



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:48 AM UTC

PDB ID : 5MP2 / pdb_00005mp2
Title : XcpQN012 in complex with VHH04
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Deposited on : 2016-12-15
Resolution : 2.90 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
Xtrriage (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

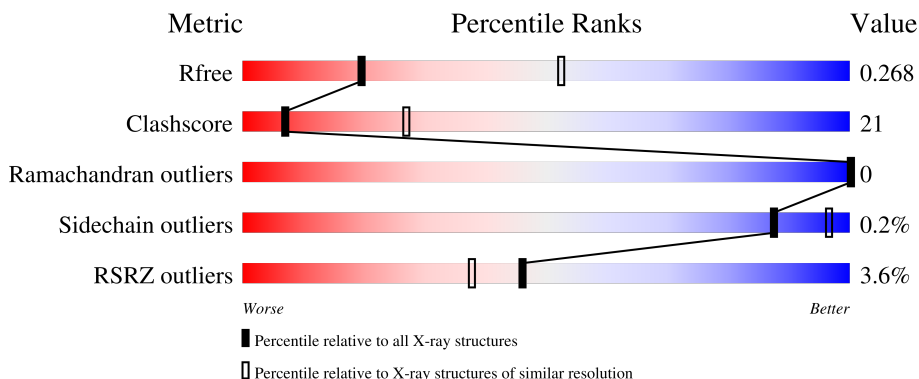
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	240	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 74%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: grey;"></div> </div>
1	B	240	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 63%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 23%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: grey;"></div> </div>
2	C	153	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 58%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 21%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 20%; height: 10px; background-color: grey;"></div> </div>
2	D	153	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 59%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 22%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 18%; height: 10px; background-color: grey;"></div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5028 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 Type II secretion system protein D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	214	Total 1539	C 973	N 270	O 293	S 3	0	0	0
1	B	212	Total 1505	C 958	N 252	O 292	S 3	0	0	0

- Molecule 2 is a protein called Camelid nanobody VHH04.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	123	Total 924	C 580	N 157	O 182	S 5	0	0	0
2	D	125	Total 936	C 587	N 160	O 184	S 5	0	0	0

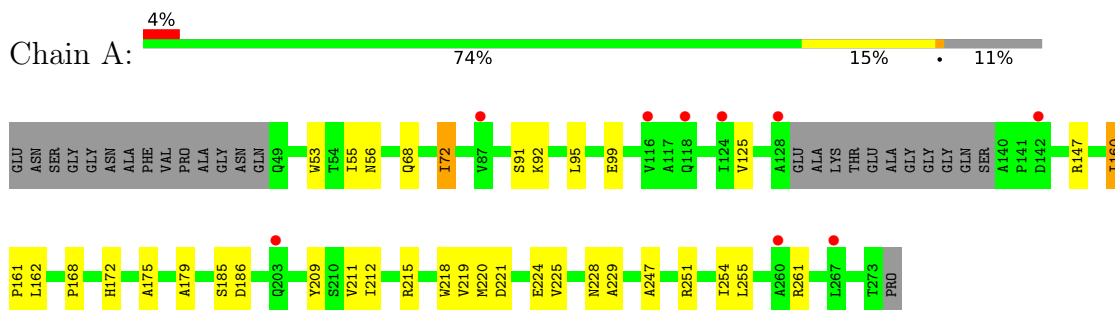
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	36	Total 36	O 36	0	0
3	B	42	Total 42	O 42	0	0
3	C	16	Total 16	O 16	0	0
3	D	30	Total 30	O 30	0	0

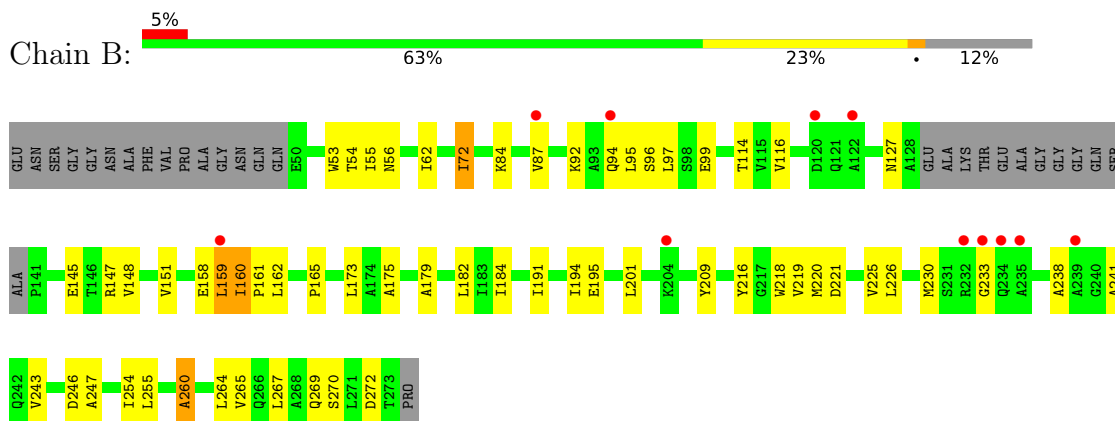
3 Residue-property plots [i](#)

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.

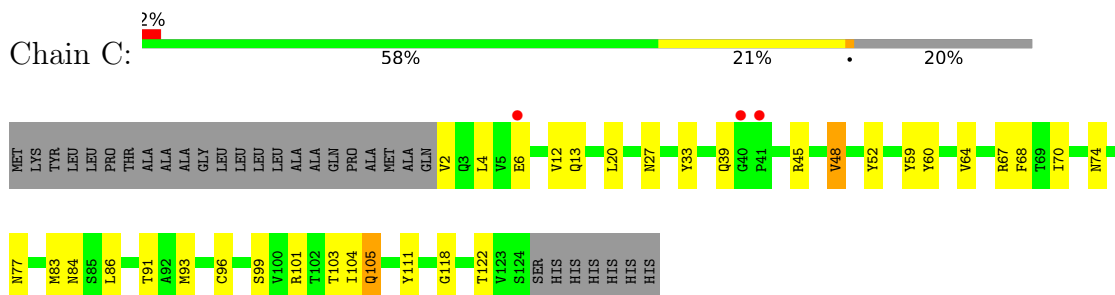
- Molecule 1: Type II secretion system protein D



- Molecule 1: Type II secretion system protein D

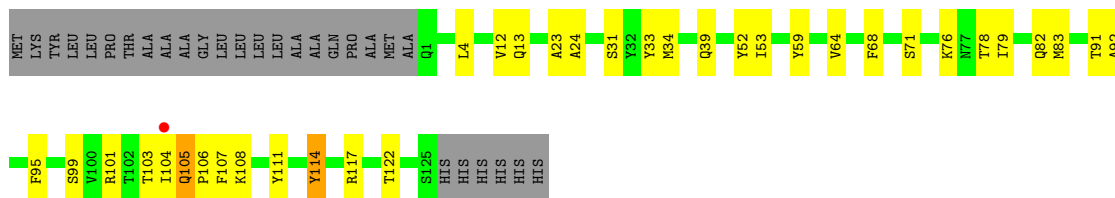


- Molecule 2: Camelid nanobody VHH04



- Molecule 2: Camelid nanobody VHH04





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	40.06Å 63.79Å 76.05Å 104.37° 100.61° 108.04°	Depositor
Resolution (Å)	38.75 – 2.90 38.75 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.2 (38.75-2.90) 95.2 (38.75-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 2.90Å)	Xtrriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.213 , 0.256 0.227 , 0.268	Depositor DCC
R_{free} test set	647 reflections (4.39%)	wwPDB-VP
Wilson B-factor (Å ²)	60.9	Xtrriage
Anisotropy	0.723	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 89.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5028	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.99% 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	0.95	0/1563	1.39	5/2142 (0.2%)
1	B	0.92	0/1529	1.36	6/2100 (0.3%)
2	C	0.94	0/945	1.21	2/1282 (0.2%)
2	D	0.96	0/957	1.27	5/1298 (0.4%)
All	All	0.94	0/4994	1.33	18/6822 (0.3%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	105	GLN	CA-C-N	8.13	127.79	119.82
2	D	105	GLN	C-N-CA	8.13	127.79	119.82
2	C	105	GLN	N-CA-C	-6.87	96.17	108.47
1	A	229	ALA	CA-C-N	6.42	128.79	120.44
1	A	229	ALA	C-N-CA	6.42	128.79	120.44
1	B	260	ALA	CA-C-N	5.86	128.14	120.28
1	B	260	ALA	C-N-CA	5.86	128.14	120.28
2	D	52	TYR	CA-C-N	5.82	128.39	120.77
2	D	52	TYR	C-N-CA	5.82	128.39	120.77
1	A	72	ILE	CA-C-N	5.80	128.05	120.28
1	A	72	ILE	C-N-CA	5.80	128.05	120.28
2	D	105	GLN	N-CA-C	-5.59	97.44	109.01
2	C	48	VAL	CB-CA-C	-5.56	105.46	110.91
1	B	160	ILE	CB-CA-C	-5.45	108.57	114.35
1	B	72	ILE	CA-C-N	5.31	127.39	120.28
1	B	72	ILE	C-N-CA	5.31	127.39	120.28
1	A	160	ILE	CB-CA-C	-5.10	108.94	114.35
1	B	159	LEU	N-CA-C	5.00	117.62	111.82

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	1539	0	1487	32	1
1	B	1505	0	1442	62	1
2	C	924	0	860	67	0
2	D	936	0	874	62	0
3	A	36	0	0	1	0
3	B	42	0	0	0	0
3	C	16	0	0	0	0
3	D	30	0	0	0	0
All	All	5028	0	4663	201	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 21.

All (201) 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)
2:C:64:VAL:HB	2:C:68:PHE:CD2	1.71	1.26
2:D:104:ILE:HG23	2:D:111:TYR:CG	1.76	1.20
2:C:6:GLU:OE2	2:C:96:CYS:CB	1.96	1.13
2:C:104:ILE:HG23	2:C:111:TYR:CG	1.84	1.12
2:C:6:GLU:OE2	2:C:96:CYS:HB3	1.48	1.12
1:A:209:TYR:HB3	1:A:255:LEU:CD2	1.79	1.11
2:D:104:ILE:HG22	2:D:105:GLN:H	1.11	1.10
2:C:104:ILE:CG2	2:C:111:TYR:HB3	1.79	1.10
2:C:101:ARG:CB	2:C:104:ILE:HD11	1.81	1.10
1:B:160:ILE:HD11	1:B:175:ALA:HB2	1.29	1.09
2:C:101:ARG:HB2	2:C:104:ILE:CD1	1.82	1.09
2:C:68:PHE:CD1	2:C:83:MET:HG2	1.90	1.05
1:A:209:TYR:HB3	1:A:255:LEU:HD23	1.32	1.05
2:D:104:ILE:HG22	2:D:105:GLN:N	1.62	1.05
1:B:94:GLN:NE2	2:D:106:PRO:HG2	1.74	1.01
2:C:103:THR:O	2:C:104:ILE:HG13	1.59	1.01
2:C:104:ILE:CG2	2:C:111:TYR:CB	2.38	1.01
1:A:160:ILE:HD11	1:A:175:ALA:HB2	1.42	1.00
2:D:104:ILE:CG2	2:D:111:TYR:CB	2.40	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:104:ILE:HG21	2:C:111:TYR:HB3	1.40	0.99
2:D:104:ILE:CG2	2:D:105:GLN:H	1.74	0.99
2:D:104:ILE:HG21	2:D:111:TYR:HB3	1.44	0.98
2:C:68:PHE:CE1	2:C:83:MET:HG2	1.98	0.98
2:C:104:ILE:HG23	2:C:111:TYR:CB	1.90	0.98
2:D:104:ILE:CG2	2:D:111:TYR:HB3	1.92	0.98
1:B:94:GLN:NE2	2:D:106:PRO:CG	2.27	0.97
1:B:94:GLN:OE1	2:D:107:PHE:HE2	1.49	0.96
1:B:55:ILE:HA	2:D:103:THR:HG22	1.49	0.95
1:B:94:GLN:OE1	2:D:107:PHE:CE2	2.19	0.94
2:C:64:VAL:HB	2:C:68:PHE:HD2	1.17	0.93
2:D:104:ILE:HG21	2:D:111:TYR:CB	1.96	0.92
1:B:151:VAL:HG11	1:B:159:LEU:HD11	1.53	0.90
1:B:94:GLN:HE21	2:D:106:PRO:CG	1.86	0.89
2:D:34:MET:HB2	2:D:79:ILE:CD1	2.02	0.89
1:B:267:LEU:O	1:B:270:SER:OG	1.90	0.88
1:B:94:GLN:HE21	2:D:106:PRO:HG3	1.38	0.87
1:A:209:TYR:CB	1:A:255:LEU:HD23	2.04	0.86
2:D:33:TYR:HB2	2:D:99:SER:OG	1.73	0.86
1:B:55:ILE:HA	2:D:103:THR:CG2	2.05	0.86
2:D:104:ILE:CG2	2:D:111:TYR:CG	2.58	0.85
2:D:104:ILE:HG23	2:D:111:TYR:CB	2.04	0.85
2:D:101:ARG:HB2	2:D:104:ILE:HD11	1.55	0.85
2:C:104:ILE:HG23	2:C:111:TYR:HB3	1.53	0.84
2:D:104:ILE:HG23	2:D:111:TYR:CD1	2.14	0.82
2:C:101:ARG:HB2	2:C:104:ILE:HD11	0.89	0.82
2:D:34:MET:CB	2:D:79:ILE:HD12	2.11	0.81
1:B:209:TYR:HB3	1:B:255:LEU:HD23	1.61	0.81
2:D:31:SER:HB3	2:D:114:TYR:OH	1.81	0.81
2:C:68:PHE:CE1	2:C:83:MET:CG	2.63	0.81
2:C:68:PHE:HD1	2:C:83:MET:HG2	1.45	0.81
2:C:103:THR:C	2:C:104:ILE:HG13	2.06	0.79
2:D:31:SER:CB	2:D:114:TYR:OH	2.29	0.79
1:A:209:TYR:HB3	1:A:255:LEU:HD21	1.64	0.79
1:B:151:VAL:HG11	1:B:159:LEU:CD1	2.12	0.78
2:C:6:GLU:OE2	2:C:96:CYS:HB2	1.81	0.78
1:A:212:ILE:HD11	1:A:254:ILE:HD12	1.66	0.78
2:C:64:VAL:CB	2:C:68:PHE:CD2	2.62	0.77
2:C:68:PHE:HE1	2:C:83:MET:CG	1.98	0.76
2:C:48:VAL:CG1	2:C:64:VAL:HG21	2.16	0.75
2:D:34:MET:CB	2:D:79:ILE:CD1	2.64	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:TYR:HB3	1:B:255:LEU:CD2	2.16	0.74
2:C:48:VAL:CG1	2:C:64:VAL:HG11	2.16	0.74
2:C:104:ILE:HG21	2:C:111:TYR:CB	2.10	0.74
2:C:104:ILE:HG23	2:C:111:TYR:CD1	2.24	0.72
1:A:160:ILE:HB	1:A:161:PRO:HD3	1.72	0.71
2:D:34:MET:HB3	2:D:79:ILE:HD12	1.73	0.71
2:C:99:SER:OG	2:C:104:ILE:CD1	2.38	0.70
2:D:53:ILE:HD11	2:D:79:ILE:HD11	1.75	0.69
2:D:39:GLN:O	2:D:92:ALA:HB1	1.93	0.69
1:B:94:GLN:NE2	2:D:107:PHE:CZ	2.61	0.68
1:B:54:THR:CG2	1:B:94:GLN:HG2	2.24	0.67
2:C:48:VAL:HG12	2:C:64:VAL:HG21	1.76	0.66
1:B:96:SER:OG	1:B:99:GLU:CG	2.44	0.66
2:C:103:THR:O	2:C:104:ILE:CG1	2.38	0.66
2:C:104:ILE:HG22	2:C:105:GLN:N	2.10	0.66
1:A:209:TYR:CB	1:A:255:LEU:CD2	2.65	0.65
1:B:94:GLN:NE2	2:D:107:PHE:CE2	2.65	0.65
2:C:48:VAL:HG13	2:C:64:VAL:HG11	1.77	0.64
1:B:96:SER:OG	1:B:99:GLU:HG3	1.97	0.64
1:B:94:GLN:HE22	2:D:106:PRO:HG2	1.63	0.63
1:B:94:GLN:CD	2:D:107:PHE:CE2	2.77	0.63
1:B:94:GLN:NE2	2:D:106:PRO:HG3	2.03	0.63
1:A:212:ILE:CD1	1:A:254:ILE:HD12	2.29	0.62
2:C:60:TYR:CE1	2:C:70:ILE:HG22	2.34	0.62
1:B:230:MET:HE1	1:B:241:ALA:HB3	1.81	0.62
2:D:104:ILE:HD13	2:D:111:TYR:HB3	1.80	0.62
1:B:56:ASN:H	2:D:103:THR:HG22	1.65	0.61
1:B:96:SER:OG	1:B:99:GLU:CD	2.44	0.61
1:B:216:TYR:HD2	1:B:272:ASP:OD1	1.85	0.60
2:C:48:VAL:HG11	2:C:64:VAL:HG11	1.81	0.60
2:D:33:TYR:HB2	2:D:99:SER:HG	1.67	0.60
2:D:53:ILE:CD1	2:D:79:ILE:HD11	2.32	0.59
1:B:238:ALA:HB1	1:B:260:ALA:HA	1.83	0.59
2:C:64:VAL:CB	2:C:68:PHE:HD2	2.03	0.58
2:D:105:GLN:NE2	2:D:108:LYS:HG3	2.19	0.58
1:B:221:ASP:O	1:B:225:VAL:HG23	2.03	0.58
1:A:221:ASP:O	1:A:225:VAL:HG23	2.03	0.57
2:C:48:VAL:HG13	2:C:64:VAL:CG1	2.35	0.56
2:D:114:TYR:N	2:D:114:TYR:CD2	2.73	0.56
2:C:68:PHE:CE1	2:C:83:MET:CB	2.88	0.55
2:C:104:ILE:CG2	2:C:105:GLN:N	2.69	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:104:ILE:HG22	2:C:105:GLN:O	2.06	0.55
1:A:219:VAL:HG11	1:A:247:ALA:HA	1.88	0.55
1:A:215:ARG:HD2	3:A:311:HOH:O	2.06	0.54
1:B:92:LYS:HE3	1:B:147:ARG:HD3	1.89	0.54
1:B:162:LEU:HD13	1:B:255:LEU:HD11	1.88	0.54
1:A:162:LEU:HD21	1:A:209:TYR:CD1	2.43	0.54
2:D:104:ILE:HG21	2:D:111:TYR:HB2	1.84	0.54
1:B:218:TRP:NE1	1:B:220:MET:HB2	2.23	0.54
2:D:68:PHE:CD1	2:D:83:MET:HA	2.42	0.53
1:B:62:ILE:HG12	1:B:87:VAL:CG2	2.39	0.53
2:D:104:ILE:HG22	2:D:105:GLN:C	2.33	0.53
1:A:92:LYS:HE3	1:A:147:ARG:HD3	1.91	0.53
2:C:6:GLU:OE2	2:C:96:CYS:N	2.41	0.53
1:A:212:ILE:CD1	1:A:254:ILE:CD1	2.87	0.53
1:A:125:VAL:HG11	1:B:116:VAL:HG13	1.91	0.53
2:D:4:LEU:HD23	2:D:24:ALA:HA	1.91	0.53
1:A:95:LEU:HD22	1:A:99:GLU:HB3	1.91	0.52
1:B:95:LEU:HD21	1:B:148:VAL:HG11	1.91	0.52
2:C:104:ILE:CG2	2:C:111:TYR:CG	2.72	0.51
1:A:160:ILE:HD11	1:A:175:ALA:CB	2.28	0.51
2:C:68:PHE:CD1	2:C:83:MET:HA	2.45	0.51
1:B:216:TYR:CD2	1:B:272:ASP:OD1	2.62	0.51
2:C:83:MET:HB3	2:C:86:LEU:HD21	1.92	0.51
2:C:68:PHE:CD1	2:C:83:MET:CG	2.80	0.51
1:B:54:THR:HG23	1:B:94:GLN:HG2	1.93	0.50
1:A:254:ILE:HG22	1:A:261:ARG:HG2	1.93	0.50
2:D:71:SER:O	2:D:79:ILE:HG23	2.11	0.50
2:C:39:GLN:HG2	2:C:45:ARG:HG2	1.93	0.50
2:C:91:THR:HG23	2:C:122:THR:HA	1.94	0.50
1:B:147:ARG:HD2	1:B:195:GLU:OE2	2.11	0.50
1:B:219:VAL:HG11	1:B:247:ALA:HA	1.93	0.50
1:A:218:TRP:NE1	1:A:220:MET:HB2	2.26	0.49
1:B:173:LEU:CD1	1:B:184:ILE:HG12	2.42	0.49
2:C:68:PHE:HE1	2:C:83:MET:CB	2.26	0.49
2:C:104:ILE:HG22	2:C:105:GLN:C	2.38	0.49
2:D:91:THR:HG23	2:D:122:THR:HA	1.94	0.48
2:C:68:PHE:CE1	2:C:83:MET:HB3	2.48	0.48
1:A:172:HIS:HB3	1:A:185:SER:HB2	1.95	0.48
1:B:226:LEU:O	1:B:230:MET:HG3	2.14	0.48
1:B:145:GLU:HB3	1:B:191:ILE:HD12	1.96	0.48
1:B:201:LEU:HD21	1:B:255:LEU:CD1	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:12:VAL:HG12	2:D:13:GLN:O	2.14	0.48
2:D:95:PHE:CE2	2:D:117:ARG:O	2.67	0.48
2:C:74:ASN:O	2:C:77:ASN:OD1	2.32	0.47
1:B:173:LEU:HD11	1:B:182:LEU:HD22	1.95	0.47
1:A:224:GLU:O	1:A:228:ASN:HB2	2.14	0.47
2:C:67:ARG:O	2:C:84:ASN:HB2	2.15	0.47
2:C:12:VAL:HG12	2:C:13:GLN:O	2.15	0.47
2:C:33:TYR:HB2	2:C:99:SER:HB3	1.97	0.47
1:B:62:ILE:HG12	1:B:87:VAL:HG23	1.96	0.47
1:B:99:GLU:CD	1:B:179:ALA:HB2	2.39	0.47
2:C:48:VAL:HG13	2:C:64:VAL:HG21	1.95	0.47
1:B:184:ILE:HD13	1:B:194:ILE:HG21	1.97	0.46
2:C:99:SER:OG	2:C:104:ILE:HD13	2.15	0.46
1:B:158:GLU:O	1:B:161:PRO:HD2	2.16	0.46
1:B:265:VAL:O	1:B:269:GLN:HG2	2.15	0.46
1:A:99:GLU:CD	1:A:179:ALA:HB2	2.40	0.46
1:B:254:ILE:HG23	1:B:264:LEU:HD23	1.98	0.46
1:B:209:TYR:HB3	1:B:255:LEU:HD21	1.98	0.46
2:D:104:ILE:CG2	2:D:105:GLN:N	2.33	0.45
1:A:53:TRP:CD2	1:A:72:ILE:HG12	2.52	0.45
2:D:23:ALA:HA	2:D:78:THR:HG22	1.99	0.45
1:B:94:GLN:CD	2:D:59:TYR:HB3	2.42	0.45
1:A:91:SER:O	2:C:59:TYR:OH	2.29	0.44
1:B:53:TRP:CD2	1:B:72:ILE:HG12	2.52	0.44
2:D:39:GLN:C	2:D:92:ALA:HB1	2.42	0.44
1:B:173:LEU:HD13	1:B:184:ILE:HG12	2.00	0.44
2:C:6:GLU:CD	2:C:96:CYS:HB3	2.35	0.44
2:D:76:LYS:O	2:D:78:THR:HG23	2.17	0.44
1:A:254:ILE:HG22	1:A:261:ARG:CG	2.48	0.44
1:A:211:VAL:HG11	1:A:251:ARG:HH21	1.81	0.44
1:B:72:ILE:HG22	1:B:97:LEU:HD23	2.00	0.44
2:D:64:VAL:HB	2:D:68:PHE:CD2	2.53	0.43
1:B:56:ASN:N	2:D:103:THR:HG22	2.33	0.43
2:C:4:LEU:HD13	2:C:96:CYS:SG	2.59	0.43
2:C:48:VAL:HG13	2:C:64:VAL:CG2	2.49	0.43
1:B:114:THR:HG23	1:B:127:ASN:OD1	2.19	0.43
2:C:105:GLN:NE2	2:C:111:TYR:CE2	2.87	0.43
1:A:68:GLN:OE1	2:C:103:THR:HG21	2.19	0.43
1:A:55:ILE:HA	2:C:103:THR:HG22	2.00	0.43
1:A:160:ILE:N	1:A:161:PRO:CD	2.81	0.43
2:D:101:ARG:HB2	2:D:104:ILE:CD1	2.38	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:101:ARG:HD2	2:C:104:ILE:HG12	2.01	0.42
1:B:94:GLN:NE2	2:D:59:TYR:HB3	2.33	0.42
2:C:2:VAL:HG23	2:C:27:ASN:H	1.85	0.42
2:D:68:PHE:HA	2:D:82:GLN:O	2.20	0.41
2:D:104:ILE:HG23	2:D:111:TYR:HB3	1.77	0.41
1:B:243:VAL:HG23	1:B:264:LEU:HD21	2.02	0.41
1:B:114:THR:CG2	1:B:127:ASN:HD21	2.33	0.41
1:A:168:PRO:HD3	1:A:186:ASP:OD1	2.20	0.41
2:D:12:VAL:HG12	2:D:13:GLN:N	2.36	0.41
2:D:34:MET:HB2	2:D:79:ILE:HD11	1.93	0.41
1:B:62:ILE:HD12	1:B:84:LYS:C	2.46	0.41
1:A:56:ASN:HB3	2:C:52:TYR:CZ	2.55	0.41
2:C:93:MET:HE3	2:C:118:GLY:HA3	2.02	0.41
1:B:95:LEU:HD11	1:B:148:VAL:HG21	2.02	0.40
2:C:12:VAL:HG12	2:C:13:GLN:N	2.35	0.40
2:C:20:LEU:HG	2:C:83:MET:HE2	2.03	0.40
1:B:165:PRO:HB3	1:B:246:ASP:HA	2.03	0.40
2:D:34:MET:HB2	2:D:53:ILE:HD11	2.03	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:A:215:ARG:O	1:B:233:GLY:O[1_655]	1.99	0.21

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	210/240 (88%)	202 (96%)	8 (4%)	0	100	100
1	B	208/240 (87%)	200 (96%)	8 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	121/153 (79%)	117 (97%)	4 (3%)	0	100	100
2	D	123/153 (80%)	120 (98%)	3 (2%)	0	100	100
All	All	662/786 (84%)	639 (96%)	23 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [\(i\)](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	153/194 (79%)	153 (100%)	0	100	100
1	B	150/194 (77%)	150 (100%)	0	100	100
2	C	94/121 (78%)	94 (100%)	0	100	100
2	D	95/121 (78%)	94 (99%)	1 (1%)	65	88
All	All	492/630 (78%)	491 (100%)	1 (0%)	87	96

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	114	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	GLN
1	A	190	ASN
1	A	213	ASN
1	A	227	ASN
1	A	242	GLN
1	B	94	GLN
1	B	102	GLN
1	B	213	ASN

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Mol	Chain	Res	Type
1	B	234	GLN
1	B	242	GLN
2	C	105	GLN
2	D	13	GLN
2	D	105	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 [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	214/240 (89%)	0.39	9 (4%) 40 32	38, 68, 111, 124	0
1	B	212/240 (88%)	0.53	11 (5%) 33 26	54, 73, 105, 128	0
2	C	123/153 (80%)	0.56	3 (2%) 59 50	45, 78, 113, 121	0
2	D	125/153 (81%)	0.27	1 (0%) 82 77	46, 65, 90, 113	0
All	All	674/786 (85%)	0.44	24 (3%) 46 38	38, 72, 109, 128	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	234	GLN	3.7
1	B	235	ALA	3.5
1	A	128	ALA	3.3
2	C	41	PRO	3.3
2	C	40	GLY	3.2
1	B	232	ARG	2.9
1	B	120	ASP	2.8
1	B	233	GLY	2.8
1	A	116	VAL	2.7
1	B	239	ALA	2.6
1	A	203	GLN	2.5
1	B	94	GLN	2.5
1	B	122	ALA	2.4
2	C	6	GLU	2.2
1	A	118	GLN	2.2
1	B	87	VAL	2.2
2	D	104	ILE	2.1
1	A	87	VAL	2.1
1	B	204	LYS	2.1
1	A	267	LEU	2.1
1	B	159	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	124	ILE	2.1
1	A	142	ASP	2.1
1	A	260	ALA	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.