



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 07:50 PM EDT

PDB ID : 4HGK / pdb_00004hgk
Title : Shark IgNAR variable domain
Authors : Olland, A.O.; Kovalenko, O.V.; Svenson, K.; King, D.
Deposited on : 2012-10-08
Resolution : 3.04 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

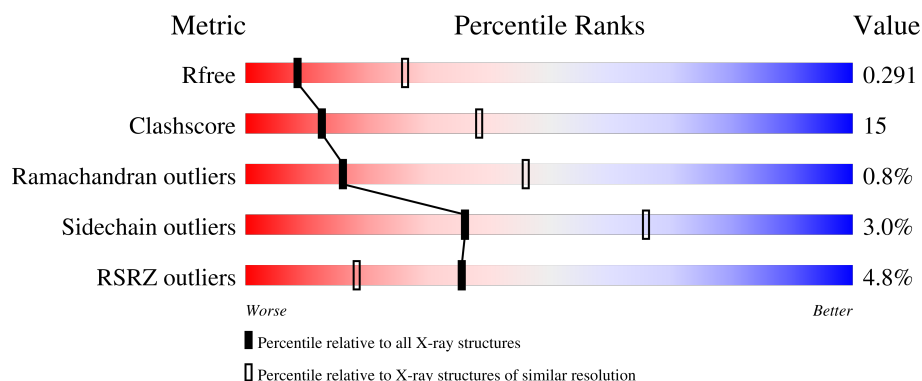
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.04 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	3685 (3.08-3.00)
Clashscore	190562	4007 (3.08-3.00)
Ramachandran outliers	187476	3834 (3.08-3.00)
Sidechain outliers	187428	3836 (3.08-3.00)
RSRZ outliers	180081	3684 (3.08-3.00)

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	585	 4% 63% 18% 17%
1	B	585	 4% 63% 17% 18%
2	C	128	 2% 60% 19% 19%
2	D	128	 0% 63% 16% 19%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9335 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 Serum albumin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	483	Total	C	N	O	S	0	0	0
			3846	2428	652	730	36			
1	B	481	Total	C	N	O	S	0	0	0
			3838	2423	651	728	36			

- Molecule 2 is a protein called shark V-NAR antibody.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	104	Total	C	N	O	S	0	0	0
			806	498	142	163	3			
2	D	104	Total	C	N	O	S	0	0	0
			806	498	142	163	3			

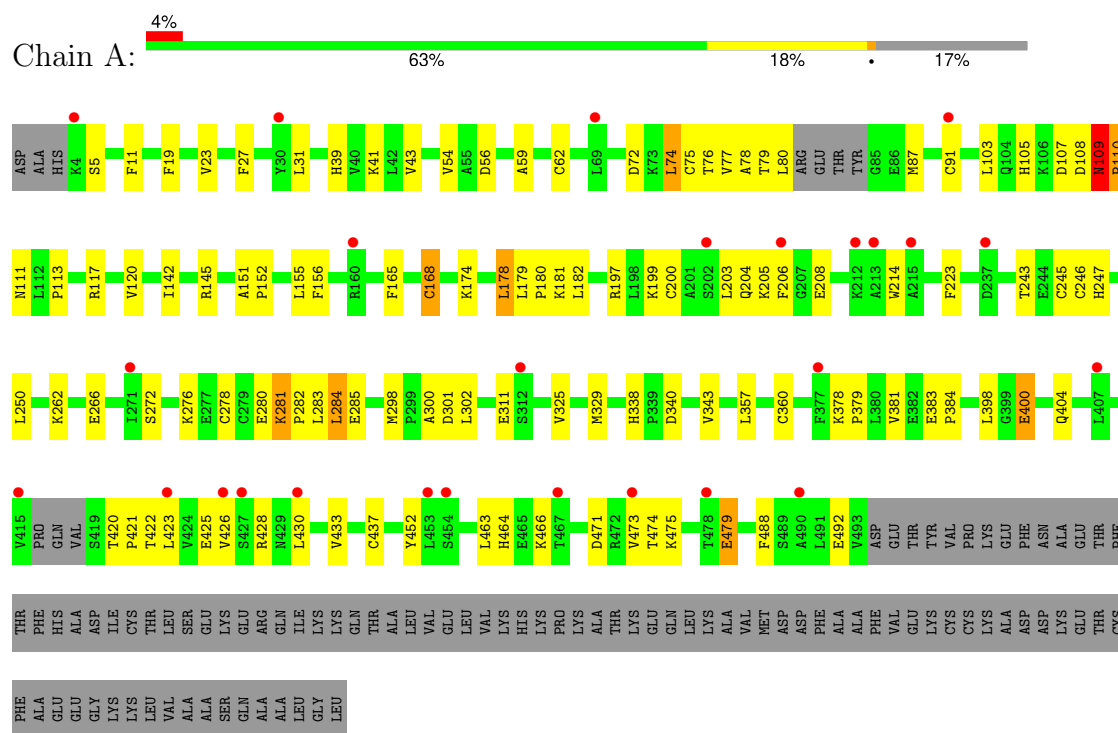
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total	O	0	0
			6	6		
3	B	19	Total	O	0	0
			19	19		
3	C	8	Total	O	0	0
			8	8		
3	D	6	Total	O	0	0
			6	6		

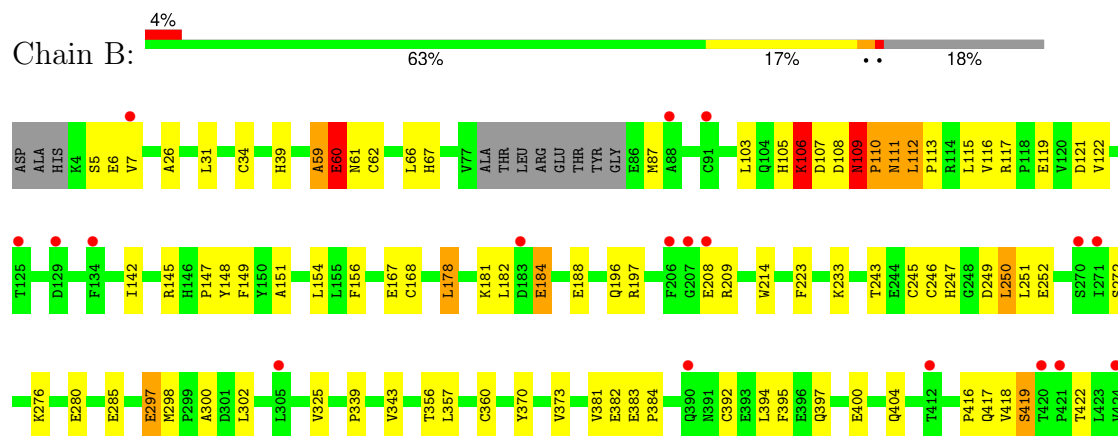
3 Residue-property plots [i](#)

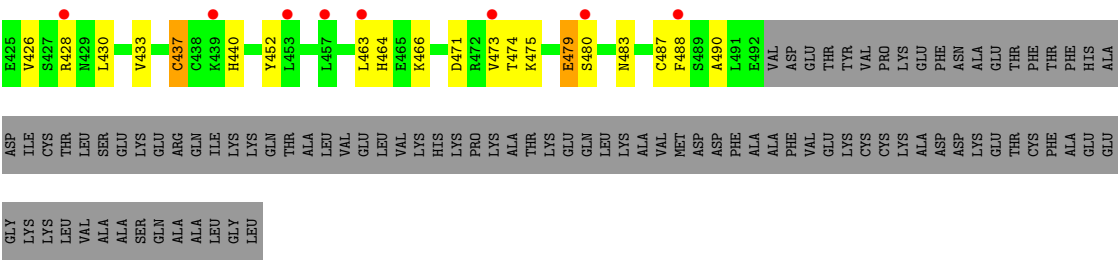
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Serum albumin

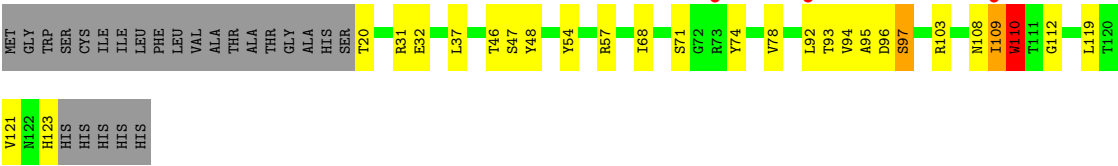


• Molecule 1: Serum albumin





• Molecule 2: shark V-NAR antibody



• Molecule 2: shark V-NAR antibody



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	127.98Å 127.98Å 151.76Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	52.05 – 3.04 52.05 – 3.04	Depositor EDS
% Data completeness (in resolution range)	92.1 (52.05-3.04) 92.2 (52.05-3.04)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 3.07Å)	Xtriage
Refinement program	BUSTER 2.9.3	Depositor
R, R_{free}	0.237 , 0.266 0.259 , 0.291	Depositor DCC
R_{free} test set	1317 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	68.3	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 65.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9335	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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](#)

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	0.67	0/3920	1.30	15/5287 (0.3%)
1	B	0.73	1/3914 (0.0%)	1.45	24/5281 (0.5%)
2	C	0.81	0/820	1.17	5/1115 (0.4%)
2	D	0.81	0/820	1.15	2/1115 (0.2%)
All	All	0.72	1/9474 (0.0%)	1.34	46/12798 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	490	ALA	C-N	5.19	1.40	1.33

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	LEU	N-CA-C	11.39	123.44	111.14
1	A	110	PRO	N-CA-C	-11.30	92.84	111.26
1	B	106	LYS	N-CA-C	-11.13	93.81	110.28
1	B	178	LEU	N-CA-C	9.46	121.59	111.28
1	B	250	LEU	N-CA-C	8.54	120.59	111.28
1	B	111	ASN	N-CA-C	7.26	120.76	110.23
1	B	110	PRO	N-CA-C	-7.16	99.59	111.26
1	A	79	THR	N-CA-C	-6.80	103.87	111.28
2	C	110	TRP	N-CA-C	-6.61	100.25	110.10
1	B	184	GLU	CA-C-N	6.44	128.91	120.28
1	B	184	GLU	C-N-CA	6.44	128.91	120.28
1	B	60	GLU	N-CA-C	6.43	118.29	111.28
1	A	437	CYS	N-CA-C	6.29	118.21	111.36
2	C	95	ALA	N-CA-C	-6.16	104.86	112.38
1	A	208	GLU	N-CA-C	6.15	117.98	111.28
1	B	416	PRO	CA-C-N	6.13	128.77	120.38
1	B	416	PRO	C-N-CA	6.13	128.77	120.38
1	A	56	ASP	CA-CB-CG	6.07	118.67	112.60
1	B	339	PRO	CA-C-N	6.02	129.47	120.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	339	PRO	C-N-CA	6.02	129.47	120.31
2	C	48	TYR	N-CA-C	5.98	117.35	109.93
1	B	121	ASP	CA-CB-CG	5.85	118.45	112.60
1	A	109	ASN	CA-C-N	-5.70	113.83	119.93
1	A	109	ASN	C-N-CA	-5.70	113.83	119.93
1	B	34	CYS	N-CA-C	5.70	117.07	109.83
2	D	48	TYR	N-CA-C	5.59	117.11	110.07
1	B	168	CYS	N-CA-C	5.58	118.29	111.82
1	B	437	CYS	N-CA-C	5.55	118.52	111.69
1	A	168	CYS	N-CA-C	5.48	118.18	111.82
1	B	417	GLN	N-CA-C	5.43	117.90	111.33
1	A	250	LEU	N-CA-C	5.41	117.60	111.11
2	D	47	SER	N-CA-C	-5.40	106.73	113.38
2	C	47	SER	N-CA-C	-5.39	106.75	113.38
1	B	6	GLU	CB-CG-CD	5.37	121.73	112.60
1	B	382	GLU	N-CA-C	5.36	117.12	111.28
1	A	301	ASP	CA-CB-CG	5.27	117.87	112.60
1	A	398	LEU	CD1-CG-CD2	-5.22	99.32	110.80
1	A	360	CYS	N-CA-C	5.16	116.90	111.28
2	C	32	GLU	N-CA-C	-5.14	102.45	110.42
1	A	109	ASN	N-CA-C	5.08	121.03	109.81
1	B	59	ALA	N-CA-C	-5.06	103.80	110.33
1	B	487	CYS	CA-C-N	5.04	127.29	120.44
1	B	487	CYS	C-N-CA	5.04	127.29	120.44
1	B	360	CYS	N-CA-C	5.02	116.75	111.28
1	A	338	HIS	N-CA-C	5.02	117.77	108.94
1	B	109	ASN	N-CA-C	5.01	120.89	109.81

There are no chirality outliers.

There are no planarity outliers.

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	3846	0	3775	117	0
1	B	3838	0	3763	118	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	806	0	790	33	0
2	D	806	0	790	35	1
3	A	6	0	0	0	0
3	B	19	0	0	0	0
3	C	8	0	0	0	0
3	D	6	0	0	0	0
All	All	9335	0	9118	282	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (282) 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:B:325:VAL:HG22	2:C:109:ILE:CD1	1.52	1.38
1:B:59:ALA:HB3	1:B:62:CYS:SG	1.74	1.25
1:A:74:LEU:O	1:A:77:VAL:HG12	1.33	1.24
1:A:72:ASP:O	1:A:76:THR:HG23	1.34	1.20
1:A:27:PHE:HD1	1:A:74:LEU:HD21	1.06	1.20
1:B:325:VAL:CA	2:C:109:ILE:HD11	1.70	1.20
1:A:27:PHE:CD1	1:A:74:LEU:HD21	1.77	1.18
1:A:75:CYS:SG	1:A:91:CYS:SG	1.29	1.15
1:A:325:VAL:HA	2:D:109:ILE:HD11	1.19	1.14
1:B:107:ASP:HB2	1:B:110:PRO:HG2	1.25	1.14
1:A:325:VAL:HG22	2:D:109:ILE:CD1	1.78	1.13
1:B:178:LEU:HD21	1:B:182:LEU:CD1	1.82	1.10
1:B:325:VAL:HA	2:C:109:ILE:HD11	1.08	1.08
1:A:325:VAL:HG22	2:D:109:ILE:HD13	1.32	1.06
1:A:325:VAL:CA	2:D:109:ILE:HD11	1.87	1.04
1:B:325:VAL:CG2	2:C:109:ILE:CD1	2.35	1.04
1:B:106:LYS:HD3	1:B:147:PRO:HB2	1.40	1.02
2:D:20:THR:HG22	2:D:110:TRP:CH2	1.95	1.02
1:B:325:VAL:HA	2:C:109:ILE:CD1	1.90	1.01
1:B:325:VAL:CB	2:C:109:ILE:HD11	1.92	0.99
1:B:325:VAL:CG2	2:C:109:ILE:HD11	1.95	0.97
1:B:26:ALA:HB2	1:B:250:LEU:HD12	1.45	0.97
1:B:325:VAL:HG22	2:C:109:ILE:HD13	1.00	0.97
1:B:107:ASP:HB2	1:B:110:PRO:CG	1.95	0.96
1:B:325:VAL:HG22	2:C:109:ILE:HD11	1.48	0.95
1:A:75:CYS:CB	1:A:91:CYS:SG	2.55	0.94
1:B:178:LEU:HD21	1:B:182:LEU:CG	1.98	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:CYS:O	1:A:204:GLN:HG3	1.67	0.93
1:B:243:THR:O	1:B:247:HIS:HB2	1.68	0.93
1:B:178:LEU:C	1:B:178:LEU:HD23	1.93	0.93
1:A:107:ASP:HB2	1:A:110:PRO:CG	1.98	0.92
1:A:27:PHE:CD1	1:A:74:LEU:CD2	2.52	0.92
1:A:179:LEU:N	1:A:180:PRO:HD3	1.84	0.92
1:B:404:GLN:HE21	1:B:428:ARG:HA	1.34	0.90
1:A:109:ASN:O	1:A:110:PRO:C	2.11	0.90
1:A:404:GLN:HE21	1:A:428:ARG:HA	1.36	0.90
1:A:200:CYS:O	1:A:204:GLN:CG	2.19	0.89
1:B:356:THR:HG21	1:B:373:VAL:HG22	1.55	0.89
1:B:178:LEU:CD2	1:B:182:LEU:HG	2.04	0.88
1:B:66:LEU:HD13	1:B:251:LEU:HD12	1.55	0.87
1:A:179:LEU:N	1:A:180:PRO:CD	2.37	0.86
1:B:178:LEU:HD21	1:B:182:LEU:HG	1.56	0.86
1:A:107:ASP:HB2	1:A:110:PRO:HG2	1.57	0.86
2:C:94:VAL:HA	2:C:121:VAL:HB	1.58	0.86
1:B:107:ASP:CB	1:B:110:PRO:CG	2.56	0.84
2:D:20:THR:HG22	2:D:110:TRP:CZ3	2.13	0.83
1:A:325:VAL:HA	2:D:109:ILE:CD1	2.06	0.83
1:A:75:CYS:SG	1:A:91:CYS:CB	2.66	0.82
1:A:283:LEU:HD12	1:A:284:LEU:N	1.95	0.82
1:B:59:ALA:CB	1:B:62:CYS:SG	2.64	0.82
1:A:107:ASP:CB	1:A:110:PRO:CG	2.58	0.81
1:B:325:VAL:CG2	2:C:109:ILE:HD13	1.97	0.81
2:D:20:THR:HB	2:D:110:TRP:CZ2	2.16	0.81
1:B:178:LEU:HD21	1:B:182:LEU:HD11	1.60	0.81
1:B:178:LEU:HD23	1:B:178:LEU:O	1.81	0.80
2:C:109:ILE:O	2:C:109:ILE:HG22	1.82	0.80
1:A:325:VAL:HG22	2:D:109:ILE:HD11	1.58	0.80
1:A:422:THR:HB	1:A:463:LEU:HD21	1.63	0.80
1:A:325:VAL:CG2	2:D:109:ILE:HD11	2.12	0.79
1:A:107:ASP:HB3	1:A:110:PRO:HD3	1.65	0.79
1:B:66:LEU:CD1	1:B:251:LEU:HD12	2.11	0.79
2:D:93:THR:HG22	2:D:95:ALA:H	1.47	0.78
1:B:107:ASP:CB	1:B:110:PRO:HG2	2.10	0.78
1:B:108:ASP:OD2	1:B:197:ARG:HD3	1.85	0.77
2:D:93:THR:HG22	2:D:94:VAL:N	1.98	0.77
1:A:107:ASP:CB	1:A:110:PRO:CD	2.63	0.77
1:B:356:THR:HG21	1:B:373:VAL:CG2	2.15	0.77
1:B:178:LEU:CD2	1:B:182:LEU:CD1	2.62	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:VAL:CB	2:D:109:ILE:HD11	2.15	0.75
1:B:243:THR:O	1:B:247:HIS:CB	2.34	0.75
1:A:77:VAL:O	1:A:80:LEU:C	2.30	0.75
2:D:20:THR:HG23	2:D:46:THR:HB	1.69	0.74
1:A:178:LEU:C	1:A:180:PRO:HD2	2.12	0.74
1:A:120:VAL:CG1	1:A:174:LYS:HB3	2.18	0.74
1:B:247:HIS:HD2	1:B:249:ASP:OD2	1.71	0.74
1:B:115:LEU:HD12	1:B:116:VAL:H	1.53	0.73
2:C:20:THR:HG23	2:C:46:THR:HB	1.71	0.73
1:A:27:PHE:CE1	1:A:74:LEU:CD2	2.71	0.72
1:A:31:LEU:HD11	1:A:74:LEU:HD22	1.70	0.72
1:B:26:ALA:CB	1:B:250:LEU:HD12	2.20	0.72
1:B:5:SER:HB2	1:B:62:CYS:O	1.90	0.71
1:B:370:TYR:O	1:B:373:VAL:HG23	1.90	0.71
2:D:20:THR:CG2	2:D:110:TRP:CH2	2.73	0.71
1:A:107:ASP:HB2	1:A:110:PRO:CD	2.21	0.70
1:A:325:VAL:CG2	2:D:109:ILE:CD1	2.62	0.70
1:A:27:PHE:CE1	1:A:74:LEU:HD23	2.26	0.70
1:A:178:LEU:C	1:A:180:PRO:CD	2.63	0.70
1:A:471:ASP:O	1:A:474:THR:HG22	1.91	0.70
1:B:107:ASP:O	1:B:110:PRO:HD2	1.92	0.70
2:D:93:THR:CG2	2:D:94:VAL:N	2.54	0.70
2:C:110:TRP:HE3	2:C:110:TRP:O	1.75	0.69
2:D:109:ILE:HG22	2:D:109:ILE:O	1.90	0.69
1:A:77:VAL:HG13	1:A:78:ALA:N	2.08	0.67
1:A:120:VAL:HG11	1:A:174:LYS:HB3	1.75	0.67
1:A:107:ASP:HB3	1:A:110:PRO:CD	2.23	0.67
1:A:204:GLN:HE21	1:A:246:CYS:HB3	1.59	0.67
1:B:87:MET:HE1	1:B:105:HIS:HB3	1.78	0.66
1:A:298:MET:HE3	1:A:302:LEU:HD12	1.76	0.66
1:B:298:MET:HE3	1:B:302:LEU:HD12	1.76	0.66
1:B:464:HIS:NE2	1:B:474:THR:HB	2.12	0.65
1:A:433:VAL:HG22	1:A:452:TYR:CE2	2.32	0.65
1:A:87:MET:HE1	1:A:105:HIS:HB3	1.78	0.65
1:A:168:CYS:HB2	1:A:178:LEU:HD13	1.78	0.64
1:B:109:ASN:HB3	1:B:466:LYS:HE2	1.78	0.64
1:A:117:ARG:NH1	1:A:182:LEU:HB3	2.13	0.64
1:A:109:ASN:O	1:A:111:ASN:N	2.30	0.64
2:C:110:TRP:CE3	2:C:110:TRP:C	2.76	0.64
1:B:178:LEU:CD2	1:B:182:LEU:CG	2.69	0.63
1:B:433:VAL:HG22	1:B:452:TYR:CE2	2.33	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:471:ASP:O	1:B:474:THR:HG22	1.98	0.63
1:A:464:HIS:NE2	1:A:474:THR:HB	2.14	0.62
2:D:20:THR:CG2	2:D:110:TRP:CZ3	2.82	0.62
2:D:108:ASN:O	2:D:109:ILE:HB	2.00	0.61
1:B:109:ASN:HB3	1:B:466:LYS:HZ3	1.66	0.61
1:A:77:VAL:CG1	1:A:78:ALA:N	2.64	0.61
1:A:381:VAL:O	1:A:384:PRO:HD2	2.00	0.60
1:B:115:LEU:HD12	1:B:116:VAL:N	2.15	0.60
2:C:31:ARG:HB2	2:C:92:LEU:CD2	2.31	0.60
1:B:109:ASN:O	1:B:110:PRO:C	2.42	0.60
2:C:110:TRP:CH2	2:C:112:GLY:HA3	2.37	0.60
1:B:381:VAL:O	1:B:384:PRO:HD2	2.01	0.60
1:B:107:ASP:HB3	1:B:110:PRO:CG	2.32	0.60
2:C:109:ILE:O	2:C:109:ILE:CG2	2.50	0.60
1:A:433:VAL:HG22	1:A:452:TYR:CD2	2.37	0.59
2:D:54:TYR:HB2	2:D:103:ARG:HB3	1.84	0.59
1:A:205:LYS:HD3	1:A:206:PHE:CE2	2.37	0.59
1:B:249:ASP:HB3	1:B:252:GLU:CD	2.27	0.59
1:B:151:ALA:CB	1:B:250:LEU:HD22	2.33	0.59
2:C:93:THR:O	2:C:96:ASP:HB2	2.03	0.58
1:B:106:LYS:CD	1:B:147:PRO:HB2	2.24	0.58
2:C:54:TYR:HB2	2:C:103:ARG:HB3	1.84	0.58
1:B:433:VAL:HG22	1:B:452:TYR:CD2	2.39	0.58
2:D:108:ASN:OD1	2:D:108:ASN:N	2.33	0.57
1:B:109:ASN:CB	1:B:466:LYS:HE2	2.35	0.57
1:A:107:ASP:CB	1:A:110:PRO:HD3	2.31	0.56
1:B:178:LEU:HD23	1:B:182:LEU:HG	1.85	0.56
1:B:31:LEU:HB2	1:B:39:HIS:HE1	1.69	0.56
1:B:67:HIS:NE2	1:B:249:ASP:OD1	2.38	0.56
1:A:75:CYS:CA	1:A:91:CYS:SG	2.93	0.56
1:B:109:ASN:HB3	1:B:466:LYS:CE	2.36	0.56
1:A:168:CYS:HB2	1:A:178:LEU:CD1	2.35	0.55
1:A:120:VAL:HG12	1:A:174:LYS:HB3	1.86	0.55
1:A:200:CYS:O	1:A:204:GLN:HG2	2.04	0.55
1:A:27:PHE:CE1	1:A:74:LEU:HD21	2.35	0.55
1:B:107:ASP:CB	1:B:110:PRO:CD	2.85	0.55
1:B:107:ASP:C	1:B:110:PRO:HD2	2.31	0.55
1:A:109:ASN:HB3	1:A:466:LYS:HE2	1.89	0.54
1:A:426:VAL:O	1:A:430:LEU:HG	2.07	0.54
1:B:103:LEU:HD11	1:B:247:HIS:O	2.08	0.54
1:A:404:GLN:NE2	1:A:428:ARG:HA	2.16	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:110:TRP:O	2:C:110:TRP:CE3	2.59	0.54
1:B:60:GLU:HG2	1:B:61:ASN:OD1	2.08	0.54
2:C:92:LEU:HD12	2:C:96:ASP:OD2	2.06	0.54
1:B:107:ASP:HB3	1:B:110:PRO:CD	2.37	0.54
1:B:66:LEU:HD13	1:B:251:LEU:CD1	2.34	0.54
1:A:200:CYS:C	1:A:204:GLN:HG3	2.32	0.54
1:A:278:CYS:HA	1:A:281:LYS:HG3	1.90	0.54
1:B:178:LEU:C	1:B:178:LEU:CD2	2.68	0.54
1:B:418:VAL:HG12	1:B:419:SER:N	2.23	0.54
1:B:404:GLN:NE2	1:B:428:ARG:HA	2.14	0.53
1:B:178:LEU:CD2	1:B:182:LEU:HD12	2.38	0.53
1:A:19:PHE:O	1:A:23:VAL:HG23	2.08	0.53
1:A:281:LYS:HB3	1:A:282:PRO:HD2	1.90	0.53
1:B:60:GLU:CG	1:B:61:ASN:OD1	2.57	0.53
2:D:20:THR:CB	2:D:110:TRP:CZ2	2.90	0.53
1:B:109:ASN:HB3	1:B:466:LYS:NZ	2.24	0.53
1:A:281:LYS:HB3	1:A:282:PRO:CD	2.38	0.52
1:A:109:ASN:HB3	1:A:466:LYS:NZ	2.25	0.52
2:D:96:ASP:O	2:D:97:SER:C	2.49	0.52
1:A:103:LEU:HD21	1:A:247:HIS:C	2.33	0.52
2:C:96:ASP:O	2:C:97:SER:C	2.50	0.52
2:D:109:ILE:O	2:D:109:ILE:CG2	2.58	0.52
2:C:20:THR:CG2	2:C:46:THR:HB	2.39	0.52
1:A:109:ASN:HB3	1:A:466:LYS:CE	2.40	0.52
1:B:109:ASN:CG	1:B:466:LYS:NZ	2.68	0.52
1:A:243:THR:O	1:A:247:HIS:CD2	2.63	0.51
2:C:92:LEU:HD11	2:C:119:LEU:HD21	1.92	0.51
1:A:107:ASP:HB3	1:A:110:PRO:CG	2.40	0.51
1:B:67:HIS:CE1	1:B:249:ASP:OD1	2.63	0.51
1:A:72:ASP:O	1:A:76:THR:CG2	2.30	0.51
1:B:109:ASN:CB	1:B:466:LYS:NZ	2.73	0.51
1:B:422:THR:HB	1:B:463:LEU:HD21	1.92	0.51
2:D:20:THR:CG2	2:D:46:THR:HB	2.38	0.51
1:A:180:PRO:HD2	1:A:181:LYS:H	1.76	0.50
1:B:109:ASN:CB	1:B:466:LYS:CE	2.90	0.50
1:B:109:ASN:CB	1:B:466:LYS:HZ3	2.24	0.50
2:C:57:ARG:HB2	2:C:68:ILE:HD11	1.92	0.50
1:A:77:VAL:O	1:A:80:LEU:O	2.29	0.50
1:A:430:LEU:HA	1:A:433:VAL:HG23	1.93	0.50
1:A:473:VAL:HG12	1:A:488:PHE:CE1	2.47	0.50
2:D:57:ARG:HB2	2:D:68:ILE:HD11	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:31:ARG:HB2	2:C:92:LEU:HD23	1.94	0.49
2:C:37:LEU:HD22	2:C:119:LEU:HD13	1.95	0.49
1:A:109:ASN:HB3	1:A:466:LYS:HZ3	1.76	0.49
2:D:108:ASN:O	2:D:109:ILE:CB	2.60	0.49
2:D:20:THR:HB	2:D:110:TRP:CE2	2.47	0.49
1:B:430:LEU:HA	1:B:433:VAL:HG23	1.92	0.49
2:D:37:LEU:HD22	2:D:119:LEU:HD13	1.94	0.49
1:A:422:THR:CB	1:A:463:LEU:HD21	2.40	0.49
1:B:151:ALA:HB2	1:B:250:LEU:HD22	1.94	0.49
1:B:473:VAL:HG12	1:B:488:PHE:CE1	2.47	0.49
1:A:283:LEU:HD12	1:A:283:LEU:C	2.38	0.48
1:B:113:PRO:O	1:B:145:ARG:NH1	2.45	0.48
1:A:404:GLN:NE2	1:A:428:ARG:HG3	2.28	0.48
1:A:31:LEU:HB2	1:A:39:HIS:HE1	1.78	0.48
1:A:108:ASP:OD1	1:A:197:ARG:HD3	2.13	0.48
1:B:475:LYS:O	1:B:479:GLU:HB2	2.13	0.48
1:A:262:LYS:O	1:A:266:GLU:HG3	2.13	0.47
1:A:107:ASP:CB	1:A:110:PRO:HG3	2.41	0.47
2:C:108:ASN:O	2:C:109:ILE:HB	2.13	0.47
1:A:475:LYS:O	1:A:479:GLU:HB2	2.14	0.47
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.96	0.47
1:B:395:PHE:HE1	1:B:400:GLU:HA	1.80	0.47
1:B:103:LEU:HD21	1:B:247:HIS:O	2.15	0.47
1:B:243:THR:O	1:B:247:HIS:N	2.47	0.47
1:B:437:CYS:HA	1:B:440:HIS:HD2	1.80	0.47
2:C:94:VAL:C	2:C:96:ASP:N	2.71	0.47
1:B:106:LYS:NZ	1:B:147:PRO:O	2.37	0.46
1:A:282:PRO:HD2	1:A:285:GLU:HB2	1.97	0.46
1:B:184:GLU:O	1:B:188:GLU:HG3	2.15	0.46
1:A:156:PHE:HE1	1:A:285:GLU:HG2	1.81	0.46
1:A:75:CYS:SG	1:A:91:CYS:CA	3.04	0.46
1:B:383:GLU:HB3	1:B:384:PRO:HD3	1.96	0.46
1:A:179:LEU:H	1:A:180:PRO:HD3	1.77	0.45
1:B:119:GLU:HG3	1:B:122:VAL:HG23	1.97	0.45
1:A:31:LEU:CD1	1:A:74:LEU:HD22	2.42	0.45
1:B:247:HIS:CD2	1:B:249:ASP:OD2	2.61	0.45
1:B:109:ASN:CG	1:B:466:LYS:HZ3	2.25	0.45
1:A:5:SER:OG	1:A:62:CYS:HB3	2.17	0.45
1:A:113:PRO:O	1:A:145:ARG:NH1	2.50	0.45
1:B:418:VAL:HG12	1:B:419:SER:H	1.82	0.45
1:B:404:GLN:NE2	1:B:428:ARG:HG3	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:474:THR:HG23	1:B:475:LYS:N	2.32	0.45
1:A:41:LYS:HD2	1:B:122:VAL:HG22	1.99	0.44
1:A:474:THR:HG23	1:A:475:LYS:N	2.31	0.44
1:B:117:ARG:NH1	1:B:182:LEU:HB3	2.33	0.44
1:A:204:GLN:NE2	1:A:246:CYS:HB3	2.29	0.44
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.99	0.44
1:A:276:LYS:O	1:A:280:GLU:HG2	2.18	0.44
1:B:149:PHE:CD1	1:B:154:LEU:HB2	2.52	0.44
1:B:208:GLU:O	1:B:209:ARG:C	2.59	0.44
1:A:214:TRP:CD1	1:A:343:VAL:HG11	2.53	0.43
1:A:31:LEU:HB2	1:A:39:HIS:CE1	2.53	0.43
2:D:92:LEU:HD23	2:D:92:LEU:HA	1.48	0.43
1:B:214:TRP:CD1	1:B:343:VAL:HG11	2.54	0.43
1:A:325:VAL:O	1:A:329:MET:HG3	2.19	0.43
2:D:93:THR:CG2	2:D:94:VAL:H	2.29	0.43
1:A:223:PHE:HD1	1:A:272:SER:HB2	1.84	0.42
1:A:109:ASN:C	1:A:110:PRO:O	2.53	0.42
1:A:400:GLU:O	1:A:404:GLN:HG3	2.19	0.42
1:B:394:LEU:O	1:B:397:GLN:HG2	2.18	0.42
1:A:165:PHE:CE1	1:A:178:LEU:HD21	2.53	0.42
1:B:111:ASN:C	1:B:112:LEU:HG	2.44	0.42
1:B:196:GLN:NE2	1:B:246:CYS:SG	2.92	0.42
2:C:110:TRP:CZ3	2:C:112:GLY:N	2.87	0.42
1:B:276:LYS:O	1:B:280:GLU:HG2	2.20	0.42
2:D:93:THR:C	2:D:121:VAL:HG11	2.45	0.42
1:A:281:LYS:CB	1:A:282:PRO:CD	2.98	0.42
1:B:156:PHE:HE1	1:B:285:GLU:HG2	1.84	0.42
1:B:223:PHE:HD1	1:B:272:SER:HB2	1.85	0.42
1:B:325:VAL:HA	2:C:109:ILE:CG1	2.49	0.42
1:A:103:LEU:HD11	1:A:247:HIS:O	2.20	0.41
1:B:115:LEU:HD12	1:B:115:LEU:HA	1.69	0.41
1:A:156:PHE:HD1	1:A:284:LEU:HD12	1.85	0.41
1:B:223:PHE:CD1	1:B:272:SER:HB2	2.55	0.41
2:D:20:THR:CG2	2:D:110:TRP:CZ2	3.03	0.41
1:A:11:PHE:CD1	1:A:54:VAL:HG21	2.55	0.41
1:A:199:LYS:O	1:A:203:LEU:HD12	2.20	0.41
1:A:223:PHE:CD1	1:A:272:SER:HB2	2.56	0.41
1:B:108:ASP:OD1	1:B:148:TYR:HD1	2.04	0.41
1:B:109:ASN:N	1:B:110:PRO:CD	2.84	0.41
1:A:39:HIS:O	1:A:43:VAL:HG23	2.21	0.40
2:D:93:THR:HG22	2:D:95:ALA:N	2.25	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:GLU:C	1:B:61:ASN:OD1	2.65	0.40
1:B:108:ASP:OD1	1:B:148:TYR:CD1	2.74	0.40
1:B:480:SER:HB3	1:B:483:ASN:HB2	2.03	0.40
1:A:378:LYS:HB2	1:A:379:PRO:HD3	2.03	0.40
1:A:420:THR:N	1:A:421:PRO:HD2	2.37	0.40
1:B:167:GLU:CD	1:B:181:LYS:NZ	2.79	0.40
1:A:59:ALA:HB3	1:A:62:CYS:SG	2.62	0.40
1:B:107:ASP:HB3	1:B:110:PRO:HD3	2.02	0.40
2:C:71:SER:H	2:C:74:TYR:HB2	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:GLU:OE2	2:D:31:ARG:NH1[6_654]	1.65	0.55

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	477/585 (82%)	463 (97%)	11 (2%)	3 (1%)	21	53
1	B	477/585 (82%)	463 (97%)	10 (2%)	4 (1%)	16	46
2	C	102/128 (80%)	95 (93%)	6 (6%)	1 (1%)	12	41
2	D	102/128 (80%)	95 (93%)	6 (6%)	1 (1%)	12	41
All	All	1158/1426 (81%)	1116 (96%)	33 (3%)	9 (1%)	16	46

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	419	SER

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Mol	Chain	Res	Type
2	D	109	ILE
1	A	479	GLU
1	B	479	GLU
2	C	109	ILE
1	A	300	ALA
1	B	300	ALA
1	A	109	ASN
1	B	109	ASN

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	425/511 (83%)	412 (97%)	13 (3%)	35	65
1	B	425/511 (83%)	414 (97%)	11 (3%)	40	69
2	C	91/110 (83%)	87 (96%)	4 (4%)	25	56
2	D	91/110 (83%)	88 (97%)	3 (3%)	33	64
All	All	1032/1242 (83%)	1001 (97%)	31 (3%)	36	66

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	LEU
1	A	142	ILE
1	A	155	LEU
1	A	245	CYS
1	A	281	LYS
1	A	284	LEU
1	A	311	GLU
1	A	340	ASP
1	A	357	LEU
1	A	400	GLU
1	A	423	LEU
1	A	425	GLU
1	A	492	GLU

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Mol	Chain	Res	Type
1	B	7	VAL
1	B	60	GLU
1	B	106	LYS
1	B	112	LEU
1	B	142	ILE
1	B	233	LYS
1	B	245	CYS
1	B	297	GLU
1	B	357	LEU
1	B	392	CYS
1	B	426	VAL
2	C	78	VAL
2	C	97	SER
2	C	110	TRP
2	C	123	HIS
2	D	25	THR
2	D	78	VAL
2	D	108	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	130	ASN
1	A	204	GLN
1	A	247	HIS
1	A	288	HIS
1	A	367	HIS
1	A	404	GLN
1	A	483	ASN
1	B	39	HIS
1	B	109	ASN
1	B	130	ASN
1	B	196	GLN
1	B	242	HIS
1	B	247	HIS
1	B	338	HIS
1	B	367	HIS
1	B	404	GLN
1	B	417	GLN
1	B	440	HIS
1	B	483	ASN
2	C	64	ASN

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Mol	Chain	Res	Type
2	D	64	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	483/585 (82%)	0.56	26 (5%)	31 15	44, 83, 149, 175	0
1	B	481/585 (82%)	0.42	26 (5%)	31 15	41, 80, 151, 179	0
2	C	104/128 (81%)	0.15	3 (2%)	53 30	39, 64, 90, 101	0
2	D	104/128 (81%)	0.03	1 (0%)	79 58	46, 65, 95, 102	0
All	All	1172/1426 (82%)	0.42	56 (4%)	35 18	39, 78, 147, 179	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	427	SER	5.8
1	B	424	VAL	5.7
2	C	109	ILE	4.8
1	B	428	ARG	4.6
1	A	426	VAL	4.4
1	B	125	THR	4.3
1	B	271	ILE	4.2
1	B	439	LYS	3.8
1	B	7	VAL	3.7
1	A	430	LEU	3.7
1	A	407	LEU	3.6
1	B	129	ASP	3.5
1	A	453	LEU	3.5
1	B	183	ASP	3.4
1	A	423	LEU	3.4
1	B	88	ALA	3.1
1	B	206	PHE	3.1
1	A	415	VAL	3.0
1	B	463	LEU	3.0
1	B	270	SER	3.0
1	B	420	THR	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	206	PHE	2.9
1	A	69	LEU	2.8
1	B	91	CYS	2.8
1	B	453	LEU	2.7
2	C	92	LEU	2.7
1	A	467	THR	2.6
1	A	312	SER	2.6
1	A	271	ILE	2.6
1	B	488	PHE	2.6
1	A	237	ASP	2.6
1	A	212	LYS	2.5
1	A	377	PHE	2.5
1	A	4	LYS	2.4
1	A	454	SER	2.4
1	A	213	ALA	2.3
1	B	457	LEU	2.3
1	B	390	GLN	2.3
1	B	480	SER	2.2
1	A	160	ARG	2.2
1	A	478	THR	2.2
1	A	30	TYR	2.2
1	A	490	ALA	2.2
1	B	208	GLU	2.1
1	B	421	PRO	2.1
1	B	412	THR	2.1
1	B	207	GLY	2.1
1	B	305	LEU	2.1
1	A	215	ALA	2.1
1	B	473	VAL	2.1
1	B	134	PHE	2.1
2	C	73	ARG	2.1
2	D	91	ASP	2.0
1	A	202	SER	2.0
1	A	91	CYS	2.0
1	A	473	VAL	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 [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.