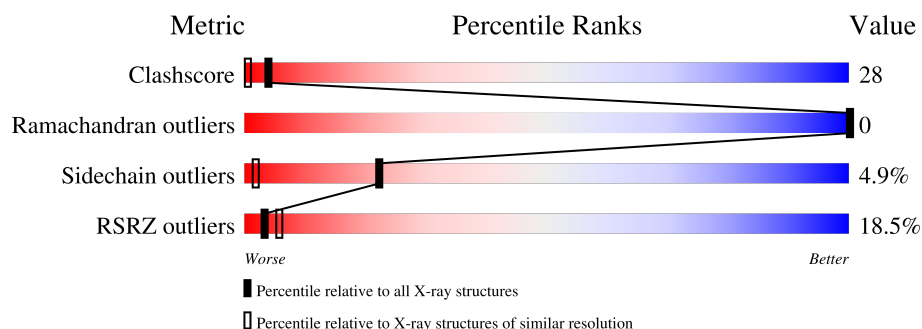
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.36 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	190562	1232 (1.36-1.36)
Ramachandran outliers	187476	1220 (1.36-1.36)
Sidechain outliers	187428	1220 (1.36-1.36)
RSRZ outliers	180081	1214 (1.36-1.36)

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	296	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2733 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 SULFUR-SUBSTITUTED RHODANESE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	293	2326	1486	405	425	10	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	247	CSS	CYS	conflict	UNP P00586

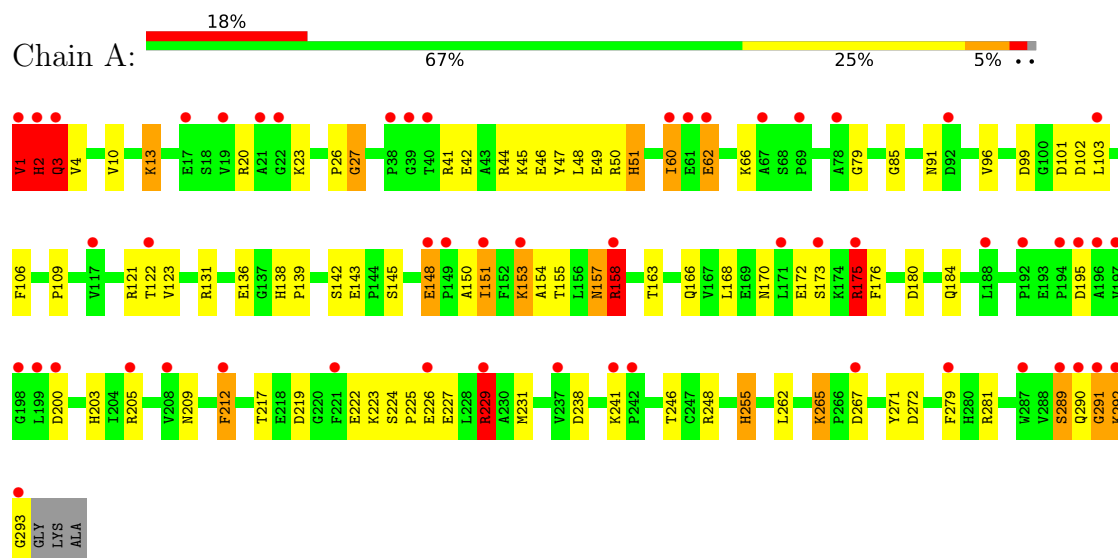
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	407	Total	O	0	0
			407	407		

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: SULFUR-SUBSTITUTED RHODANESE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	154.36Å 49.31Å 41.68Å 90.00° 99.81° 90.00°	Depositor
Resolution (Å)	25.00 – 1.36 25.00 – 1.36	Depositor EDS
% Data completeness (in resolution range)	74.0 (25.00-1.36) 70.6 (25.00-1.36)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.63 (at 1.30Å)	Xtriage
Refinement program	SHELXL-93	Depositor
R, R_{free}	0.169 , 0.229 0.200 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	5.8	Xtriage
Anisotropy	0.955	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.52 , 57.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2733	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.55% 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: CSS

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	1.65	21/2385 (0.9%)	1.79	51/3235 (1.6%)

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	1	6

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	109	PRO	CA-CB	-7.76	1.43	1.53
1	A	3	GLN	CA-CB	-7.64	1.41	1.53
1	A	122	THR	CB-OG1	-7.25	1.32	1.43
1	A	151	ILE	CA-CB	6.75	1.62	1.54
1	A	27	GLY	C-O	6.74	1.31	1.24
1	A	51	HIS	CD2-NE2	-6.58	1.30	1.37
1	A	203	HIS	CG-ND1	6.21	1.45	1.38
1	A	10	VAL	CA-CB	-6.00	1.46	1.54
1	A	4	VAL	N-CA	5.93	1.53	1.46
1	A	123	VAL	C-O	5.91	1.29	1.24
1	A	3	GLN	CA-C	5.90	1.58	1.53
1	A	175	ARG	CZ-NH1	5.74	1.40	1.32
1	A	79	GLY	C-N	5.68	1.41	1.33
1	A	145	SER	C-N	-5.62	1.25	1.33
1	A	255	HIS	CE1-NE2	5.48	1.38	1.32
1	A	195	ASP	CG-OD2	5.46	1.35	1.25
1	A	289	SER	C-O	5.39	1.30	1.23
1	A	271	TYR	C-O	5.30	1.30	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	219	ASP	CG-OD2	5.23	1.35	1.25
1	A	262	LEU	N-CA	5.18	1.53	1.46
1	A	145	SER	C-O	5.08	1.29	1.23

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	148	GLU	CB-CG-CD	24.18	153.71	112.60
1	A	205	ARG	CD-NE-CZ	10.43	139.00	124.40
1	A	175	ARG	NE-CZ-NH2	9.76	127.98	119.20
1	A	281	ARG	NE-CZ-NH1	-9.04	112.46	121.50
1	A	158	ARG	NE-CZ-NH2	8.77	127.09	119.20
1	A	1	VAL	CB-CA-C	8.54	127.63	111.40
1	A	281	ARG	NE-CZ-NH2	8.03	126.42	119.20
1	A	3	GLN	O-C-N	7.90	131.51	122.19
1	A	122	THR	CA-CB-OG1	7.50	120.84	109.60
1	A	195	ASP	CA-CB-CG	7.43	120.03	112.60
1	A	139	PRO	CB-CA-C	-7.15	102.24	111.46
1	A	289	SER	CA-C-O	-7.14	111.76	120.54
1	A	175	ARG	NE-CZ-NH1	-7.08	114.42	121.50
1	A	205	ARG	NE-CZ-NH1	7.01	128.51	121.50
1	A	158	ARG	NE-CZ-NH1	-7.01	114.49	121.50
1	A	2	HIS	CB-CA-C	-6.91	96.67	110.42
1	A	3	GLN	N-CA-CB	6.85	122.53	110.27
1	A	158	ARG	CD-NE-CZ	6.60	133.64	124.40
1	A	96	VAL	CA-C-O	-6.25	113.89	120.39
1	A	279	PHE	CA-C-O	6.25	127.17	120.55
1	A	3	GLN	N-CA-C	-6.03	103.46	111.96
1	A	272	ASP	CA-CB-CG	6.01	118.61	112.60
1	A	44	ARG	CD-NE-CZ	5.82	132.54	124.40
1	A	23	LYS	N-CA-C	5.78	119.59	112.54
1	A	20	ARG	O-C-N	5.64	128.10	122.12
1	A	168	LEU	O-C-N	-5.62	116.29	122.07
1	A	150	ALA	CA-C-N	-5.61	116.23	123.19
1	A	150	ALA	C-N-CA	-5.61	116.23	123.19
1	A	212	PHE	N-CA-C	5.60	118.15	111.71
1	A	102	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	3	GLN	CA-C-N	-5.50	114.08	121.66
1	A	3	GLN	C-N-CA	-5.50	114.08	121.66
1	A	290	GLN	CA-C-N	-5.41	110.80	121.41
1	A	290	GLN	C-N-CA	-5.41	110.80	121.41
1	A	265	LYS	CG-CD-CE	5.41	123.75	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	200	ASP	CA-CB-CG	5.38	117.98	112.60
1	A	62	GLU	CB-CG-CD	-5.31	103.58	112.60
1	A	23	LYS	CA-C-O	5.29	125.89	119.38
1	A	291	GLY	CA-C-N	5.29	131.22	121.70
1	A	291	GLY	C-N-CA	5.29	131.22	121.70
1	A	60	ILE	CB-CG1-CD1	5.28	124.88	113.80
1	A	99	ASP	CA-C-O	-5.25	115.83	121.45
1	A	13	LYS	O-C-N	5.19	127.41	122.07
1	A	101	ASP	CA-C-N	5.15	127.70	120.28
1	A	101	ASP	C-N-CA	5.15	127.70	120.28
1	A	1	VAL	N-CA-CB	5.13	120.22	111.50
1	A	51	HIS	ND1-CE1-NE2	-5.08	103.32	108.40
1	A	101	ASP	O-C-N	-5.08	116.78	122.97
1	A	279	PHE	O-C-N	-5.08	116.74	122.12
1	A	85	GLY	CA-C-O	5.06	125.96	120.75
1	A	20	ARG	NE-CZ-NH1	5.02	126.52	121.50

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	1	VAL	CA

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	VAL	Peptide
1	A	131	ARG	Sidechain
1	A	158	ARG	Sidechain
1	A	175	ARG	Sidechain
1	A	229	ARG	Sidechain
1	A	248	ARG	Sidechain

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	A	2326	0	2262	129	9

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	407	0	0	57	8
All	All	2733	0	2262	129	10

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 28.

All (129) 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:2:HIS:NE2	1:A:121:ARG:CD	1.81	1.41
1:A:2:HIS:NE2	1:A:121:ARG:HD2	1.11	1.38
1:A:229:ARG:HH11	1:A:229:ARG:CB	1.49	1.24
1:A:223:LYS:HB3	2:A:346:HOH:O	1.34	1.24
1:A:229:ARG:HB3	1:A:229:ARG:NH1	1.52	1.24
1:A:227:GLU:HB2	2:A:346:HOH:O	1.38	1.22
1:A:13:LYS:HA	2:A:671:HOH:O	1.40	1.21
1:A:2:HIS:CD2	2:A:461:HOH:O	1.88	1.20
1:A:153:LYS:HD2	1:A:153:LYS:C	1.49	1.19
1:A:292:LYS:H	1:A:293:GLY:HA3	1.00	1.15
1:A:2:HIS:CE1	1:A:121:ARG:HD2	1.82	1.12
1:A:47:TYR:HE2	2:A:446:HOH:O	1.31	1.12
1:A:50:ARG:N	2:A:462:HOH:O	1.83	1.09
1:A:46:GLU:HA	2:A:391:HOH:O	1.48	1.08
1:A:50:ARG:HG3	2:A:462:HOH:O	1.53	1.08
1:A:13:LYS:CA	2:A:671:HOH:O	1.98	1.06
1:A:226:GLU:HA	1:A:229:ARG:HH12	0.95	1.06
1:A:2:HIS:NE2	2:A:461:HOH:O	1.86	1.05
1:A:265:LYS:NZ	1:A:267:ASP:OD1	1.88	1.04
1:A:3:GLN:HA	1:A:3:GLN:OE1	1.43	1.03
1:A:46:GLU:CA	2:A:391:HOH:O	2.04	1.02
1:A:13:LYS:CB	2:A:671:HOH:O	2.06	1.02
1:A:170:ASN:HD22	1:A:175:ARG:HH11	1.09	0.99
1:A:2:HIS:NE2	1:A:121:ARG:NE	2.10	0.99
1:A:170:ASN:HB2	1:A:175:ARG:NH1	1.77	0.99
1:A:292:LYS:N	1:A:293:GLY:HA3	1.75	0.99
1:A:224:SER:N	2:A:346:HOH:O	1.96	0.96
1:A:153:LYS:C	1:A:153:LYS:CD	2.40	0.95
1:A:2:HIS:CE1	1:A:121:ARG:CD	2.45	0.95
1:A:226:GLU:HA	1:A:229:ARG:NH1	1.80	0.95
1:A:103:LEU:CD1	2:A:658:HOH:O	2.13	0.94
1:A:292:LYS:H	1:A:293:GLY:CA	1.79	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:GLU:CA	1:A:229:ARG:HH12	1.80	0.93
1:A:217:THR:HG23	2:A:348:HOH:O	0.74	0.92
1:A:238:ASP:OD2	1:A:241:LYS:NZ	2.03	0.91
1:A:49:GLU:C	2:A:462:HOH:O	2.09	0.90
1:A:173:SER:OG	1:A:175:ARG:HG2	1.72	0.90
1:A:170:ASN:HD21	1:A:176:PHE:H	1.11	0.89
1:A:49:GLU:HA	2:A:467:HOH:O	1.73	0.89
1:A:229:ARG:HH11	1:A:229:ARG:HB3	0.72	0.88
1:A:170:ASN:HB2	1:A:175:ARG:HH12	1.36	0.86
1:A:26:PRO:C	2:A:411:HOH:O	2.19	0.83
1:A:170:ASN:ND2	1:A:175:ARG:HH11	1.75	0.83
1:A:229:ARG:CB	1:A:229:ARG:NH1	2.24	0.82
1:A:49:GLU:CB	2:A:462:HOH:O	2.28	0.81
1:A:153:LYS:HD2	1:A:153:LYS:O	1.79	0.81
1:A:222:GLU:HG2	2:A:347:HOH:O	1.80	0.80
1:A:170:ASN:CB	1:A:175:ARG:NH1	2.44	0.80
1:A:175:ARG:HH11	1:A:175:ARG:HG3	1.46	0.79
1:A:91:ASN:HD21	1:A:153:LYS:H	1.27	0.79
1:A:142:SER:HB2	2:A:467:HOH:O	1.82	0.79
1:A:106:PHE:CE2	2:A:443:HOH:O	2.36	0.78
1:A:41:ARG:CB	2:A:380:HOH:O	2.31	0.77
1:A:163:THR:H	1:A:166:GLN:HE21	1.33	0.77
1:A:175:ARG:HH11	1:A:175:ARG:CG	1.98	0.76
1:A:103:LEU:HD13	2:A:658:HOH:O	1.80	0.75
1:A:2:HIS:CE1	1:A:121:ARG:NE	2.54	0.75
1:A:48:LEU:O	2:A:467:HOH:O	2.04	0.74
1:A:175:ARG:HG3	1:A:176:PHE:HD1	1.54	0.73
1:A:49:GLU:HB2	2:A:462:HOH:O	1.88	0.71
1:A:153:LYS:HD2	1:A:154:ALA:N	2.03	0.71
1:A:13:LYS:HB2	2:A:671:HOH:O	1.81	0.71
1:A:170:ASN:ND2	1:A:175:ARG:NH1	2.40	0.70
1:A:27:GLY:N	2:A:411:HOH:O	2.26	0.69
1:A:175:ARG:HG3	1:A:176:PHE:CD1	2.27	0.68
1:A:46:GLU:CB	2:A:391:HOH:O	2.38	0.68
1:A:170:ASN:HD22	1:A:175:ARG:NH1	1.87	0.67
1:A:41:ARG:HB3	2:A:380:HOH:O	1.94	0.67
1:A:175:ARG:NH1	1:A:175:ARG:HG3	2.09	0.67
1:A:170:ASN:ND2	1:A:176:PHE:H	1.90	0.66
1:A:103:LEU:HD11	2:A:658:HOH:O	1.84	0.66
1:A:27:GLY:N	2:A:636:HOH:O	2.24	0.65
1:A:50:ARG:CB	2:A:462:HOH:O	2.45	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:PHE:CD2	2:A:443:HOH:O	2.46	0.65
1:A:47:TYR:O	1:A:51:HIS:HD2	1.80	0.65
1:A:227:GLU:OE1	2:A:346:HOH:O	2.15	0.65
1:A:291:GLY:N	1:A:292:LYS:HA	2.11	0.64
1:A:170:ASN:HD22	1:A:175:ARG:CG	2.11	0.64
1:A:106:PHE:O	1:A:255:HIS:HE1	1.82	0.63
1:A:153:LYS:HE2	1:A:155:THR:OG1	1.99	0.63
1:A:157:ASN:C	1:A:157:ASN:HD22	2.05	0.63
1:A:289:SER:O	1:A:292:LYS:HA	1.99	0.63
1:A:222:GLU:CG	2:A:347:HOH:O	2.43	0.62
1:A:138:HIS:HD2	2:A:448:HOH:O	1.83	0.62
1:A:238:ASP:CG	1:A:241:LYS:HZ1	2.07	0.61
1:A:142:SER:CB	2:A:467:HOH:O	2.45	0.59
1:A:46:GLU:C	2:A:391:HOH:O	2.40	0.58
1:A:223:LYS:C	2:A:346:HOH:O	2.37	0.56
1:A:2:HIS:CE1	2:A:461:HOH:O	2.44	0.55
1:A:222:GLU:HB3	2:A:347:HOH:O	2.07	0.55
1:A:229:ARG:NH1	1:A:229:ARG:HB2	2.19	0.54
1:A:170:ASN:HD22	1:A:175:ARG:HG2	1.71	0.54
1:A:173:SER:OG	1:A:175:ARG:CG	2.52	0.54
1:A:175:ARG:NH1	1:A:176:PHE:HD1	2.06	0.54
1:A:212:PHE:HZ	2:A:443:HOH:O	1.91	0.53
1:A:136:GLU:OE1	1:A:138:HIS:HE1	1.91	0.52
1:A:225:PRO:O	1:A:229:ARG:NH1	2.42	0.52
1:A:170:ASN:HD22	1:A:175:ARG:HG3	1.74	0.52
1:A:229:ARG:HB2	1:A:229:ARG:CZ	2.41	0.51
1:A:292:LYS:N	1:A:293:GLY:CA	2.55	0.50
1:A:2:HIS:CD2	1:A:121:ARG:HE	2.29	0.50
1:A:46:GLU:HB3	2:A:391:HOH:O	2.07	0.50
1:A:153:LYS:CE	1:A:155:THR:OG1	2.60	0.50
1:A:175:ARG:NH1	1:A:176:PHE:CD1	2.80	0.49
1:A:231:MET:N	2:A:397:HOH:O	2.27	0.49
1:A:170:ASN:CG	1:A:175:ARG:NH1	2.71	0.49
1:A:46:GLU:O	2:A:391:HOH:O	2.20	0.48
1:A:222:GLU:CB	2:A:347:HOH:O	2.62	0.47
1:A:289:SER:HB3	1:A:292:LYS:O	2.13	0.47
1:A:143:GLU:HB3	2:A:452:HOH:O	2.14	0.47
1:A:106:PHE:HE2	2:A:443:HOH:O	1.87	0.47
1:A:50:ARG:CG	2:A:462:HOH:O	2.28	0.46
1:A:26:PRO:HB2	2:A:411:HOH:O	2.15	0.46
1:A:217:THR:CG2	2:A:348:HOH:O	1.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:ASN:ND2	1:A:175:ARG:HG3	2.30	0.46
1:A:291:GLY:H	1:A:292:LYS:HA	1.78	0.46
1:A:158:ARG:NH2	1:A:166:GLN:HE22	2.14	0.45
1:A:13:LYS:HG3	2:A:671:HOH:O	2.17	0.44
1:A:238:ASP:OD2	1:A:241:LYS:CE	2.65	0.44
1:A:180:ASP:HA	1:A:246:THR:O	2.18	0.43
1:A:170:ASN:ND2	1:A:175:ARG:CG	2.80	0.43
1:A:27:GLY:CA	2:A:411:HOH:O	2.65	0.43
1:A:62:GLU:HG3	2:A:694:HOH:O	2.19	0.43
1:A:2:HIS:NE2	1:A:121:ARG:CB	2.82	0.42
1:A:184:GLN:HA	1:A:209:ASN:HD21	1.86	0.41
1:A:292:LYS:O	1:A:292:LYS:HG2	2.15	0.41
1:A:223:LYS:CA	2:A:346:HOH:O	2.60	0.41
1:A:26:PRO:HB2	2:A:636:HOH:O	2.20	0.40
1:A:175:ARG:HG2	1:A:175:ARG:H	1.58	0.40

All (10) 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)
1:A:172:GLU:OE2	1:A:172:GLU:OE2[2_657]	0.93	1.27
1:A:42:GLU:CG	2:A:527:HOH:O[4_557]	1.01	1.19
1:A:45:LYS:CB	2:A:663:HOH:O[4_557]	1.07	1.13
1:A:42:GLU:CB	2:A:527:HOH:O[4_557]	1.36	0.84
1:A:45:LYS:CG	2:A:663:HOH:O[4_557]	1.36	0.84
1:A:42:GLU:CD	2:A:527:HOH:O[4_557]	1.49	0.71
1:A:45:LYS:CD	2:A:663:HOH:O[4_557]	1.67	0.53
2:A:533:HOH:O	2:A:668:HOH:O[1_554]	2.00	0.20
1:A:42:GLU:OE1	2:A:527:HOH:O[4_557]	2.14	0.06
1:A:172:GLU:CD	1:A:172:GLU:OE2[2_657]	2.18	0.02

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	290/296 (98%)	277 (96%)	13 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	244/245 (100%)	232 (95%)	12 (5%)	22	2

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	VAL
1	A	2	HIS
1	A	3	GLN
1	A	60	ILE
1	A	66	LYS
1	A	148	GLU
1	A	151	ILE
1	A	153	LYS
1	A	157	ASN
1	A	158	ARG
1	A	229	ARG
1	A	292	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	HIS
1	A	91	ASN
1	A	138	HIS
1	A	157	ASN
1	A	166	GLN

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Mol	Chain	Res	Type
1	A	170	ASN
1	A	191	GLN
1	A	209	ASN
1	A	255	HIS
1	A	290	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	CSS	A	247	1	4,6,7	0.95	0	2,6,8	1.18	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
1	CSS	A	247	1	-	0/1/5/7	-

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	292/296 (98%)	1.30	54 (18%) 3 6	11, 16, 26, 114	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	VAL	12.3
1	A	2	HIS	9.0
1	A	293	GLY	8.7
1	A	291	GLY	7.0
1	A	292	LYS	7.0
1	A	3	GLN	5.4
1	A	175	ARG	4.7
1	A	290	GLN	4.4
1	A	60	ILE	4.4
1	A	267	ASP	4.3
1	A	148	GLU	4.0
1	A	151	ILE	3.8
1	A	229	ARG	3.8
1	A	39	GLY	3.7
1	A	197	VAL	3.5
1	A	171	LEU	3.4
1	A	38	PRO	3.3
1	A	208	VAL	3.2
1	A	40	THR	3.2
1	A	67	ALA	3.1
1	A	62	GLU	3.1
1	A	287	TRP	2.9
1	A	241	LYS	2.8
1	A	149	PRO	2.8
1	A	195	ASP	2.8
1	A	226	GLU	2.8
1	A	196	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	21	ALA	2.6
1	A	242	PRO	2.6
1	A	158	ARG	2.5
1	A	200	ASP	2.5
1	A	22	GLY	2.5
1	A	153	LYS	2.4
1	A	289	SER	2.4
1	A	103	LEU	2.4
1	A	237	VAL	2.4
1	A	212	PHE	2.4
1	A	192	PRO	2.4
1	A	19	VAL	2.3
1	A	198	GLY	2.3
1	A	199	LEU	2.3
1	A	69	PRO	2.3
1	A	194	PRO	2.3
1	A	92	ASP	2.2
1	A	17	GLU	2.1
1	A	205	ARG	2.1
1	A	188	LEU	2.1
1	A	122	THR	2.1
1	A	173	SER	2.1
1	A	221	PHE	2.1
1	A	279	PHE	2.1
1	A	117	VAL	2.1
1	A	78	ALA	2.0
1	A	61	GLU	2.0

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
1	CSS	A	247	7/8	0.99	0.06	11,11,12,13	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.