



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 02:27 AM BST

PDB ID : 6OED  
Title : CRYSTAL STRUCTURE OF THE RV144 C1-C2 SPECIFIC ANTIBODY  
CH55 FAB  
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Deposited on : 2019-03-27  
Resolution : 2.46 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

<https://www.wwpdb.org/validation/2017/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)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

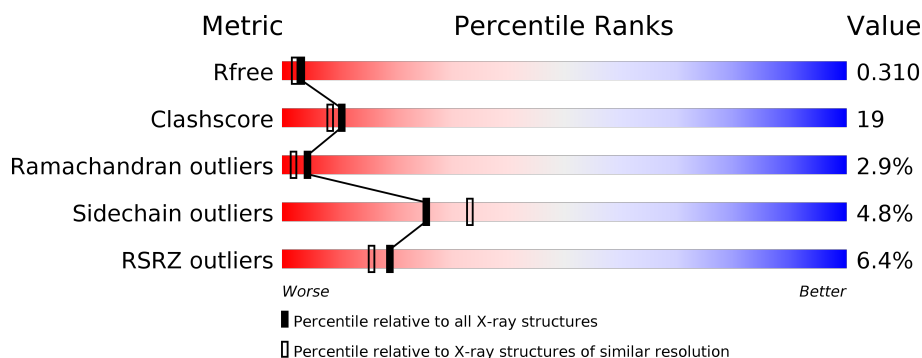
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.46 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))
$R_{free}$	130704	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 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	223	<div> <div>4%</div> <div>65%</div> <div>28%</div> <div>5%</div> </div>
1	C	223	<div> <div>14%</div> <div>52%</div> <div>38%</div> <div>5%</div> </div>
1	H	223	<div> <div>%</div> <div>63%</div> <div>30%</div> <div>.</div> </div>
2	B	215	<div> <div>3%</div> <div>65%</div> <div>30%</div> <div>.</div> </div>
2	D	215	<div> <div>14%</div> <div>61%</div> <div>32%</div> <div>6%</div> </div>
2	L	215	<div> <div>66%</div> <div>30%</div> <div>.</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9840 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 CH55 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	214	Total	C	N	O	S	0	0	0
			1634	1031	282	313	8			
1	A	211	Total	C	N	O	S	0	0	0
			1613	1019	278	308	8			
1	C	211	Total	C	N	O	S	0	0	0
			1613	1019	278	308	8			

- Molecule 2 is a protein called CH55 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	213	Total	C	N	O	S	0	0	0
			1632	1017	280	330	5			
2	B	213	Total	C	N	O	S	0	0	0
			1632	1017	280	330	5			
2	D	213	Total	C	N	O	S	0	0	0
			1632	1017	280	330	5			

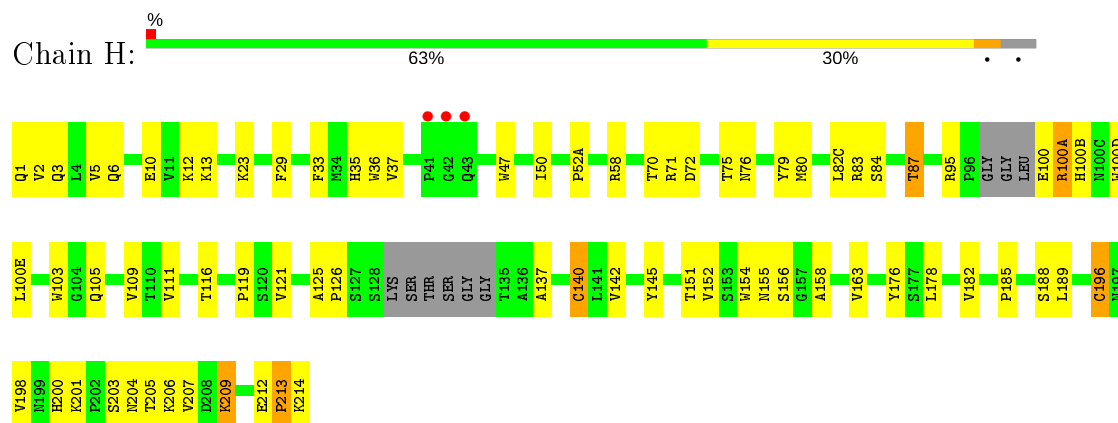
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	18	Total	O	0	0
			18	18		
3	L	16	Total	O	0	0
			16	16		
3	A	11	Total	O	0	0
			11	11		
3	B	16	Total	O	0	0
			16	16		
3	C	9	Total	O	0	0
			9	9		
3	D	14	Total	O	0	0
			14	14		

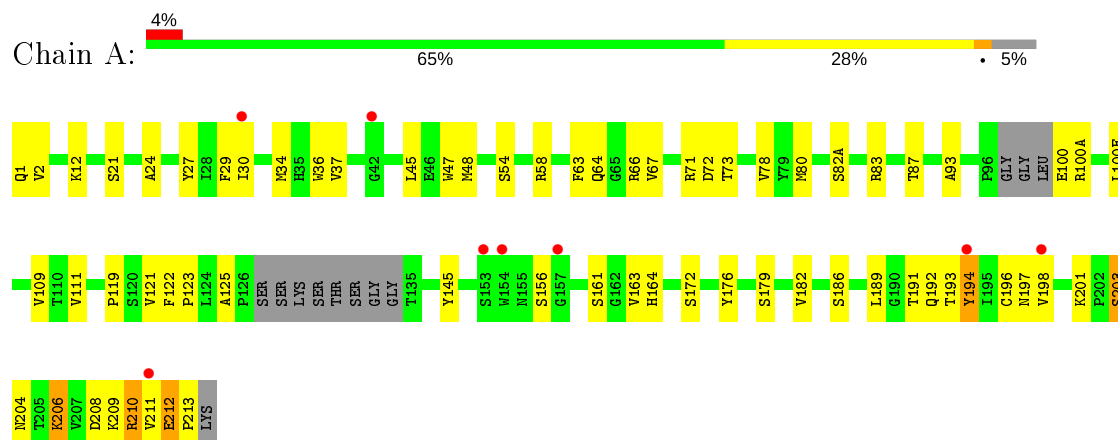
### 3 Residue-property plots [i](#)

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.

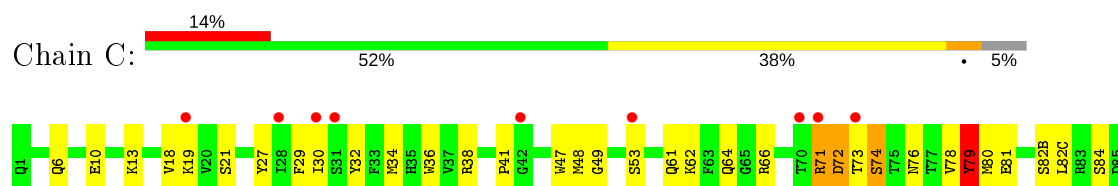
- Molecule 1: CH55 Fab heavy chain

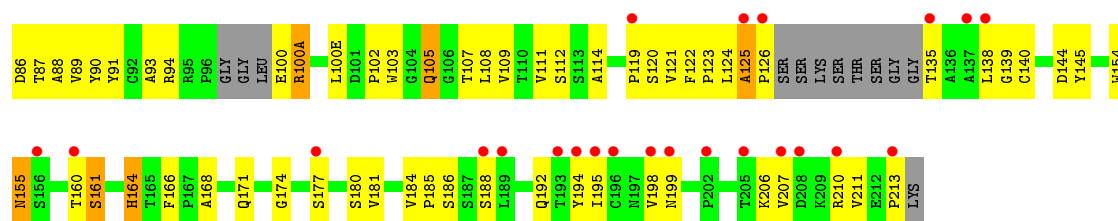


- Molecule 1: CH55 Fab heavy chain



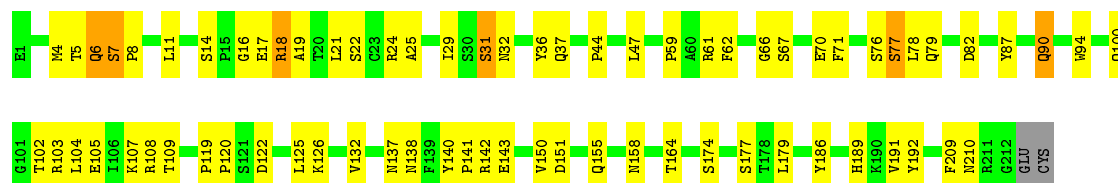
- Molecule 1: CH55 Fab heavy chain





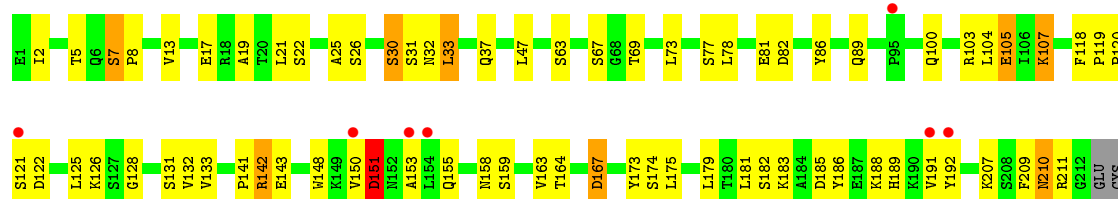
- Molecule 2: CH55 Fab light chain

Chain L: 66% 30% . .



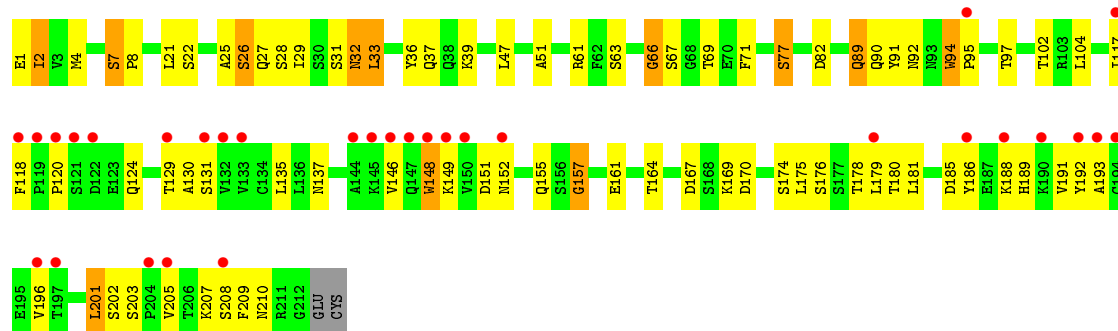
- Molecule 2: CH55 Fab light chain

Chain B: 3% 65% 30% . .



- Molecule 2: CH55 Fab light chain

Chain D: 14% 61% 32% 6% .



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.99 Å 74.03 Å 75.93 Å 106.84° 108.43° 91.78°	Depositor
Resolution (Å)	42.12 – 2.46 60.00 – 2.46	Depositor EDS
% Data completeness (in resolution range)	65.6 (42.12-2.46) 76.8 (60.00-2.46)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.56 (at 2.45 Å)	Xtriage
Refinement program	PHENIX (1.13 _2998: ???)	Depositor
R, $R_{free}$	0.256 , 0.312 0.257 , 0.310	Depositor DCC
$R_{free}$ test set	1991 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.6	Xtriage
Anisotropy	0.641	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 45.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.011 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	9840	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.85% 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

### 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	0.28	0/1653	0.52	0/2252
1	C	0.34	0/1653	0.58	1/2252 (0.0%)
1	H	0.30	0/1674	0.53	0/2279
2	B	0.28	0/1667	0.53	0/2265
2	D	0.31	0/1667	0.57	0/2265
2	L	0.30	0/1667	0.54	1/2265 (0.0%)
All	All	0.30	0/9981	0.55	2/13578 (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	C	79	TYR	CA-CB-CG	5.76	124.34	113.40
2	L	6	GLN	C-N-CA	5.11	134.47	121.70

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	1613	0	1586	46	1
1	C	1613	0	1586	94	0
1	H	1634	0	1609	53	1
2	B	1632	0	1584	56	0
2	D	1632	0	1584	69	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	1632	0	1584	54	0
3	A	11	0	0	1	0
3	B	16	0	0	1	0
3	C	9	0	0	2	0
3	D	14	0	0	0	0
3	H	18	0	0	2	0
3	L	16	0	0	1	0
All	All	9840	0	9533	357	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 19.

All (357) 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:C:19:LYS:HD3	1:C:81:GLU:HG2	1.13	1.12
1:C:126:PRO:HD3	1:C:213:PRO:HA	1.37	1.07
1:C:19:LYS:CD	1:C:81:GLU:HG2	1.92	0.99
1:C:135:THR:N	1:C:186:SER:HG	1.61	0.97
1:C:19:LYS:HD3	1:C:81:GLU:CG	1.96	0.96
1:C:125:ALA:HB3	1:C:126:PRO:CD	1.98	0.93
2:B:5:THR:HA	2:B:100:GLN:HE22	1.41	0.84
1:C:19:LYS:HE2	1:C:79:TYR:HB2	1.58	0.84
1:C:125:ALA:CB	1:C:126:PRO:CD	2.56	0.84
1:C:125:ALA:HB3	1:C:126:PRO:HD2	1.59	0.84
2:B:21:LEU:HD12	2:B:73:LEU:HD23	1.60	0.82
1:C:171:GLN:O	1:C:174:GLY:N	2.13	0.82
1:C:61:GLN:HG2	1:C:62:LYS:HD3	1.62	0.81
1:A:37:VAL:HG12	1:A:47:TRP:HA	1.62	0.81
2:D:28:SER:HA	2:D:69:THR:HG22	1.63	0.81
2:B:158:ASN:HD22	2:B:181:LEU:HD11	1.45	0.80
1:H:5:VAL:HG13	1:H:105:GLN:HE22	1.46	0.80
1:C:121:VAL:HG21	1:C:198:VAL:HG11	1.63	0.80
2:D:151:ASP:OD2	2:D:191:VAL:N	2.14	0.79
2:D:7:SER:HB2	2:D:22:SER:H	1.49	0.78
1:C:19:LYS:HE2	1:C:79:TYR:CB	2.14	0.78
2:L:61:ARG:NH1	2:L:82:ASP:OD1	2.16	0.78
2:L:6:GLN:HE22	2:L:87:TYR:HA	1.48	0.78
1:A:34:MET:HG3	1:A:78:VAL:HG21	1.66	0.78
2:B:81:GLU:OE2	3:B:301:HOH:O	2.01	0.78
2:B:105:GLU:OE1	2:B:142:ARG:NH1	2.17	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:LYS:CE	1:C:79:TYR:HB2	2.12	0.77
1:C:87:THR:HG22	1:C:111:VAL:H	1.48	0.77
2:D:21:LEU:HD22	2:D:102:THR:HG21	1.66	0.77
1:C:29:PHE:HZ	1:C:78:VAL:HG23	1.47	0.77
2:D:191:VAL:HG22	2:D:210:ASN:HD21	1.50	0.77
2:B:7:SER:HB3	2:B:8:PRO:HD3	1.67	0.76
1:C:100:GLU:HG3	1:C:100(A):ARG:HG2	1.66	0.76
2:D:135:LEU:HD21	2:D:137:ASN:HB2	1.70	0.73
1:C:124:LEU:HD22	2:D:118:PHE:HB3	1.70	0.73
1:C:66:ARG:NH1	1:C:86:ASP:OD2	2.22	0.72
1:C:171:GLN:HE21	1:C:177:SER:HB2	1.54	0.72
2:L:6:GLN:NE2	2:L:87:TYR:HA	2.05	0.71
1:A:1:GLN:HG2	1:A:2:VAL:HG23	1.73	0.71
1:A:121:VAL:HG21	1:A:198:VAL:HG11	1.73	0.70
2:D:146:VAL:HG13	2:D:196:VAL:HG12	1.72	0.70
2:B:122:ASP:HA	2:B:125:LEU:HD12	1.74	0.70
1:C:125:ALA:CB	1:C:126:PRO:HD3	2.22	0.69
1:A:156:SER:H	1:A:197:ASN:HD21	1.38	0.69
1:H:126:PRO:HD2	1:H:213:PRO:HB3	1.74	0.69
1:H:70:THR:HG23	1:H:79:TYR:HB2	1.74	0.69
1:H:33:PHE:HB3	1:H:50:ILE:HD11	1.76	0.68
1:A:54:SER:O	1:A:71:ARG:NH2	2.26	0.68
1:C:82(C):LEU:HB3	1:C:111:VAL:HG21	1.75	0.68
1:H:155:ASN:HB3	1:H:158:ALA:HB3	1.76	0.68
2:B:30:SER:OG	2:B:31:SER:N	2.23	0.67
1:H:152:VAL:HG22	1:H:198:VAL:HG12	1.77	0.67
2:D:151:ASP:N	2:D:191:VAL:O	2.23	0.67
1:H:13:LYS:NZ	3:H:302:HOH:O	2.27	0.67
1:C:29:PHE:CZ	1:C:78:VAL:HG23	2.30	0.67
1:H:121:VAL:O	1:H:209:LYS:NZ	2.27	0.67
2:L:24:ARG:NH1	3:L:301:HOH:O	2.22	0.67
2:B:120:PRO:HD3	2:B:132:VAL:HG22	1.77	0.66
1:H:116:THR:HG22	1:H:203:SER:HB3	1.77	0.66
2:L:17:GLU:O	2:L:78:LEU:HD22	1.95	0.66
2:D:1:GLU:O	2:D:2:ILE:HG22	1.95	0.66
2:D:181:LEU:HD13	2:D:185:ASP:HB3	1.78	0.66
2:D:193:ALA:HB2	2:D:208:SER:HB3	1.78	0.66
1:H:142:VAL:HG22	1:H:178:LEU:HG	1.78	0.65
2:B:118:PHE:HB2	2:B:133:VAL:HG23	1.79	0.65
1:H:100(D):TRP:O	2:L:36:TYR:OH	2.14	0.65
1:C:126:PRO:HG3	1:C:213:PRO:O	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:ALA:HB3	1:C:126:PRO:HD3	1.78	0.65
1:C:198:VAL:HG12	1:C:207:VAL:O	1.97	0.65
1:H:154:TRP:CZ3	1:H:196:CYS:HB2	2.32	0.64
2:D:26:SER:O	2:D:27:GLN:HB3	1.96	0.64
2:L:7:SER:HB2	2:L:8:PRO:CD	2.28	0.64
1:C:6:GLN:NE2	1:C:90:TYR:O	2.27	0.64
2:D:36:TYR:OH	2:D:89:GLN:OE1	2.05	0.63
2:L:16:GLY:H	2:L:78:LEU:HB2	1.64	0.63
2:D:124:GLN:HE22	2:D:131:SER:HB3	1.62	0.63
2:B:33:LEU:HA	2:B:89:GLN:O	1.98	0.63
1:C:192:GLN:NE2	3:C:303:HOH:O	2.32	0.63
1:C:195:ILE:HA	1:C:210:ARG:HA	1.81	0.63
2:L:191:VAL:HG12	2:L:210:ASN:ND2	2.14	0.63
1:A:196:CYS:O	1:A:208:ASP:HB2	1.98	0.62
1:A:194:TYR:H	1:A:211:VAL:HG13	1.64	0.62
2:B:189:HIS:O	2:B:211:ARG:NH1	2.32	0.62
2:L:5:THR:HG23	2:L:100:GLN:HE22	1.64	0.62
1:C:125:ALA:HA	1:C:138:LEU:HD22	1.82	0.62
2:D:37:GLN:HB2	2:D:47:LEU:HD11	1.81	0.62
1:H:185:PRO:O	1:H:188:SER:OG	2.15	0.62
1:C:120:SER:HB3	1:C:122:PHE:CE2	2.34	0.61
1:C:66:ARG:NH2	1:C:82(B):SER:O	2.26	0.61
2:L:122:ASP:O	2:L:126:LYS:HE3	2.00	0.61
2:D:7:SER:HB2	2:D:22:SER:O	2.00	0.61
1:H:87:THR:HA	1:H:109:VAL:O	1.99	0.61
2:B:163:VAL:HG12	2:B:175:LEU:HD12	1.82	0.61
2:B:150:VAL:HG23	2:B:153:ALA:HB3	1.82	0.61
1:C:139:GLY:HA2	1:C:154:TRP:HH2	1.66	0.61
1:C:155:ASN:HA	1:C:195:ILE:CG2	2.30	0.60
1:C:19:LYS:CE	1:C:79:TYR:CB	2.78	0.60
1:C:36:TRP:HB3	1:C:48:MET:HE2	1.82	0.60
2:L:29:ILE:O	2:L:32:ASN:HB2	2.00	0.60
1:A:163:VAL:HG22	1:A:182:VAL:HG12	1.84	0.59
2:L:186:TYR:O	2:L:192:TYR:OH	2.20	0.59
1:A:58:ARG:NH1	1:C:100:GLU:OE2	2.35	0.59
1:A:123:PRO:HG3	1:A:210:ARG:HD2	1.84	0.59
1:H:6:GLN:N	1:H:105:GLN:OE1	2.29	0.59
2:L:29:ILE:HG21	2:L:90:GLN:HG3	1.84	0.59
1:H:201:LYS:O	1:H:201:LYS:NZ	2.28	0.59
2:L:143:GLU:N	2:L:143:GLU:OE2	2.34	0.58
2:L:137:ASN:ND2	2:L:138:ASN:OD1	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:17:GLU:O	2:B:78:LEU:HD12	2.03	0.58
1:C:168:ALA:O	3:C:301:HOH:O	2.17	0.57
2:B:19:ALA:HB2	2:B:78:LEU:HD11	1.85	0.57
1:C:122:PHE:CE2	2:D:124:GLN:HB2	2.39	0.57
2:D:130:ALA:N	2:D:181:LEU:O	2.37	0.57
1:H:203:SER:O	1:H:205:THR:N	2.38	0.57
2:L:191:VAL:HG12	2:L:210:ASN:HD21	1.69	0.57
2:B:158:ASN:ND2	2:B:181:LEU:HD11	2.19	0.56
1:C:184:VAL:HG22	1:C:185:PRO:HD2	1.86	0.56
2:L:19:ALA:HB2	2:L:78:LEU:HD11	1.87	0.56
1:C:6:GLN:HB2	1:C:105:GLN:HG3	1.88	0.56
1:C:155:ASN:HA	1:C:195:ILE:HG21	1.88	0.56
1:A:93:ALA:HB3	1:A:100(E):LEU:HD23	1.87	0.56
1:C:89:VAL:HG23	1:C:108:LEU:HD13	1.86	0.55
2:L:59:PRO:HG2	2:L:62:PHE:CE2	2.41	0.55
1:A:1:GLN:OE1	1:A:1:GLN:N	2.33	0.55
1:C:195:ILE:HD12	1:C:210:ARG:CZ	2.36	0.55
2:D:164:THR:HG22	2:D:174:SER:H	1.72	0.55
2:D:29:ILE:HD11	2:D:33:LEU:HB2	1.88	0.55
2:D:7:SER:OG	2:D:8:PRO:HD3	2.05	0.55
1:H:119:PRO:HD2	1:H:205:THR:HG21	1.87	0.55
2:L:14:SER:O	2:L:17:GLU:HG2	2.07	0.55
1:A:210:ARG:HG3	1:A:211:VAL:N	2.22	0.55
2:L:132:VAL:HG23	2:L:179:LEU:HB3	1.89	0.55
2:B:191:VAL:HG23	2:B:210:ASN:HB3	1.89	0.54
2:L:21:LEU:HD22	2:L:102:THR:HG21	1.88	0.54
2:B:192:TYR:HB2	2:B:209:PHE:CE2	2.43	0.54
2:B:37:GLN:HB2	2:B:47:LEU:HD11	1.88	0.54
2:L:120:PRO:HB2	2:L:125:LEU:HD21	1.88	0.54
1:C:53:SER:O	1:C:71:ARG:NH2	2.40	0.54
2:D:155:GLN:OE1	2:D:157:GLY:N	2.39	0.54
1:C:125:ALA:HB2	1:C:211:VAL:HG22	1.90	0.54
1:C:195:ILE:CD1	1:C:210:ARG:CZ	2.86	0.53
2:B:179:LEU:HG	2:B:181:LEU:HD13	1.89	0.53
2:B:186:TYR:HD1	2:B:192:TYR:HH	1.55	0.53
1:A:123:PRO:HB3	1:A:210:ARG:HG2	1.90	0.53
2:D:7:SER:CB	2:D:22:SER:H	2.19	0.53
1:C:199:ASN:OD1	1:C:206:LYS:HD3	2.08	0.53
1:A:123:PRO:O	2:B:121:SER:OG	2.26	0.53
1:C:171:GLN:NE2	1:C:177:SER:HB2	2.24	0.53
1:A:189:LEU:HD23	1:A:189:LEU:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:7:SER:CB	2:B:8:PRO:HD3	2.37	0.53
1:C:87:THR:HA	1:C:109:VAL:O	2.09	0.53
1:H:37:VAL:HG12	1:H:47:TRP:HD1	1.74	0.53
1:C:166:PHE:HE2	2:D:176:SER:HB3	1.74	0.52
2:L:126:LYS:N	2:L:126:LYS:HE2	2.24	0.52
1:A:145:TYR:CE1	1:A:176:TYR:HB2	2.43	0.52
1:A:29:PHE:HE1	1:A:73:THR:HG22	1.73	0.52
1:C:72:ASP:OD1	1:C:73:THR:N	2.42	0.52
2:D:185:ASP:HA	2:D:188:LYS:HG3	1.91	0.52
1:H:140:CYS:HB2	1:H:154:TRP:CH2	2.44	0.52
1:H:87:THR:HB	1:H:111:VAL:H	1.73	0.52
1:H:5:VAL:CG1	1:H:105:GLN:HE22	2.20	0.52
2:L:4:MET:HE1	2:L:25:ALA:HB2	1.90	0.52
2:B:148:TRP:CE2	2:B:179:LEU:HB2	2.45	0.52
2:L:59:PRO:HG2	2:L:62:PHE:HE2	1.74	0.52
2:L:77:SER:O	2:L:77:SER:OG	2.28	0.51
1:H:200:HIS:HB3	1:H:205:THR:HB	1.92	0.51
1:H:72:ASP:HB3	1:H:75:THR:HG22	1.92	0.51
2:L:103:ARG:NH1	2:L:105:GLU:OE2	2.44	0.51
1:C:119:PRO:HB3	1:C:145:TYR:HB3	1.91	0.51
1:H:100(A):ARG:O	1:H:100(A):ARG:HG2	2.11	0.51
1:H:12:LYS:HG3	1:H:82(C):LEU:HD23	1.93	0.51
1:A:211:VAL:HG12	1:A:212:GLU:HB3	1.92	0.51
2:B:164:THR:HG22	2:B:174:SER:H	1.76	0.50
2:D:151:ASP:HA	2:D:191:VAL:HB	1.92	0.50
2:B:185:ASP:O	2:B:189:HIS:ND1	2.44	0.50
2:B:132:VAL:HG11	2:B:209:PHE:HE2	1.77	0.50
1:C:19:LYS:HE3	1:C:79:TYR:HB2	1.93	0.50
2:B:158:ASN:OD1	2:B:158:ASN:N	2.45	0.50
1:H:100(B):HIS:HB3	1:H:100(D):TRP:CZ3	2.47	0.50
2:D:2:ILE:HG21	2:D:90:GLN:NE2	2.27	0.50
1:H:145:TYR:OH	1:H:178:LEU:HD23	2.11	0.50
1:C:18:VAL:HG22	1:C:82(C):LEU:HD21	1.93	0.50
1:C:123:PRO:HA	1:C:140:CYS:HA	1.93	0.49
1:H:2:VAL:O	1:H:3:GLN:HG2	2.12	0.49
1:A:36:TRP:CE2	1:A:80:MET:HB2	2.47	0.49
1:A:64:GLN:O	1:A:64:GLN:NE2	2.40	0.49
2:B:141:PRO:HB2	2:B:143:GLU:OE2	2.13	0.49
2:D:129:THR:HB	2:D:181:LEU:O	2.11	0.49
1:A:24:ALA:HB1	1:A:27:TYR:CE2	2.47	0.49
1:C:27:TYR:CD1	1:C:32:TYR:HD2	2.31	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:103:TRP:CE3	2:L:44:PRO:HD2	2.47	0.49
1:H:151:THR:O	1:H:198:VAL:HA	2.12	0.49
2:L:66:GLY:HA3	2:L:71:PHE:CD1	2.48	0.49
2:L:16:GLY:O	2:L:77:SER:HA	2.12	0.49
2:D:36:TYR:CE2	2:D:89:GLN:OE1	2.66	0.49
1:H:5:VAL:HA	1:H:105:GLN:OE1	2.12	0.49
1:A:100:GLU:O	1:A:100(A):ARG:HB2	2.13	0.49
1:C:29:PHE:CD2	1:C:76:ASN:HA	2.48	0.48
2:B:131:SER:HA	2:B:179:LEU:O	2.13	0.48
2:D:66:GLY:HA3	2:D:71:PHE:CD1	2.48	0.48
2:B:188:LYS:HE2	2:B:188:LYS:HB3	1.59	0.48
2:L:7:SER:CB	2:L:22:SER:H	2.26	0.48
2:L:37:GLN:HB2	2:L:47:LEU:HD22	1.95	0.48
1:H:71:ARG:NH1	3:H:305:HOH:O	2.45	0.48
1:A:122:PHE:HB3	2:B:121:SER:HB2	1.95	0.48
1:A:164:HIS:HE2	2:B:167:ASP:CG	2.17	0.48
1:H:163:VAL:HG22	1:H:182:VAL:HB	1.95	0.48
1:H:36:TRP:CE2	1:H:80:MET:HB2	2.48	0.48
2:D:124:GLN:HE22	2:D:131:SER:CB	2.27	0.48
1:C:181:VAL:HG21	2:D:135:LEU:CD1	2.44	0.48
2:B:105:GLU:OE1	2:B:173:TYR:OH	2.32	0.48
2:L:17:GLU:OE2	2:L:18:ARG:NH1	2.45	0.47
1:H:126:PRO:HB3	1:H:137:ALA:O	2.13	0.47
2:L:151:ASP:OD2	2:L:189:HIS:ND1	2.35	0.47
1:H:12:LYS:HG3	1:H:82(C):LEU:CD2	2.43	0.47
2:B:150:VAL:HG12	2:B:192:TYR:CD1	2.49	0.47
1:C:195:ILE:HD12	1:C:210:ARG:NH2	2.30	0.47
1:C:13:LYS:NZ	1:C:114:ALA:O	2.45	0.47
2:D:117:ILE:HG12	2:D:209:PHE:HD2	1.80	0.47
1:H:142:VAL:CG2	1:H:178:LEU:HG	2.44	0.47
1:A:83:ARG:O	1:A:111:VAL:HG11	2.15	0.47
2:L:164:THR:HG22	2:L:174:SER:H	1.79	0.47
2:B:148:TRP:CD1	2:B:159:SER:HB3	2.50	0.47
2:D:25:ALA:O	2:D:69:THR:HB	2.15	0.46
1:C:125:ALA:HB2	1:C:211:VAL:CG2	2.45	0.46
2:D:191:VAL:CG2	2:D:210:ASN:HD21	2.25	0.46
1:C:84:SER:HA	1:C:111:VAL:HG13	1.96	0.46
2:D:61:ARG:NH2	2:D:82:ASP:OD1	2.48	0.46
2:D:31:SER:OG	2:D:31:SER:O	2.30	0.46
1:A:119:PRO:HB3	1:A:145:TYR:HB3	1.97	0.46
1:H:83:ARG:HG3	1:H:84:SER:H	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:LYS:HD2	1:A:210:ARG:HB2	1.98	0.46
2:D:120:PRO:HG2	2:D:186:TYR:CE1	2.51	0.46
1:C:93:ALA:HB3	1:C:100(E):LEU:HD23	1.98	0.46
1:C:19:LYS:HE3	1:C:79:TYR:C	2.37	0.45
1:C:166:PHE:CE2	2:D:176:SER:HB3	2.50	0.45
1:A:30:ILE:HD12	1:C:30:ILE:HD11	1.98	0.45
2:B:122:ASP:O	2:B:125:LEU:HB2	2.15	0.45
1:C:19:LYS:HE2	1:C:79:TYR:HB3	1.95	0.45
2:D:1:GLU:O	2:D:1:GLU:HG2	2.15	0.45
1:H:29:PHE:CG	1:H:76:ASN:HB2	2.52	0.45
1:C:103:TRP:HZ2	2:D:36:TYR:CE1	2.33	0.45
1:H:156:SER:OG	1:H:156:SER:O	2.34	0.45
2:B:148:TRP:CD1	2:B:179:LEU:HD13	2.52	0.45
2:B:86:TYR:CE2	2:B:104:LEU:HD22	2.51	0.45
2:D:104:LEU:HA	2:D:104:LEU:HD12	1.81	0.45
2:D:7:SER:OG	2:D:8:PRO:CD	2.64	0.45
2:L:119:PRO:HB3	2:L:209:PHE:CE1	2.52	0.45
1:C:125:ALA:HB1	1:C:126:PRO:HD3	1.97	0.45
1:H:145:TYR:CE1	1:H:176:TYR:HB2	2.51	0.45
2:L:7:SER:HB3	2:L:22:SER:HB3	1.99	0.45
1:A:45:LEU:N	3:A:304:HOH:O	2.49	0.45
2:B:7:SER:HB2	2:B:22:SER:HB3	1.98	0.45
1:C:34:MET:HG3	1:C:78:VAL:HG21	1.99	0.45
1:C:168:ALA:HA	1:C:177:SER:O	2.17	0.44
1:C:94:ARG:HB3	1:C:102:PRO:HD2	1.99	0.44
2:D:94:TRP:HB3	2:D:95:PRO:HD3	1.99	0.44
2:B:150:VAL:HG12	2:B:192:TYR:CE1	2.52	0.44
2:D:164:THR:HG22	2:D:174:SER:N	2.33	0.44
2:D:32:ASN:O	2:D:51:ALA:N	2.50	0.44
2:D:39:LYS:HE3	2:D:39:LYS:HB3	1.55	0.44
1:C:18:VAL:O	1:C:81:GLU:HA	2.17	0.44
1:A:100(A):ARG:HD2	1:A:100(A):ARG:HA	1.84	0.44
1:A:192:GLN:HG3	1:A:193:THR:H	1.83	0.44
2:D:161:GLU:OE2	2:D:175:LEU:HD21	2.18	0.44
2:L:4:MET:HA	2:L:24:ARG:O	2.18	0.44
2:D:90:GLN:OE1	2:D:92:ASN:N	2.49	0.44
2:D:77:SER:OG	2:D:77:SER:O	2.34	0.44
2:L:7:SER:HB3	2:L:22:SER:H	1.81	0.44
1:C:6:GLN:N	1:C:105:GLN:HE21	2.16	0.44
2:D:131:SER:HB2	2:D:180:THR:HG22	1.99	0.44
2:D:4:MET:HG2	2:D:97:THR:HG23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:13:VAL:C	2:B:107:LYS:HB3	2.38	0.44
1:A:66:ARG:HG2	1:A:82(A):SER:O	2.18	0.43
2:D:148:TRP:HB3	2:D:179:LEU:HD11	2.00	0.43
1:A:204:ASN:OD1	1:A:206:LYS:HD3	2.18	0.43
1:A:71:ARG:HG2	1:A:72:ASP:N	2.33	0.43
1:C:89:VAL:HG11	1:C:91:TYR:CZ	2.53	0.43
1:H:10:GLU:HG2	1:H:12:LYS:HE3	1.99	0.43
2:L:17:GLU:C	2:L:78:LEU:HD22	2.38	0.43
1:C:100:GLU:O	1:C:100(A):ARG:C	2.56	0.43
2:D:131:SER:HA	2:D:180:THR:HA	1.98	0.43
1:H:35:HIS:CD2	1:H:50:ILE:HD13	2.53	0.43
2:L:158:ASN:OD1	2:L:158:ASN:N	2.52	0.43
1:A:125:ALA:HB1	1:A:213:PRO:HB3	2.01	0.43
1:C:194:TYR:HA	1:C:194:TYR:HD1	1.68	0.43
2:D:178:THR:HG22	2:D:180:THR:HG23	2.01	0.43
1:H:70:THR:OG1	1:H:71:ARG:N	2.52	0.43
2:B:186:TYR:HD1	2:B:192:TYR:OH	2.02	0.43
1:C:19:LYS:NZ	1:C:81:GLU:HG3	2.34	0.43
1:C:6:GLN:NE2	1:C:107:THR:HG23	2.34	0.43
2:D:191:VAL:HG22	2:D:210:ASN:ND2	2.25	0.43
2:D:192:TYR:O	2:D:208:SER:HA	2.19	0.43
2:D:36:TYR:CZ	2:D:89:GLN:OE1	2.71	0.43
2:B:155:GLN:CD	2:B:158:ASN:HD21	2.21	0.43
2:D:36:TYR:HE2	2:D:89:GLN:HG2	1.83	0.43
2:L:11:LEU:HB3	2:L:104:LEU:HD12	2.01	0.43
2:B:207:LYS:HA	2:B:207:LYS:HD3	1.88	0.42
2:D:167:ASP:HB3	2:D:170:ASP:O	2.18	0.42
2:D:151:ASP:CG	2:D:189:HIS:HB3	2.40	0.42
1:H:209:LYS:HD3	1:H:209:LYS:HA	1.77	0.42
1:A:63:PHE:O	1:A:67:VAL:HG12	2.18	0.42
2:B:82:ASP:O	2:B:86:TYR:OH	2.18	0.42
2:B:151:ASP:HA	2:B:191:VAL:HG12	2.01	0.42
1:C:195:ILE:HB	1:C:210:ARG:HG3	2.02	0.42
1:C:181:VAL:HG21	2:D:135:LEU:HD13	2.02	0.42
1:C:164:HIS:HB2	1:C:181:VAL:HG13	2.02	0.42
1:C:210:ARG:N	1:C:210:ARG:HD2	2.35	0.42
1:C:124:LEU:HB3	2:D:118:PHE:CD1	2.54	0.42
1:H:33:PHE:HA	1:H:52(A):PRO:HD3	2.02	0.42
2:B:7:SER:HB3	2:B:8:PRO:CD	2.44	0.42
2:D:201:LEU:H	2:D:201:LEU:HG	1.56	0.42
1:H:119:PRO:HB3	1:H:145:TYR:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:LYS:HE3	1:C:80:MET:N	2.35	0.42
1:H:189:LEU:HD21	1:H:212:GLU:OE1	2.20	0.42
2:L:143:GLU:H	2:L:143:GLU:CD	2.22	0.42
1:A:12:LYS:O	1:A:111:VAL:HA	2.20	0.42
2:B:104:LEU:HA	2:B:104:LEU:HD12	1.82	0.42
1:H:95:ARG:HD3	1:H:100(B):HIS:O	2.20	0.42
2:L:107:LYS:HA	2:L:140:TYR:OH	2.20	0.42
2:L:132:VAL:CG2	2:L:179:LEU:HB3	2.50	0.42
1:A:191:THR:OG1	1:A:192:GLN:N	2.53	0.42
1:A:201:LYS:O	1:A:203:SER:N	2.49	0.42
1:C:73:THR:O	1:C:74:SER:OG	2.30	0.42
1:C:41:PRO:HD3	1:C:88:ALA:HA	2.02	0.42
2:B:119:PRO:HB3	2:B:209:PHE:CE1	2.56	0.41
1:C:160:THR:HG23	1:C:161:SER:H	1.84	0.41
2:D:117:ILE:CG2	2:D:207:LYS:HG2	2.50	0.41
1:C:10:GLU:O	1:C:109:VAL:HA	2.20	0.41
1:C:13:LYS:HD3	1:C:13:LYS:HA	1.89	0.41
1:H:125:ALA:HB1	1:H:213:PRO:HB3	2.03	0.41
1:A:48:MET:HE3	1:A:48:MET:HB2	1.85	0.41
2:L:108:ARG:NH1	2:L:109:THR:O	2.54	0.41
2:L:47:LEU:HA	2:L:47:LEU:HD12	1.93	0.41
2:L:61:ARG:HB3	2:L:76:SER:O	2.20	0.41
2:B:19:ALA:CB	2:B:78:LEU:HD11	2.50	0.41
2:B:182:SER:O	2:B:185:ASP:N	2.54	0.41
2:L:140:TYR:CG	2:L:141:PRO:HA	2.55	0.41
2:L:141:PRO:HB2	2:L:143:GLU:OE2	2.20	0.41
2:L:18:ARG:HD3	2:L:18:ARG:H	1.85	0.41
1:A:87:THR:HA	1:A:109:VAL:O	2.21	0.41
2:B:151:ASP:HB3	2:B:191:VAL:HG12	2.03	0.41
2:D:196:VAL:HG22	2:D:205:VAL:HG12	2.02	0.41
2:B:25:ALA:O	2:B:69:THR:HG23	2.21	0.41
1:H:205:THR:HG23	1:H:207:VAL:HG23	2.02	0.41
2:L:150:VAL:HG22	2:L:155:GLN:NE2	2.36	0.41
1:C:19:LYS:NZ	1:C:81:GLU:CG	2.84	0.41
2:B:126:LYS:N	2:B:126:LYS:HD3	2.36	0.40
1:C:19:LYS:HZ2	1:C:81:GLU:HG3	1.86	0.40
2:D:90:GLN:HG2	2:D:97:THR:HG22	2.03	0.40
2:D:90:GLN:CG	2:D:97:THR:HG22	2.50	0.40
1:H:100:GLU:HB3	1:H:100(A):ARG:H	1.60	0.40
1:A:194:TYR:O	1:A:211:VAL:HA	2.21	0.40
1:C:47:TRP:CZ2	1:C:49:GLY:HA2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:GLN:HG2	1:A:2:VAL:N	2.36	0.40
1:A:71:ARG:HB3	1:A:71:ARG:HE	1.75	0.40
2:L:31:SER:HB3	2:L:32:ASN:H	1.75	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:H:23:LYS:NZ	1:A:192:GLN:NE2[1_566]	2.19	0.01

## 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	205/223 (92%)	185 (90%)	19 (9%)	1 (0%)	29	34
1	C	205/223 (92%)	172 (84%)	26 (13%)	7 (3%)	3	1
1	H	208/223 (93%)	186 (89%)	19 (9%)	3 (1%)	11	9
2	B	211/215 (98%)	181 (86%)	20 (10%)	10 (5%)	2	1
2	D	211/215 (98%)	177 (84%)	23 (11%)	11 (5%)	2	0
2	L	211/215 (98%)	195 (92%)	12 (6%)	4 (2%)	8	6
All	All	1251/1314 (95%)	1096 (88%)	119 (10%)	36 (3%)	4	2

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	204	ASN
2	L	7	SER
2	B	7	SER
1	C	74	SER
1	C	100(A)	ARG

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Mol	Chain	Res	Type
1	C	125	ALA
2	D	7	SER
2	D	94	TRP
1	A	203	SER
2	B	2	ILE
2	B	30	SER
2	B	32	ASN
1	C	155	ASN
2	D	152	ASN
1	H	213	PRO
2	L	67	SER
2	B	33	LEU
2	B	67	SER
2	B	77	SER
2	B	128	GLY
2	B	151	ASP
1	C	64	GLN
1	C	72	ASP
2	D	2	ILE
2	D	77	SER
2	D	169	LYS
1	H	100(E)	LEU
2	L	94	TRP
2	B	105	GLU
2	D	67	SER
2	D	33	LEU
2	D	157	GLY
2	L	77	SER
1	C	144	ASP
2	D	32	ASN
2	D	66	GLY

### 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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	182/190 (96%)	173 (95%)	9 (5%)	25	32
1	C	182/190 (96%)	172 (94%)	10 (6%)	21	27
1	H	185/190 (97%)	176 (95%)	9 (5%)	25	32
2	B	183/185 (99%)	174 (95%)	9 (5%)	25	32
2	D	183/185 (99%)	174 (95%)	9 (5%)	25	32
2	L	183/185 (99%)	176 (96%)	7 (4%)	33	43
All	All	1098/1125 (98%)	1045 (95%)	53 (5%)	25	33

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1	GLN
1	H	58	ARG
1	H	87	THR
1	H	100(A)	ARG
1	H	140	CYS
1	H	196	CYS
1	H	206	LYS
1	H	209	LYS
1	H	214	LYS
2	L	18	ARG
2	L	31	SER
2	L	70	GLU
2	L	79	GLN
2	L	90	GLN
2	L	142	ARG
2	L	177	SER
1	A	21	SER
1	A	161	SER
1	A	172	SER
1	A	179	SER
1	A	186	SER
1	A	194	TYR
1	A	206	LYS
1	A	210	ARG
1	A	212	GLU
2	B	26	SER
2	B	63	SER

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Mol	Chain	Res	Type
2	B	103	ARG
2	B	107	LYS
2	B	142	ARG
2	B	151	ASP
2	B	167	ASP
2	B	183	LYS
2	B	210	ASN
1	C	21	SER
1	C	38	ARG
1	C	71	ARG
1	C	79	TYR
1	C	105	GLN
1	C	112	SER
1	C	161	SER
1	C	164	HIS
1	C	180	SER
1	C	188	SER
2	D	26	SER
2	D	63	SER
2	D	89	GLN
2	D	91	TYR
2	D	148	TRP
2	D	149	LYS
2	D	201	LEU
2	D	202	SER
2	D	203	SER

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

Mol	Chain	Res	Type
1	H	39	GLN
2	L	6	GLN
2	L	210	ASN
2	B	100	GLN
2	B	158	ASN
1	C	39	GLN
1	C	105	GLN
1	C	171	GLN
2	D	124	GLN
2	D	210	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 monosaccharides 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	211/223 (94%)	0.30	8 (3%) 40 37	36, 61, 109, 126	0
1	C	211/223 (94%)	0.84	32 (15%) 2 1	46, 84, 150, 173	0
1	H	214/223 (95%)	-0.02	3 (1%) 75 74	29, 56, 84, 101	0
2	B	213/215 (99%)	0.24	7 (3%) 46 43	27, 50, 108, 124	0
2	D	213/215 (99%)	0.66	31 (14%) 2 1	37, 67, 181, 219	1 (0%)
2	L	213/215 (99%)	0.03	0 100 100	31, 50, 70, 89	0
All	All	1275/1314 (97%)	0.34	81 (6%) 19 16	27, 59, 130, 219	1 (0%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	193	ALA	6.7
2	D	179	LEU	5.6
1	C	19	LYS	5.4
1	C	193	THR	4.9
1	C	126	PRO	4.9
2	D	150	VAL	4.6
1	C	137	ALA	4.4
1	C	125	ALA	4.4
2	D	132	VAL	4.3
1	C	205	THR	4.3
1	C	198	VAL	4.0
2	D	133	VAL	3.9
2	D	146	VAL	3.9
1	C	196	CYS	3.8
2	B	154	LEU	3.6
1	C	70	THR	3.5
2	D	121	SER	3.5
1	A	30	ILE	3.4
1	C	202	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	198	VAL	3.3
1	C	71	ARG	3.2
1	A	154	TRP	3.2
1	C	194	TYR	3.1
2	D	148	TRP	3.1
1	C	199	ASN	3.1
1	C	73	THR	3.1
2	D	117	ILE	3.0
2	D	197	THR	3.0
1	C	30	ILE	3.0
1	C	31	SER	3.0
1	C	119	PRO	3.0
2	D	119	PRO	2.9
2	D	196	VAL	2.8
2	D	190	LYS	2.8
1	C	135	THR	2.8
2	D	152	ASN	2.8
1	A	194	TYR	2.8
2	B	192	TYR	2.8
1	H	43	GLN	2.7
2	D	188	LYS	2.7
1	C	156	SER	2.7
2	D	192	TYR	2.7
2	D	186	TYR	2.6
2	D	149	LYS	2.6
1	C	138	LEU	2.6
2	D	131	SER	2.5
2	D	118	PHE	2.5
2	D	129	THR	2.5
2	D	208	SER	2.5
1	A	42	GLY	2.5
2	D	120	PRO	2.5
2	D	194	CYS	2.4
1	A	211	VAL	2.4
2	D	144	ALA	2.4
1	C	28	ILE	2.4
1	C	53	SER	2.4
2	D	122	ASP	2.3
1	C	42	GLY	2.3
1	C	213	PRO	2.3
2	D	145	LYS	2.3
2	B	191	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	H	42	GLY	2.2
1	C	208	ASP	2.2
1	A	157	GLY	2.2
1	C	177	SER	2.2
1	H	41	PRO	2.2
1	C	207	VAL	2.2
1	A	153	SER	2.1
1	C	160	THR	2.1
2	D	147	GLN	2.1
2	B	153	ALA	2.1
2	D	205	VAL	2.1
2	B	95	PRO	2.1
1	C	188	SER	2.1
2	D	204	PRO	2.1
1	C	195	ILE	2.1
1	C	210	ARG	2.1
2	B	150	VAL	2.0
2	D	95	PRO	2.0
2	B	121	SER	2.0
1	C	189	LEU	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 monosaccharides 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.