



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 03:16 AM UTC

PDB ID : 3G4E / pdb_00003g4e
Title : Crystal structure of human senescence marker protein-30(SMP30)(Calcium bound)
Authors : Chakraborti, S.; Bahnson, B.J.
Deposited on : 2009-02-03
Resolution : 1.42 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
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

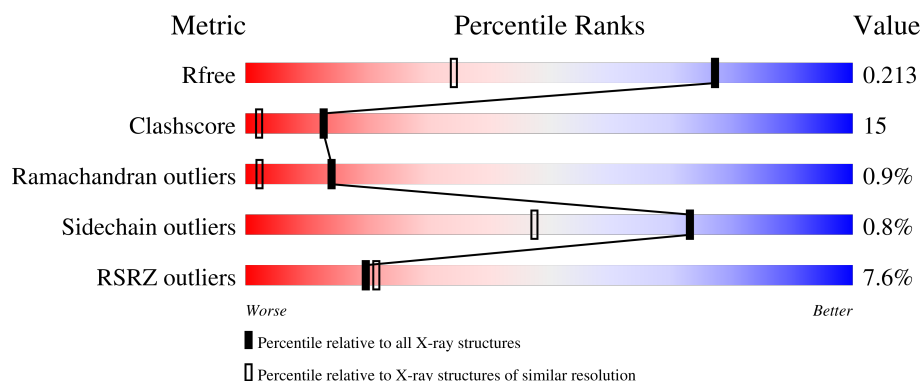
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.42 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	4041 (1.44-1.40)
Clashscore	190562	4154 (1.44-1.40)
Ramachandran outliers	187476	4083 (1.44-1.40)
Sidechain outliers	187428	4082 (1.44-1.40)
RSRZ outliers	180081	4039 (1.44-1.40)

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	297	 8% 74% 21% .
1	B	297	 7% 78% 17% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5217 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 Regucalcin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	13	0
			2365	1495	402	452	16			
1	B	297	Total	C	N	O	S	8	11	0
			2353	1488	399	449	17			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		

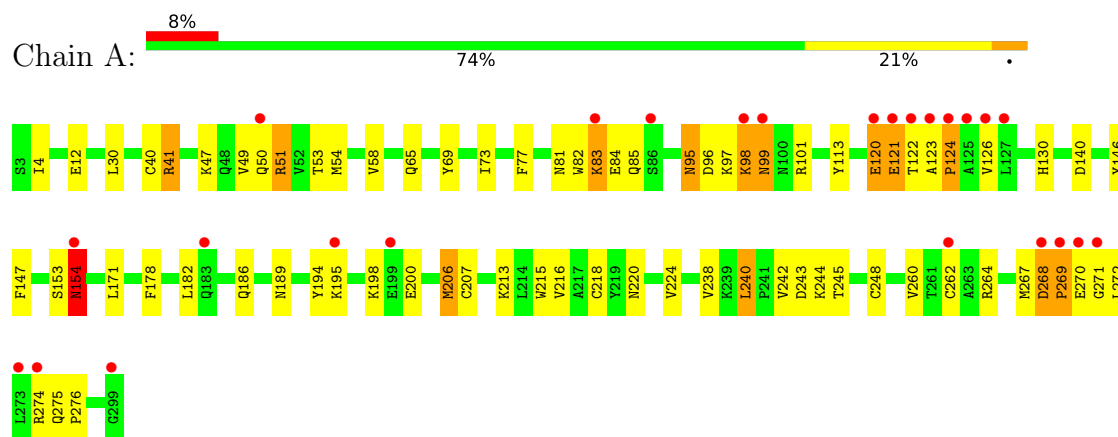
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	234	Total	O	0	0
			234	234		
3	B	263	Total	O	0	0
			263	263		

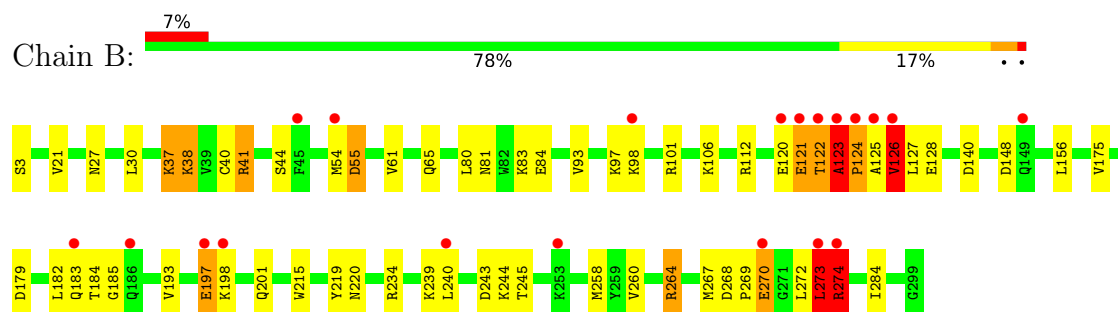
3 Residue-property plots

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: Regucalcin



• Molecule 1: Regucalcin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.47Å 51.00Å 86.12Å 90.00° 100.11° 90.00°	Depositor
Resolution (Å)	8.00 – 1.42 8.00 – 1.42	Depositor EDS
% Data completeness (in resolution range)	95.1 (8.00-1.42) 99.3 (8.00-1.42)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 1.42Å)	Xtriage
Refinement program	SHELX, SHELXL-97	Depositor
R, R_{free}	0.129 , 0.191 0.171 , 0.213	Depositor DCC
R_{free} test set	5215 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	15.6	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.59 , 84.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5217	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.91% 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

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	1.62	20/2423 (0.8%)	1.79	46/3278 (1.4%)
1	B	1.40	20/2409 (0.8%)	1.80	39/3258 (1.2%)
All	All	1.51	40/4832 (0.8%)	1.80	85/6536 (1.3%)

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	0	4
1	B	0	2
All	All	0	6

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	98	LYS	C-N	30.26	1.76	1.33
1	A	272	LEU	C-N	-28.23	0.95	1.33
1	B	122[A]	THR	C-O	27.95	1.60	1.24
1	A	269	PRO	C-N	-24.04	1.00	1.33
1	B	148	ASP	C-N	-22.67	1.02	1.33
1	B	269	PRO	C-N	-20.91	1.05	1.33
1	A	120	GLU	C-N	-19.30	1.06	1.33
1	A	270	GLU	C-N	17.03	1.58	1.33
1	B	182	LEU	C-N	-16.91	1.09	1.33
1	B	123[A]	ALA	C-O	16.31	1.45	1.24
1	A	83	LYS	C-N	15.61	1.57	1.33
1	A	124[A]	PRO	C-O	15.35	1.40	1.23
1	A	154	ASN	C-N	15.15	1.52	1.33
1	A	268	ASP	C-N	-14.71	1.11	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	185	GLY	C-N	-13.39	1.15	1.33
1	A	182	LEU	C-N	-11.25	1.17	1.33
1	B	197	GLU	C-N	-11.03	1.19	1.33
1	A	82	TRP	C-N	-10.98	1.20	1.33
1	A	41	ARG	C-N	-10.84	1.18	1.33
1	A	195	LYS	C-N	-9.97	1.20	1.33
1	A	240	LEU	C-N	-9.77	1.24	1.34
1	A	122[A]	THR	C-O	-9.42	1.12	1.24
1	B	37	LYS	C-N	-9.35	1.20	1.33
1	B	121[A]	GLU	C-O	-8.81	1.12	1.24
1	B	123[A]	ALA	CA-C	8.79	1.64	1.52
1	A	49	VAL	C-N	8.74	1.46	1.33
1	B	183	GLN	C-N	-8.02	1.22	1.33
1	A	50	GLN	C-N	-7.79	1.22	1.33
1	B	123[A]	ALA	N-CA	7.42	1.56	1.46
1	B	121[A]	GLU	C-N	7.29	1.44	1.33
1	A	124[A]	PRO	CA-C	-6.94	1.45	1.52
1	A	198	LYS	C-N	-6.92	1.23	1.33
1	B	198	LYS	C-N	-6.49	1.25	1.33
1	B	274	ARG	C-O	-5.63	1.16	1.24
1	A	96	ASP	C-N	5.49	1.41	1.33
1	B	140	ASP	C-N	-5.49	1.25	1.33
1	B	270	GLU	C-N	-5.37	1.26	1.33
1	B	38	LYS	C-N	5.36	1.41	1.33
1	B	197	GLU	C-O	5.10	1.30	1.23
1	B	272	LEU	C-O	-5.00	1.18	1.24

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	122[A]	THR	CA-C-O	25.95	150.06	119.18
1	B	123[A]	ALA	CA-C-O	-20.67	91.84	120.16
1	B	97	LYS	CA-C-N	-20.36	89.61	122.79
1	B	97	LYS	C-N-CA	-20.36	89.61	122.79
1	A	154	ASN	O-C-N	-20.16	95.78	122.59
1	B	121[A]	GLU	O-C-N	-19.98	97.25	123.23
1	A	153	SER	CA-C-N	-17.95	87.26	121.54
1	A	153	SER	C-N-CA	-17.95	87.26	121.54
1	A	124[A]	PRO	CA-C-O	-17.34	100.34	122.19
1	B	197	GLU	O-C-N	-16.75	103.75	122.85
1	A	122[A]	THR	O-C-N	13.56	136.47	122.10
1	B	122[A]	THR	O-C-N	-13.37	103.42	122.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	LYS	O-C-N	12.94	138.93	122.37
1	B	121[A]	GLU	CA-C-O	-11.97	107.78	120.70
1	A	153	SER	O-C-N	11.74	137.38	122.93
1	A	122[A]	THR	CA-C-O	-11.37	105.58	118.55
1	B	123[A]	ALA	O-C-N	-10.95	108.73	121.32
1	A	269	PRO	CA-C-N	10.71	142.00	121.54
1	A	269	PRO	C-N-CA	10.71	142.00	121.54
1	A	113	TYR	CB-CG-CD1	10.68	136.82	120.80
1	B	197	GLU	CA-C-N	10.53	134.39	120.28
1	B	197	GLU	C-N-CA	10.53	134.39	120.28
1	A	154	ASN	CA-C-N	10.47	130.29	120.34
1	A	154	ASN	C-N-CA	10.47	130.29	120.34
1	A	113	TYR	CG-CD1-CE1	10.32	136.68	121.20
1	A	98	LYS	CA-C-N	-10.13	102.19	121.54
1	A	98	LYS	C-N-CA	-10.13	102.19	121.54
1	B	184	THR	CA-C-N	9.81	135.25	122.00
1	B	184	THR	C-N-CA	9.81	135.25	122.00
1	A	270	GLU	CA-C-N	-9.61	108.12	120.14
1	A	270	GLU	C-N-CA	-9.61	108.12	120.14
1	A	124[A]	PRO	O-C-N	-9.24	110.67	123.05
1	B	123[A]	ALA	N-CA-C	9.23	130.22	109.81
1	A	267	MET	O-C-N	-9.22	112.46	123.16
1	A	121[A]	GLU	CA-C-O	-8.26	112.45	121.94
1	B	126	VAL	CA-C-O	8.23	131.07	120.78
1	A	82	TRP	CA-C-N	8.22	133.15	120.82
1	A	82	TRP	C-N-CA	8.22	133.15	120.82
1	A	272	LEU	O-C-N	8.22	133.34	122.33
1	B	97	LYS	O-C-N	8.10	132.13	123.06
1	A	194	TYR	O-C-N	-7.70	114.64	123.42
1	A	206	MET	O-C-N	-7.62	114.80	123.48
1	B	269	PRO	O-C-N	7.58	130.95	122.24
1	A	274	ARG	CA-C-N	-7.57	114.72	122.74
1	A	274	ARG	C-N-CA	-7.57	114.72	122.74
1	A	240	LEU	CA-C-O	-7.47	112.58	120.87
1	B	185	GLY	N-CA-C	-7.46	104.45	115.72
1	B	121[A]	GLU	N-CA-C	7.33	121.35	109.40
1	A	194	TYR	CA-C-N	7.32	132.51	122.77
1	A	194	TYR	C-N-CA	7.32	132.51	122.77
1	B	184	THR	O-C-N	-7.11	113.87	122.48
1	A	270	GLU	O-C-N	7.06	131.98	122.59
1	A	113	TYR	CD1-CG-CD2	-6.79	107.91	118.10
1	A	140	ASP	CA-CB-CG	-6.69	105.91	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	148	ASP	CA-C-N	6.52	132.13	122.74
1	B	148	ASP	C-N-CA	6.52	132.13	122.74
1	A	77	PHE	CA-CB-CG	6.46	120.25	113.80
1	B	264	ARG	CG-CD-NE	-6.25	98.25	112.00
1	A	268	ASP	CA-C-N	-6.15	112.31	119.32
1	A	268	ASP	C-N-CA	-6.15	112.31	119.32
1	B	55	ASP	CA-CB-CG	6.09	118.69	112.60
1	B	234	ARG	CD-NE-CZ	5.98	132.77	124.40
1	B	185	GLY	CA-C-N	5.97	129.96	121.42
1	B	185	GLY	C-N-CA	5.97	129.96	121.42
1	B	182	LEU	CA-C-O	-5.92	113.41	120.10
1	A	51	ARG	CD-NE-CZ	5.91	132.67	124.40
1	B	3	SER	CA-C-N	-5.82	114.87	122.37
1	B	3	SER	C-N-CA	-5.82	114.87	122.37
1	A	40	CYS	CA-C-N	5.78	130.92	122.77
1	A	40	CYS	C-N-CA	5.78	130.92	122.77
1	A	269	PRO	O-C-N	-5.75	115.62	122.24
1	A	198	LYS	O-C-N	-5.72	116.06	122.12
1	B	179	ASP	CA-CB-CG	-5.70	106.90	112.60
1	B	201	GLN	CA-C-N	-5.66	117.83	123.04
1	B	201	GLN	C-N-CA	-5.66	117.83	123.04
1	B	243	ASP	N-CA-C	5.49	117.34	111.36
1	A	47	LYS	CB-CA-C	-5.46	102.14	111.26
1	B	112	ARG	NE-CZ-NH2	5.41	124.07	119.20
1	B	41	ARG	NE-CZ-NH1	-5.38	116.12	121.50
1	A	53	THR	O-C-N	5.37	129.34	123.22
1	B	273	LEU	N-CA-C	-5.37	101.72	110.20
1	B	268	ASP	CB-CA-C	5.22	117.79	109.96
1	A	178	PHE	CA-CB-CG	5.17	118.97	113.80
1	B	183	GLN	CA-C-O	-5.16	114.51	120.24
1	A	113	TYR	CD1-CE1-CZ	-5.03	110.55	119.60

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	121[A]	GLU	Mainchain
1	A	124[A]	PRO	Peptide,Mainchain
1	A	154	ASN	Mainchain
1	B	121[A]	GLU	Mainchain
1	B	123[A]	ALA	Mainchain

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	2365	0	2295	76	7
1	B	2353	0	2279	61	9
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	234	0	0	17	0
3	B	263	0	0	17	0
All	All	5217	0	4574	137	16

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 15.

All (137) 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:98:LYS:C	1:A:99:ASN:N	1.76	1.42
1:B:156:LEU:HG	3:B:743:HOH:O	1.16	1.30
1:A:97:LYS:C	1:A:98:LYS:N	2.17	1.03
1:A:123[A]:ALA:HB3	1:A:126:VAL:HG21	1.46	0.95
1:A:98:LYS:O	1:A:99:ASN:HB2	1.71	0.91
1:A:243:ASP:H	1:A:275[B]:GLN:HE22	1.15	0.90
1:A:98:LYS:C	1:A:99:ASN:HB2	1.97	0.89
1:A:243:ASP:H	1:A:275[A]:GLN:HE22	1.21	0.89
1:B:273:LEU:O	1:B:274:ARG:HB2	1.74	0.87
1:B:156:LEU:HB3	3:B:849:HOH:O	1.75	0.87
1:A:98:LYS:C	1:A:99:ASN:CB	2.48	0.86
1:A:123[A]:ALA:C	1:A:126:VAL:HG22	2.01	0.84
1:A:98:LYS:C	1:A:99:ASN:CA	2.51	0.83
1:A:123[A]:ALA:HB3	1:A:126:VAL:CG2	2.09	0.82
1:B:273:LEU:O	1:B:274:ARG:CB	2.29	0.80
1:A:224[A]:VAL:HB	3:A:793:HOH:O	1.84	0.76
1:A:240:LEU:HD22	1:A:245:THR:CG2	2.16	0.76
1:A:58[A]:VAL:HG22	1:A:73:ILE:HG12	1.69	0.74
1:A:240:LEU:HD22	1:A:245:THR:HG21	1.67	0.74
1:A:123[A]:ALA:O	1:A:126:VAL:HG22	1.86	0.73
1:B:93:VAL:HA	3:B:802:HOH:O	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:LYS:O	1:B:98:LYS:HG3	1.89	0.72
1:B:270:GLU:O	1:B:273:LEU:O	2.08	0.72
1:A:58[B]:VAL:HG12	1:A:73:ILE:HG12	1.72	0.72
1:B:61:VAL:HG13	3:B:792:HOH:O	1.91	0.71
1:A:99:ASN:ND2	1:A:120:GLU:HB3	2.06	0.70
1:A:95:ASN:H	1:A:95:ASN:ND2	1.88	0.69
1:B:106:LYS:HB3	3:B:743:HOH:O	1.91	0.69
1:A:98:LYS:CA	1:A:99:ASN:N	2.56	0.69
1:A:271:GLY:O	1:A:275[B]:GLN:HG2	1.91	0.68
1:A:243:ASP:H	1:A:275[B]:GLN:NE2	1.91	0.67
1:A:85[B]:GLN:NE2	3:A:411:HOH:O	2.27	0.67
1:A:248:CYS:O	3:A:804:HOH:O	2.12	0.66
1:A:99:ASN:HD21	1:A:120:GLU:HB3	1.59	0.66
1:A:240:LEU:HD21	1:A:260:VAL:HG11	1.78	0.66
1:A:101:ARG:HD3	1:A:120:GLU:OE1	1.96	0.65
1:A:123[A]:ALA:N	1:A:126:VAL:CG2	2.61	0.64
1:A:123[A]:ALA:H	1:A:126:VAL:CG2	2.11	0.63
1:B:21[A]:VAL:HB	3:B:792:HOH:O	1.97	0.63
1:A:123[A]:ALA:CA	1:A:126:VAL:HG22	2.29	0.62
1:B:264:ARG:HD3	3:B:762:HOH:O	1.99	0.61
1:A:4:ILE:HD13	1:A:238:VAL:HG13	1.83	0.61
1:A:240:LEU:CD2	1:A:260:VAL:HG11	2.30	0.61
1:A:123[A]:ALA:H	1:A:126:VAL:HG23	1.65	0.60
1:B:54[A]:MET:HE1	1:B:80:LEU:HD13	1.83	0.60
1:A:30:LEU:HD23	1:A:41:ARG:HG2	1.85	0.59
1:A:98:LYS:O	1:A:98:LYS:HG2	2.02	0.59
1:A:207:CYS:HB3	3:A:804:HOH:O	2.03	0.59
1:A:130:HIS:HD2	3:A:597:HOH:O	1.86	0.58
1:A:123[A]:ALA:CB	1:A:126:VAL:CG2	2.82	0.57
1:A:98:LYS:N	1:A:99:ASN:N	2.53	0.56
1:A:215:TRP:HB3	3:A:793:HOH:O	2.06	0.55
1:A:84:GLU:O	1:A:85[B]:GLN:HB2	2.05	0.55
1:A:268:ASP:O	1:A:269:PRO:C	2.39	0.55
1:B:21[B]:VAL:HG13	1:B:30:LEU:HB2	1.88	0.55
1:A:243:ASP:N	1:A:275[B]:GLN:HE22	1.97	0.54
1:A:264:ARG:HB3	3:A:681:HOH:O	2.07	0.54
1:A:186:GLN:HG3	3:A:799:HOH:O	2.07	0.54
1:A:275[A]:GLN:NE2	3:A:380:HOH:O	2.40	0.53
1:A:220:ASN:HD21	1:A:244:LYS:NZ	2.06	0.53
1:A:123[A]:ALA:CB	1:A:126:VAL:HG21	2.31	0.53
1:A:262[A]:CYS:HB3	3:A:760:HOH:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LEU:HD11	1:A:69:TYR:CE1	2.43	0.53
1:A:240:LEU:HD22	1:A:245:THR:HG22	1.90	0.53
1:B:126:VAL:HG22	1:B:126:VAL:O	2.09	0.53
1:A:123[A]:ALA:CA	1:A:126:VAL:CG2	2.86	0.52
1:A:242:VAL:HA	1:A:275[B]:GLN:NE2	2.26	0.51
1:A:243:ASP:H	1:A:275[A]:GLN:NE2	1.98	0.51
1:B:240:LEU:CD2	1:B:260:VAL:HG11	2.40	0.51
1:B:124[A]:PRO:O	1:B:127:LEU:HD22	2.10	0.50
1:B:30:LEU:HB2	3:B:792:HOH:O	2.11	0.50
1:A:54:MET:HG3	1:A:58[A]:VAL:HG21	1.92	0.50
1:A:81:ASN:OD1	1:A:83:LYS:HB2	2.11	0.50
1:A:101:ARG:CD	1:A:120:GLU:OE1	2.59	0.49
1:B:273:LEU:O	1:B:274:ARG:CG	2.61	0.49
1:B:101:ARG:HD3	1:B:120:GLU:OE1	2.12	0.48
1:B:83:LYS:HB2	3:B:629:HOH:O	2.13	0.48
1:A:189:ASN:HA	3:A:593:HOH:O	2.14	0.47
1:A:244:LYS:O	1:A:262[B]:CYS:HB3	2.13	0.47
1:A:275[A]:GLN:HA	1:A:276:PRO:HD2	1.73	0.46
1:A:54:MET:HG3	1:A:58[B]:VAL:HG11	1.97	0.46
1:B:215:TRP:CD2	1:B:258:MET:HE1	2.51	0.46
1:B:41:ARG:NH1	3:B:653:HOH:O	2.49	0.45
1:B:65:GLN:NE2	3:B:850:HOH:O	2.49	0.45
1:A:275[B]:GLN:NE2	3:A:380:HOH:O	2.50	0.45
1:B:81:ASN:ND2	1:B:84:GLU:HG3	2.32	0.45
1:A:65:GLN:NE2	3:A:374:HOH:O	2.49	0.45
1:A:81:ASN:ND2	1:A:84:GLU:OE1	2.50	0.45
1:A:200:GLU:O	1:A:218:CYS:HB3	2.17	0.45
1:A:216:VAL:N	3:A:793:HOH:O	2.50	0.44
1:A:213:LYS:NZ	3:A:834:HOH:O	2.49	0.44
1:B:175:VAL:O	1:B:193:VAL:HG22	2.17	0.44
1:B:240:LEU:HD21	1:B:260:VAL:HG11	2.00	0.44
1:A:126:VAL:HG23	1:A:126:VAL:O	2.17	0.44
1:A:98:LYS:O	1:A:98:LYS:CG	2.66	0.44
1:B:239:LYS:NZ	3:B:618:HOH:O	2.50	0.44
1:B:27:ASN:ND2	3:B:815:HOH:O	2.50	0.44
1:B:37:LYS:HE3	1:B:54[A]:MET:O	2.18	0.43
1:B:240:LEU:HD11	1:B:284:ILE:HD11	1.99	0.43
1:B:55:ASP:HB2	3:B:641:HOH:O	2.18	0.43
1:B:273:LEU:O	1:B:274:ARG:HG3	2.19	0.43
1:A:206:MET:C	3:A:804:HOH:O	2.61	0.43
1:B:220:ASN:HD22	1:B:220:ASN:HA	1.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:ARG:HH11	1:B:41:ARG:HD3	1.65	0.42
1:B:215:TRP:CE3	1:B:258:MET:HE1	2.54	0.42
1:B:240:LEU:HD22	1:B:245:THR:CG2	2.49	0.42
1:B:220:ASN:HD21	1:B:244:LYS:NZ	2.18	0.42
1:B:264:ARG:O	1:B:267[A]:MET:HG3	2.20	0.42
1:B:27:ASN:ND2	3:B:774:HOH:O	2.49	0.41
1:B:38:LYS:HE3	1:B:40:CYS:SG	2.60	0.41
1:A:12:GLU:O	1:A:12:GLU:HG3	2.21	0.41
1:A:271:GLY:O	1:A:275[A]:GLN:HG2	2.20	0.41
1:B:44:SER:HB2	3:B:845:HOH:O	2.20	0.41
1:A:220:ASN:HD22	1:A:220:ASN:HA	1.66	0.41
1:A:130:HIS:HE1	1:B:128:GLU:OE2	2.04	0.41
1:A:51:ARG:NH1	3:A:495:HOH:O	2.55	0.40
1:B:219:TYR:O	1:B:220:ASN:HB2	2.22	0.40

All (16) 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:98:LYS:NZ	1:A:146:TYR:C[2_745]	0.34	1.86
1:B:197:GLU:OE2	1:B:274:ARG:CZ[2_746]	0.38	1.82
1:A:98:LYS:CE	1:A:146:TYR:O[2_745]	0.39	1.81
1:B:197:GLU:CD	1:B:274:ARG:NH2[2_746]	0.53	1.67
1:A:98:LYS:NZ	1:A:146:TYR:O[2_745]	1.14	1.06
1:B:197:GLU:OE2	1:B:274:ARG:NH2[2_746]	1.16	1.04
1:B:197:GLU:CG	1:B:274:ARG:NH2[2_746]	1.29	0.91
1:B:197:GLU:OE2	1:B:274:ARG:NE[2_746]	1.34	0.86
1:A:98:LYS:NZ	1:A:146:TYR:CA[2_745]	1.36	0.84
1:B:197:GLU:CD	1:B:274:ARG:CZ[2_746]	1.47	0.73
1:A:98:LYS:CE	1:A:146:TYR:C[2_745]	1.58	0.62
1:B:197:GLU:OE2	1:B:274:ARG:NH1[2_746]	1.61	0.59
1:A:98:LYS:NZ	1:A:147:PHE:N[2_745]	1.64	0.56
1:A:98:LYS:CD	1:A:146:TYR:O[2_745]	1.68	0.52
1:B:197:GLU:OE1	1:B:274:ARG:NH2[2_746]	1.75	0.45
1:B:197:GLU:CD	1:B:274:ARG:NE[2_746]	2.13	0.07

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	300/297 (101%)	290 (97%)	8 (3%)	2 (1%)	18	4
1	B	300/297 (101%)	284 (95%)	13 (4%)	3 (1%)	12	2
All	All	600/594 (101%)	574 (96%)	21 (4%)	5 (1%)	14	3

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	ASN
1	B	274	ARG
1	B	124[A]	PRO
1	A	154	ASN
1	B	123[A]	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	A	259/252 (103%)	257 (99%)	2 (1%)	73	47
1	B	257/252 (102%)	255 (99%)	2 (1%)	73	47
All	All	516/504 (102%)	512 (99%)	4 (1%)	73	47

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	ASN

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Mol	Chain	Res	Type
1	A	171	LEU
1	B	126	VAL
1	B	273	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	65	GLN
1	A	95	ASN
1	A	99	ASN
1	A	130	HIS
1	A	142	HIS
1	A	220	ASN
1	A	254	ASN
1	B	27	ASN
1	B	95	ASN
1	B	99	ASN
1	B	131	GLN
1	B	141	HIS
1	B	142	HIS
1	B	220	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	10
1	B	5

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	97:LYS	C	98:LYS	N	2.17
1	A	98:LYS	C	99:ASN	N	1.76
1	A	82:TRP	C	83:LYS	N	1.20
1	A	195:LYS	C	196:LEU	N	1.20
1	B	197:GLU	C	198:LYS	N	1.19
1	A	41:ARG	C	42:TRP	N	1.18
1	A	182:LEU	C	183:GLN	N	1.17
1	B	185:GLY	C	186:GLN	N	1.15
1	A	268:ASP	C	269:PRO	N	1.11
1	B	182:LEU	C	183:GLN	N	1.09
1	A	120:GLU	C	121:GLU	N	1.06
1	B	269:PRO	C	270:GLU	N	1.05
1	B	148:ASP	C	149:GLN	N	1.02
1	A	269:PRO	C	270:GLU	N	1.00
1	A	272:LEU	C	273:LEU	N	0.95

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	297/297 (100%)	0.01	25 (8%)	17 18	8, 21, 39, 79	9 (3%)
1	B	297/297 (100%)	-0.12	20 (6%)	24 26	6, 20, 37, 73	6 (2%)
All	All	594/594 (100%)	-0.05	45 (7%)	20 21	6, 20, 39, 79	15 (2%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	123[A]	ALA	15.8
1	B	124[A]	PRO	11.8
1	A	124[A]	PRO	11.4
1	A	126	VAL	11.2
1	B	122[A]	THR	9.7
1	A	270	GLU	9.6
1	A	122[A]	THR	8.7
1	A	268	ASP	8.1
1	A	273	LEU	7.6
1	A	98	LYS	7.1
1	B	274	ARG	7.0
1	B	121[A]	GLU	6.6
1	A	123[A]	ALA	6.1
1	B	98	LYS	6.1
1	A	83	LYS	6.0
1	A	121[A]	GLU	5.8
1	A	127	LEU	5.5
1	B	253	LYS	5.1
1	A	99	ASN	4.7
1	A	274	ARG	4.6
1	B	197	GLU	4.5
1	B	126	VAL	4.4
1	A	199	GLU	4.4
1	B	45	PHE	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	270	GLU	4.2
1	B	273	LEU	4.0
1	B	198	LYS	3.6
1	A	120	GLU	3.5
1	B	186	GLN	3.4
1	B	183	GLN	3.2
1	B	125[A]	ALA	3.1
1	A	299[A]	GLY	3.0
1	A	183	GLN	3.0
1	A	125[A]	ALA	2.8
1	A	50	GLN	2.8
1	A	262[A]	CYS	2.6
1	A	269	PRO	2.6
1	A	271	GLY	2.4
1	A	154	ASN	2.4
1	B	120	GLU	2.3
1	A	86	SER	2.1
1	B	149	GLN	2.1
1	A	195	LYS	2.1
1	B	54[A]	MET	2.1
1	B	240	LEU	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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	CA	A	1	1/1	1.00	0.01	16,16,16,16	0
2	CA	B	1	1/1	1.00	0.01	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.