



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:54 PM UTC

PDB ID : 9G13 / pdb_00009g13
Title : VHH H3-2 in complex with Tau C-terminal peptide
Authors : Dupre, E.; Landrieu, I.; Danis, C.; Hanouille, X.; Morteletcque, J.
Deposited on : 2024-07-09
Resolution : 1.80 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

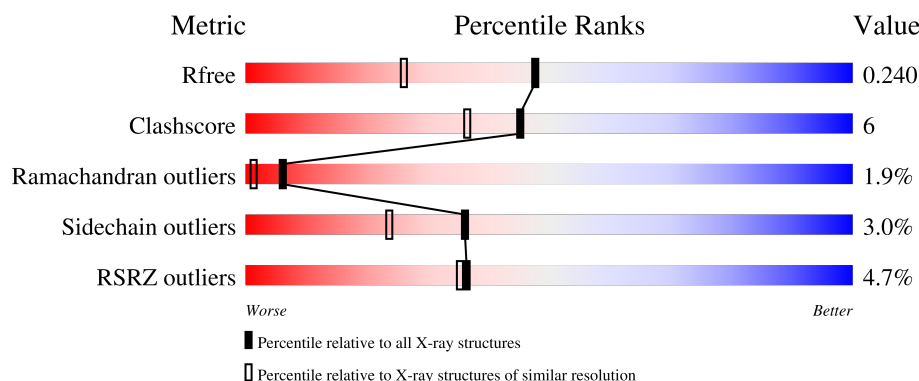
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	132	5% 73% 20% • 5%
1	C	132	5% 74% 17% • 5%
1	E	132	4% 88% 8% •
1	G	132	6% 74% 19% • 5%
2	B	13	54% 38% 8%

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Mol	Chain	Length	Quality of chain
2	D	13	8% 54% 38% 8%
2	F	13	62% 23% 8% 8%
2	H	13	54% 31% 15%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8787 atoms, of which 4072 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 VHH H3-2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	126	Total 1840	C 594	H 893	N 162	O 186	S 5	40	0	0
1	E	127	Total 1885	C 609	H 914	N 163	O 194	S 5	38	0	0
1	C	125	Total 1868	C 602	H 908	N 165	O 188	S 5	39	2	0
1	G	126	Total 1844	C 596	H 896	N 162	O 185	S 5	40	0	0

- Molecule 2 is a protein called Isoform Tau-F of Microtubule-associated protein tau.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	12	Total	C	H	N	O	4	0	0
			226	69	119	20	18			
2	F	12	Total	C	H	N	O	4	0	0
			226	69	119	20	18			
2	D	12	Total	C	H	N	O	4	0	0
			226	69	119	20	18			
2	H	11	Total	C	H	N	O	4	0	0
			202	63	104	18	17			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	107	Total	O	0	0
			107	107		
3	B	17	Total	O	0	0
			17	17		
3	E	103	Total	O	0	0
			103	103		
3	F	19	Total	O	0	0
			19	19		

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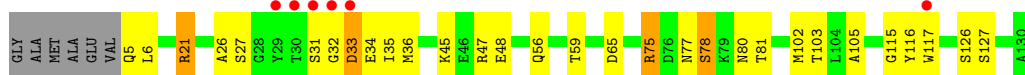
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	105	Total 105	O 105	0	0
3	D	16	Total 16	O 16	0	0
3	G	86	Total 86	O 86	0	0
3	H	17	Total 17	O 17	0	0

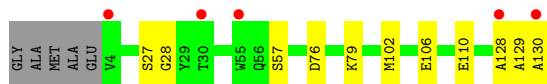
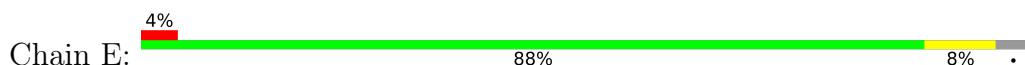
3 Residue-property plots [i](#)

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: VHH H3-2



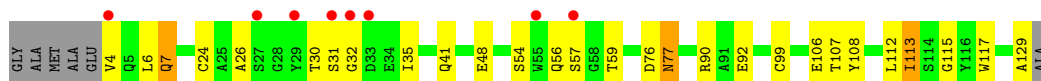
- Molecule 1: VHH H3-2



- Molecule 1: VHH H3-2



- Molecule 1: VHH H3-2



- Molecule 2: Isoform Tau-F of Microtubule-associated protein tau



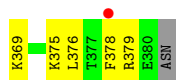
- Molecule 2: Isoform Tau-F of Microtubule-associated protein tau

Chain F:  62% 23% 8% 8%



- Molecule 2: Isoform Tau-F of Microtubule-associated protein tau

Chain D:  8% 54% 38% 8%



- Molecule 2: Isoform Tau-F of Microtubule-associated protein tau

Chain H:  54% 31% 15%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.39Å 76.35Å 95.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.55 – 1.80 45.55 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.55-1.80) 99.9 (45.55-1.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, R_{free}	0.172 , 0.231 (Not available) , 0.240	Depositor DCC
R_{free} test set	2348 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8787	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.11% 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

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	1.11	1/969 (0.1%)	1.50	10/1315 (0.8%)
1	C	0.97	1/988 (0.1%)	1.49	8/1340 (0.6%)
1	E	0.98	0/994	1.47	4/1349 (0.3%)
1	G	0.98	0/970	1.47	8/1317 (0.6%)
2	B	1.00	0/108	1.80	2/141 (1.4%)
2	D	0.89	0/108	1.61	1/141 (0.7%)
2	F	1.17	1/108 (0.9%)	1.40	1/141 (0.7%)
2	H	1.25	0/99	1.96	1/130 (0.8%)
All	All	1.02	3/4344 (0.1%)	1.50	35/5874 (0.6%)

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	0	3
1	C	0	1
1	E	0	2
All	All	0	6

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	374	HIS	CE1-NE2	-5.60	1.26	1.32
1	A	127	SER	CA-CB	-5.43	1.44	1.53
1	C	101	PRO	CA-CB	5.27	1.60	1.53

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	THR	CA-CB-OG1	-9.23	95.76	109.60
1	E	130	ALA	CA-C-O	-8.88	105.70	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	376	LEU	N-CA-CB	-7.95	97.40	110.52
1	C	71	PHE	CA-CB-CG	7.83	121.63	113.80
1	C	116	TYR	CA-C-N	-7.55	109.28	122.64
1	C	116	TYR	C-N-CA	-7.55	109.28	122.64
1	A	21	ARG	NE-CZ-NH1	7.41	128.91	121.50
1	A	56	GLN	N-CA-CB	-7.00	99.61	110.39
1	E	106	GLU	CB-CG-CD	6.91	124.34	112.60
1	G	59	THR	CA-CB-OG1	-6.46	99.91	109.60
1	A	81	THR	CA-CB-OG1	-6.34	100.09	109.60
1	C	108	TYR	N-CA-CB	6.20	120.29	110.55
1	C	62	TYR	O-C-N	6.17	130.83	123.24
1	E	106	GLU	OE1-CD-OE2	-5.95	108.62	122.90
1	G	41	GLN	OE1-CD-NE2	5.69	128.29	122.60
1	A	45	LYS	CG-CD-CE	5.61	124.19	111.30
1	C	116	TYR	N-CA-CB	5.58	118.49	110.06
1	G	112	LEU	N-CA-CB	-5.58	101.71	110.69
1	C	92	GLU	CB-CG-CD	5.57	122.06	112.60
1	G	113	ILE	CA-C-O	-5.55	114.56	120.39
1	G	115	GLY	CA-C-O	-5.53	118.15	122.52
2	F	376	LEU	N-CA-CB	-5.52	101.40	110.68
1	G	48	GLU	CB-CG-CD	-5.46	103.31	112.60
1	G	129	ALA	CA-C-O	-5.43	111.57	120.80
1	E	110	GLU	CB-CG-CD	5.39	121.76	112.60
2	B	377	THR	CA-C-O	5.34	125.90	120.24
1	A	21	ARG	NE-CZ-NH2	-5.32	114.41	119.20
1	A	77	ASN	CA-CB-CG	-5.29	107.31	112.60
2	H	377	THR	CA-CB-OG1	-5.23	101.75	109.60
1	G	113	ILE	O-C-N	5.21	128.89	123.26
1	A	103	THR	OG1-CB-CG2	-5.16	98.97	109.30
1	A	36	MET	CA-C-N	-5.11	116.07	121.48
1	A	36	MET	C-N-CA	-5.11	116.07	121.48
1	C	100	ALA	O-C-N	-5.03	117.98	121.65
2	B	373	THR	O-C-N	-5.02	117.44	123.31

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	21	ARG	Sidechain
1	A	47	ARG	Sidechain
1	A	75	ARG	Sidechain
1	C	75	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	E	128	ALA	Peptide
1	E	57	SER	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	947	893	881	12	1
1	C	960	908	901	12	0
1	E	971	914	911	4	0
1	G	948	896	882	14	1
2	B	107	119	115	3	0
2	D	107	119	115	4	0
2	F	107	119	115	3	0
2	H	98	104	102	1	0
3	A	107	0	0	8	0
3	B	17	0	0	0	1
3	C	105	0	0	2	0
3	D	16	0	0	2	0
3	E	103	0	0	1	0
3	F	19	0	0	3	0
3	G	86	0	0	1	0
3	H	17	0	0	1	1
All	All	4715	4072	4022	48	2

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 6.

All (48) 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:G:35:ILE:HG12	1:G:54:SER:HB3	1.41	0.99
1:E:79:LYS:O	3:E:201:HOH:O	2.04	0.75
2:F:370:LYS:HD3	3:F:417:HOH:O	1.86	0.74
1:A:65:ASP:CG	3:A:232:HOH:O	2.32	0.72
1:A:6:LEU:HD23	1:A:26:ALA:HA	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:GLU:OE2	3:A:201:HOH:O	2.09	0.69
1:A:48:GLU:OE2	3:A:202:HOH:O	2.11	0.68
1:A:78:SER:HB3	3:A:204:HOH:O	1.94	0.66
1:A:5:GLN:HA	1:A:5:GLN:OE1	1.94	0.65
1:A:78:SER:CB	3:A:204:HOH:O	2.48	0.62
1:C:98:TYR:OH	3:C:201:HOH:O	2.13	0.59
2:D:375:LYS:NZ	3:D:401:HOH:O	2.35	0.59
3:A:270:HOH:O	1:C:103:THR:HG23	2.02	0.59
1:G:56:GLN:O	1:G:57:SER:OG	2.21	0.59
1:C:21[A]:ARG:HH11	1:C:85:GLN:HE21	1.54	0.56
1:C:35:ILE:HG12	1:C:54:SER:HB3	1.88	0.55
2:D:369:LYS:HB3	3:D:414:HOH:O	2.06	0.55
1:C:117:TRP:CZ3	2:D:378:PHE:CE2	2.95	0.54
1:A:33:ASP:O	1:A:35:ILE:N	2.38	0.54
2:F:369:LYS:N	3:F:401:HOH:O	2.41	0.53
1:C:6:LEU:HB3	3:C:277:HOH:O	2.08	0.52
1:A:105:ALA:HB2	1:C:112:LEU:HD22	1.94	0.50
1:E:102:MET:HE2	1:G:113:ILE:HD11	1.92	0.50
1:A:75:ARG:NH2	1:A:80:ASN:OD1	2.44	0.50
3:A:236:HOH:O	2:B:370:LYS:HG2	2.12	0.49
1:G:4:VAL:HG13	3:G:286:HOH:O	2.12	0.49
1:G:24:CYS:CB	1:G:99:CYS:SG	3.01	0.48
1:G:106:GLU:HG2	1:G:108:TYR:CE1	2.47	0.48
1:G:77:ASN:OD1	1:G:77:ASN:N	2.47	0.47
2:B:378:PHE:CZ	1:C:47:ARG:HD2	2.50	0.47
1:G:106:GLU:HG3	1:G:107:THR:N	2.31	0.46
1:E:76:ASP:OD2	1:E:79:LYS:HG3	2.15	0.46
1:E:27:SER:OG	1:E:28:GLY:N	2.49	0.46
1:G:106:GLU:HG2	1:G:108:TYR:CZ	2.51	0.46
1:G:30:THR:O	1:G:32:GLY:N	2.48	0.46
1:C:54:SER:O	1:C:59:THR:HA	2.16	0.45
1:G:90:ARG:HG3	1:G:92:GLU:HG2	1.97	0.45
2:F:370:LYS:NZ	3:F:402:HOH:O	2.44	0.44
1:C:75:ARG:HH22	1:C:77:ASN:ND2	2.15	0.44
1:G:6:LEU:HD23	1:G:26:ALA:HA	2.00	0.43
2:D:379:ARG:HD3	2:D:379:ARG:HA	1.81	0.43
1:A:102:MET:HG2	1:C:113:ILE:HG12	1.99	0.42
1:G:7:GLN:HE21	1:G:7:GLN:HA	1.83	0.42
2:H:375:LYS:HE3	3:H:414:HOH:O	2.19	0.42
1:G:76:ASP:OD1	1:G:76:ASP:C	2.60	0.42
1:A:115:GLY:O	1:A:117:TRP:CD1	2.72	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:106:GLU:HG2	1:C:108:TYR:CE1	2.55	0.41
3:A:236:HOH:O	2:B:370:LYS:CG	2.69	0.40

All (2) 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:116:TYR:HH	1:G:117:TRP:H[1_455]	1.23	0.37
3:B:408:HOH:O	3:H:411:HOH:O[3_544]	1.92	0.28

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	124/132 (94%)	113 (91%)	7 (6%)	4 (3%)	3	0
1	C	125/132 (95%)	117 (94%)	4 (3%)	4 (3%)	3	0
1	E	125/132 (95%)	119 (95%)	5 (4%)	1 (1%)	16	6
1	G	124/132 (94%)	118 (95%)	5 (4%)	1 (1%)	16	6
2	B	10/13 (77%)	10 (100%)	0	0	100	100
2	D	10/13 (77%)	8 (80%)	2 (20%)	0	100	100
2	F	10/13 (77%)	10 (100%)	0	0	100	100
2	H	9/13 (69%)	9 (100%)	0	0	100	100
All	All	537/580 (93%)	504 (94%)	23 (4%)	10 (2%)	6	1

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	31	SER
1	A	32	GLY

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Mol	Chain	Res	Type
1	A	33	ASP
1	C	29	TYR
1	A	31	SER
1	C	30	THR
1	C	33	ASP
1	A	34	GLU
1	E	129	ALA
1	C	56	GLN

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	94/102 (92%)	91 (97%)	3 (3%)	34	22
1	C	97/102 (95%)	93 (96%)	4 (4%)	27	15
1	E	100/102 (98%)	100 (100%)	0	100	100
1	G	94/102 (92%)	92 (98%)	2 (2%)	47	36
2	B	12/13 (92%)	11 (92%)	1 (8%)	10	3
2	D	12/13 (92%)	12 (100%)	0	100	100
2	F	12/13 (92%)	11 (92%)	1 (8%)	10	3
2	H	11/13 (85%)	9 (82%)	2 (18%)	2	0
All	All	432/460 (94%)	419 (97%)	13 (3%)	36	24

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	SER
1	A	78	SER
1	A	126	SER
2	B	375	LYS
2	F	369	LYS
1	C	56	GLN
1	C	60	SER

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Mol	Chain	Res	Type
1	C	103	THR
1	C	112	LEU
1	G	7	GLN
1	G	77	ASN
2	H	370	LYS
2	H	380	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	56	GLN
1	C	77	ASN
1	G	5	GLN
1	G	7	GLN
1	G	15	GLN
1	G	122	GLN

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 ⓘ

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	126/132 (95%)	-0.06	6 (4%) 35 34	19, 30, 82, 108	0
1	C	125/132 (94%)	-0.02	6 (4%) 35 34	15, 29, 85, 110	2 (1%)
1	E	127/132 (96%)	-0.10	5 (3%) 43 43	19, 31, 80, 95	0
1	G	126/132 (95%)	0.04	8 (6%) 26 24	20, 33, 87, 129	0
2	B	12/13 (92%)	0.33	0 100 100	23, 39, 84, 98	0
2	D	12/13 (92%)	0.37	1 (8%) 17 15	23, 42, 93, 106	0
2	F	12/13 (92%)	0.07	0 100 100	21, 38, 66, 67	0
2	H	11/13 (84%)	0.14	0 100 100	20, 32, 77, 79	0
All	All	551/580 (95%)	-0.01	26 (4%) 36 35	15, 31, 87, 129	2 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	31	SER	4.3
1	A	117	TRP	4.2
1	C	29	TYR	3.8
1	C	30	THR	3.8
1	G	31	SER	3.6
1	A	30	THR	3.5
1	E	55	TRP	3.4
1	A	33	ASP	3.1
1	E	4	VAL	3.1
1	G	29	TYR	3.0
1	G	33	ASP	3.0
1	C	117	TRP	3.0
1	C	57	SER	2.9
1	G	27	SER	2.9
1	A	32	GLY	2.9
1	E	128	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	G	32	GLY	2.7
1	E	130	ALA	2.6
1	G	4	VAL	2.4
1	A	31	SER	2.3
1	G	55	TRP	2.3
1	G	57	SER	2.3
1	A	29	TYR	2.2
1	E	30	THR	2.2
1	C	129	ALA	2.1
2	D	378	PHE	2.0

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 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.