



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 12:27 am GMT

PDB ID : 4HGK  
Title : Shark IgNAR variable domain  
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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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

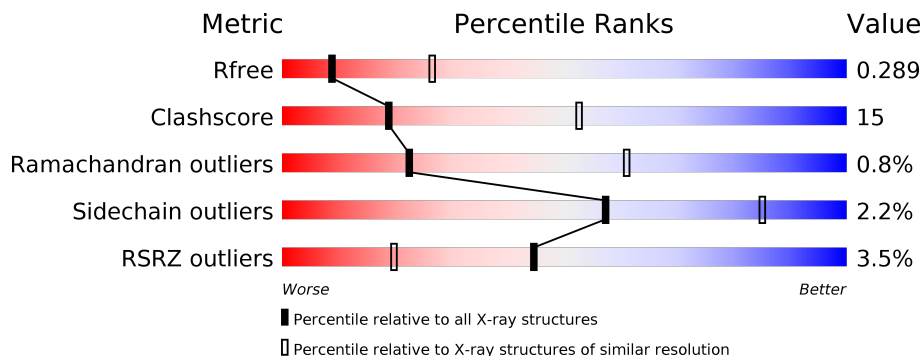
# 1 Overall quality at a glance

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))
$R_{free}$	100719	2176 (3.08-3.00)
Clashscore	112137	2542 (3.08-3.00)
Ramachandran outliers	110173	2458 (3.08-3.00)
Sidechain outliers	110143	2461 (3.08-3.00)
RSRZ outliers	101464	2202 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>3%</div> <div> <div></div> <div>65%</div> <div>16%</div> <div>•</div> <div>17%</div> </div> </div>
1	B	585	<div> <div>4%</div> <div> <div></div> <div>67%</div> <div>15%</div> <div>•</div> <div>18%</div> </div> </div>
2	C	128	<div> <div></div> <div> <div></div> <div>64%</div> <div>15%</div> <div>•</div> <div>19%</div> </div> </div>
2	D	128	<div> <div>%</div> <div> <div></div> <div>67%</div> <div>13%</div> <div>•</div> <div>19%</div> </div> </div>

## 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		



- Molecule 1: Serum albumin



- Chain D:  %

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	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) 92.3 (52.05-3.04)	Depositor EDS
$R_{merge}$	0.22	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.73 (at 3.07Å)	Xtriage
Refinement program	BUSTER 2.9.3	Depositor
R, $R_{free}$	0.237 , 0.266 0.255 , 0.289	Depositor DCC
$R_{free}$ test set	1314 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	68.3	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 69.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	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 <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.42	0/3920	0.65	1/5287 (0.0%)
1	B	0.43	0/3914	0.66	1/5281 (0.0%)
2	C	0.54	0/820	0.70	0/1115
2	D	0.51	0/820	0.68	0/1115
All	All	0.45	0/9474	0.66	2/12798 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	110	PRO	N-CA-C	-7.41	92.84	112.10
1	B	106	LYS	N-CA-C	-6.37	93.81	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3846	0	3775	115	0
1	B	3838	0	3763	110	1
2	C	806	0	790	33	0
2	D	806	0	790	34	1
3	A	6	0	0	0	0
3	B	19	0	0	0	0
3	C	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	6	0	0	0	0
All	All	9335	0	9118	271	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 (271) 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:B:325:VAL:CG2	2:C:109:ILE:CD1	2.35	1.04
1:A:325:VAL:CA	2:D:109:ILE:HD11	1.87	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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:404:GLN:HE21	1:B:428:ARG:HA	1.34	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:C	1:B:178:LEU:HD23	1.93	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:A:109:ASN:O	1:A:110:PRO:C	2.11	0.85
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:O	1:B:178:LEU:HD23	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
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
2:D:20:THR:HG23	2:D:46:THR:HB	1.69	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:370:TYR:O	1:B:373:VAL:HG23	1.90	0.71
1:B:5:SER:HB2	1:B:62:CYS:O	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:77:VAL:O	1:A:80:LEU:C	2.30	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
1:A:178:LEU:C	1:A:180:PRO:HD2	2.12	0.69
2:D:109:ILE:O	2:D:109:ILE:HG22	1.90	0.69
2:C:110:TRP:HE3	2:C:110:TRP:O	1.75	0.69
1:A:120:VAL:HG11	1:A:174:LYS:HB3	1.75	0.67
1:A:77:VAL:HG13	1:A:78:ALA:N	2.08	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:A:178:LEU:C	1:A:180:PRO:CD	2.63	0.67
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: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
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
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: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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:ASP:HB3	1:B:110:PRO:CG	2.32	0.60
1:B:109:ASN:HB3	1:B:466:LYS:HZ3	1.67	0.60
2:C:109:ILE:CG2	2:C:109:ILE:O	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
2:C:110:TRP:C	2:C:110:TRP:CE3	2.76	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:N	2:D:108:ASN:OD1	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:B:109:ASN:O	1:B:110:PRO:C	2.42	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:B:249:ASP:HB3	1:B:252:GLU:CD	2.27	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: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
2:C:110:TRP:CE3	2:C:110:TRP:O	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:278:CYS:HA	1:A:281:LYS:HG3	1.90	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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
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:B:107:ASP:C	1:B:110:PRO:HD2	2.31	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
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
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: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:103:LEU:HD21	1:A:247:HIS:C	2.33	0.48
2:D:96:ASP:O	2:D:97:SER:C	2.49	0.48
1:A:31:LEU:HB2	1:A:39:HIS:HE1	1.78	0.48
1:B:109:ASN:CB	1:B:466:LYS:HZ3	2.27	0.48
1:A:108:ASP:OD1	1:A:197:ARG:HD3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:C:96:ASP:O	2:C:97:SER:C	2.50	0.47
1:B:109:ASN:CG	1:B:466:LYS:NZ	2.68	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:243:THR:O	1:B:247:HIS:N	2.47	0.47
1:B:103:LEU:HD21	1:B:247:HIS:O	2.15	0.47
1:B:437:CYS:HA	1:B:440:HIS:HD2	1.80	0.47
1:A:200:CYS:C	1:A:204:GLN:HG3	2.32	0.46
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:A:474:THR:HG23	1:A:475:LYS:N	2.31	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:A:113:PRO:O	1:A:145:ARG:NH1	2.50	0.45
1:A:167:GLU:CD	1:A:181:LYS:HZ1	2.20	0.45
1:A:5:SER:OG	1:A:62:CYS:HB3	2.17	0.45
1:B:418:VAL:HG12	1:B:419:SER:H	1.82	0.45
1:A:109:ASN:HB3	1:A:466:LYS:HZ3	1.80	0.45
1:B:404:GLN:NE2	1:B:428:ARG:HG3	2.31	0.45
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: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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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
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
2:C:94:VAL:C	2:C:96:ASP:N	2.71	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:283:LEU:C	1:A:283:LEU:HD12	2.38	0.41
1:A:223:PHE:CD1	1:A:272:SER:HB2	2.56	0.41
1:B:109:ASN:N	1:B:110:PRO:CD	2.84	0.41
1:B:108:ASP:OD1	1:B:148:TYR:HD1	2.04	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
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: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

### 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	477/585 (82%)	463 (97%)	11 (2%)	3 (1%)	28	67
1	B	477/585 (82%)	463 (97%)	10 (2%)	4 (1%)	22	61
2	C	102/128 (80%)	95 (93%)	6 (6%)	1 (1%)	18	56
2	D	102/128 (80%)	95 (93%)	6 (6%)	1 (1%)	18	56
All	All	1158/1426 (81%)	1116 (96%)	33 (3%)	9 (1%)	22	61

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	419	SER
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%)	415 (98%)	10 (2%)	54	83
1	B	425/511 (83%)	416 (98%)	9 (2%)	59	86

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	91/110 (83%)	88 (97%)	3 (3%)	43	78
2	D	91/110 (83%)	90 (99%)	1 (1%)	78	93
All	All	1032/1242 (83%)	1009 (98%)	23 (2%)	57	85

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	TYR
1	A	142	ILE
1	A	245	CYS
1	A	281	LYS
1	A	284	LEU
1	A	340	ASP
1	A	423	LEU
1	A	425	GLU
1	A	468	PRO
1	A	492	GLU
1	B	7	VAL
1	B	106	LYS
1	B	112	LEU
1	B	138	TYR
1	B	142	ILE
1	B	245	CYS
1	B	297	GLU
1	B	314	ASP
1	B	392	CYS
2	C	97	SER
2	C	110	TRP
2	C	123	HIS
2	D	108	ASN

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

Mol	Chain	Res	Type
1	A	204	GLN
1	A	247	HIS
1	A	404	GLN
1	A	483	ASN
1	B	39	HIS
1	B	109	ASN

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Mol	Chain	Res	Type
1	B	247	HIS
1	B	404	GLN
1	B	440	HIS
1	B	483	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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	483/585 (82%)	0.06	18 (3%) 42 18	44, 83, 149, 175	0
1	B	481/585 (82%)	-0.01	22 (4%) 33 14	41, 80, 151, 179	0
2	C	104/128 (81%)	-0.43	0 100 100	39, 64, 90, 101	0
2	D	104/128 (81%)	-0.37	1 (0%) 82 58	46, 65, 95, 102	0
All	All	1172/1426 (82%)	-0.05	41 (3%) 44 20	39, 78, 147, 179	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	424	VAL	6.9
1	B	405	ASN	6.4
1	A	85	GLY	6.3
1	B	428	ARG	5.9
1	B	88	ALA	5.7
1	A	423	LEU	5.1
1	B	463	LEU	4.9
1	A	87	MET	4.9
1	B	412	THR	4.8
1	A	419	SER	4.7
1	A	407	LEU	4.5
1	B	452	TYR	4.2
1	B	87	MET	4.0
1	B	408	LEU	4.0
1	A	426	VAL	3.7
1	B	421	PRO	3.6
1	A	430	LEU	3.4
1	A	102	PHE	3.1
1	B	450	GLU	2.9
1	A	94	GLN	2.9
1	B	89	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	439	LYS	2.8
1	A	454	SER	2.8
1	A	442	GLU	2.8
1	B	425	GLU	2.8
1	A	427	SER	2.7
1	A	420	THR	2.7
1	B	409	VAL	2.7
1	A	79	THR	2.7
1	B	92	ALA	2.6
1	A	415	VAL	2.6
1	A	431	GLY	2.6
1	B	91	CYS	2.4
1	B	86	GLU	2.4
1	B	93	LYS	2.4
1	B	416	PRO	2.3
1	A	30	TYR	2.3
2	D	91	ASP	2.3
1	A	458	ASN	2.3
1	B	129	ASP	2.2
1	B	488	PHE	2.1

## 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 carbohydrates 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.