



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:56 AM UTC

PDB ID : 1XF5 / pdb_00001xf5
Title : Complex HCV core-Fab 19D9D6-Protein L mutant (H74C, Y64W)in space group P21212
Authors : Menez, R.; Housden, N.G.; Harrison, S.; Jolivet-Reynaud, C.; Gore, M.G.; Stura, E.A.
Deposited on : 2004-09-14
Resolution : 2.60 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

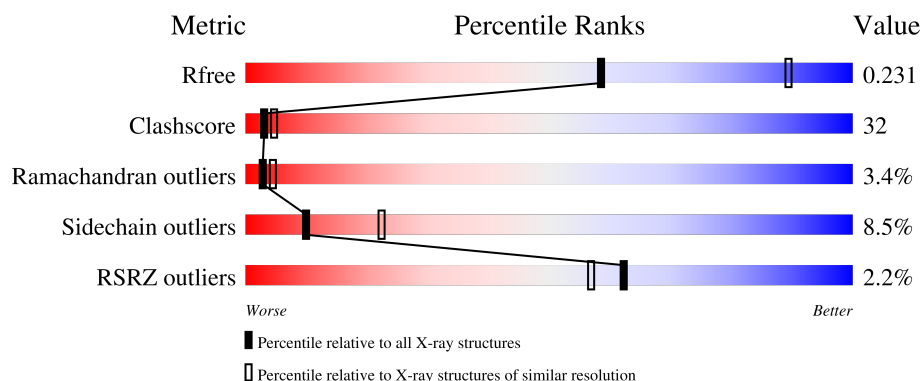
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	P	44	5% 23% 7% 5% . 64%
1	Q	44	11% 14% . 82%
2	A	220	59% 32% 5% .
2	C	220	% 55% 35% 9% .
3	B	218	2% 61% 31% 6% .

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Mol	Chain	Length	Quality of chain
3	D	218	% 58% 35% 6%
4	L	80	4% 32% 45% 8% 11%
4	M	80	5% 44% 31% 9% 16%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8420 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 Capsid protein C.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	P	16	Total	C	N	O	0	0	0
			115	74	23	18			
1	Q	8	Total	C	N	O	0	0	0
			59	40	9	10			

- Molecule 2 is a protein called Monoclonal antibody 19D9D6 Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	220	Total	C	N	O	S	0	0	0
			1708	1063	291	346	8			
2	C	220	Total	C	N	O	S	3	0	0
			1708	1063	291	346	8			

- Molecule 3 is a protein called Monoclonal antibody 19D9D6 Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	218	Total	C	N	O	S	0	0	0
			1660	1058	270	325	7			
3	D	218	Total	C	N	O	S	0	0	0
			1660	1058	270	325	7			

- Molecule 4 is a protein called Protein L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	71	Total	C	N	O	S	0	0	0
			547	346	87	112	2			
4	M	67	Total	C	N	O	S	0	0	0
			517	328	83	104	2			

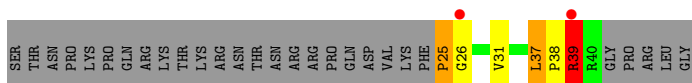
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	P	8	Total 8	O 8	0	0
5	Q	1	Total 1	O 1	0	0
5	A	141	Total 141	O 141	0	0
5	B	115	Total 115	O 115	0	0
5	L	15	Total 15	O 15	0	0
5	C	68	Total 68	O 68	0	0
5	D	75	Total 75	O 75	0	0
5	M	23	Total 23	O 23	0	0

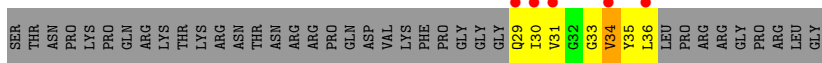
3 Residue-property plots

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.

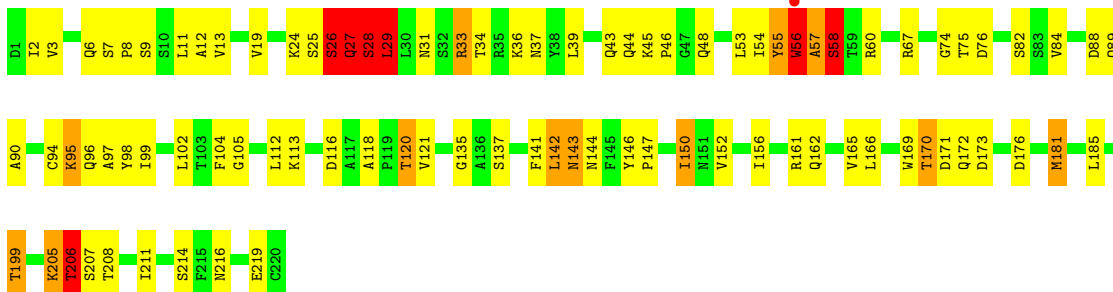
- Molecule 1: Capsid protein C



- Molecule 1: Capsid protein C

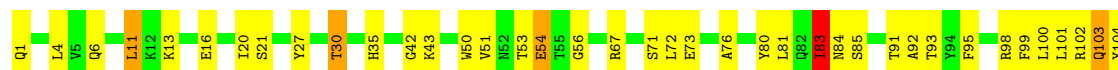


- Molecule 2: Monoclonal antibody 19D9D6 Light chain

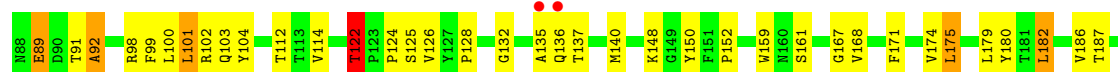
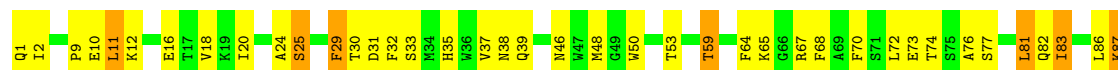




• Molecule 3: Monoclonal antibody 19D9D6 Heavy chain



• Molecule 3: Monoclonal antibody 19D9D6 Heavy chain



• Molecule 4: Protein L



• Molecule 4: Protein L





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	129.34Å 222.93Å 43.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.60 19.92 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.92-2.60) 97.5 (19.92-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.59Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.199 , 0.240 0.217 , 0.231	Depositor DCC
R_{free} test set	1951 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8420	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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	P	0.69	0/117	2.03	6/156 (3.8%)
1	Q	0.48	0/59	1.02	0/79
2	A	0.81	4/1745 (0.2%)	1.31	22/2366 (0.9%)
2	C	0.50	0/1745	1.12	16/2366 (0.7%)
3	B	0.58	1/1707 (0.1%)	1.23	23/2335 (1.0%)
3	D	0.47	0/1707	0.91	4/2335 (0.2%)
4	L	0.64	0/556	1.74	17/748 (2.3%)
4	M	0.46	0/525	0.95	2/705 (0.3%)
All	All	0.60	5/8161 (0.1%)	1.21	90/11090 (0.8%)

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
2	A	0	9
3	B	0	1
All	All	0	10

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	55	TYR	C-N	-14.94	1.12	1.33
2	A	27	GLN	C-N	-6.38	1.25	1.33
2	A	28	SER	N-CA	-5.36	1.39	1.46
2	A	26	SER	C-N	-5.30	1.26	1.33
3	B	83	ILE	CA-C	-5.29	1.46	1.52

All (90) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	27	GLN	CA-C-N	19.25	150.14	123.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	27	GLN	C-N-CA	19.25	150.14	123.00
4	L	18	LYS	N-CA-C	16.72	146.41	110.80
4	L	16	GLU	CA-C-N	-15.13	100.93	119.84
4	L	16	GLU	C-N-CA	-15.13	100.93	119.84
2	A	26	SER	CA-C-N	13.91	148.11	121.54
2	A	26	SER	C-N-CA	13.91	148.11	121.54
4	L	17	PRO	CA-C-N	11.53	143.57	121.54
4	L	17	PRO	C-N-CA	11.53	143.57	121.54
3	B	136	GLN	CA-C-N	9.49	139.66	121.54
3	B	136	GLN	C-N-CA	9.49	139.66	121.54
3	B	136	GLN	N-CA-C	9.47	125.42	110.17
2	C	32	SER	CA-C-N	-9.01	104.95	121.52
2	C	32	SER	C-N-CA	-9.01	104.95	121.52
3	D	122	THR	CA-C-N	8.88	126.05	119.66
3	D	122	THR	C-N-CA	8.88	126.05	119.66
2	C	30	LEU	N-CA-C	8.85	129.66	110.80
4	L	13	THR	N-CA-C	8.83	129.33	109.81
3	B	83	ILE	N-CA-CB	8.72	123.11	111.25
1	P	26	GLY	N-CA-C	-8.62	92.75	113.18
1	P	25	PRO	N-CA-C	8.55	133.46	112.10
3	B	135	ALA	CA-C-N	-8.26	108.60	122.64
3	B	135	ALA	C-N-CA	-8.26	108.60	122.64
3	B	136	GLN	O-C-N	7.79	132.06	123.10
2	C	29	LEU	N-CA-C	-7.79	101.98	112.26
3	B	138	ASN	N-CA-C	-7.77	97.29	109.72
3	B	122	THR	CA-C-N	7.76	125.32	119.66
3	B	122	THR	C-N-CA	7.76	125.32	119.66
2	A	9	SER	N-CA-C	-7.57	102.69	112.23
2	C	33	ARG	CB-CA-C	7.50	124.74	109.67
4	L	19	GLU	N-CA-C	7.25	126.25	110.80
2	A	27	GLN	CA-C-O	-7.15	110.28	120.51
3	B	148	LYS	N-CA-C	7.12	120.76	109.50
4	L	17	PRO	N-CA-C	7.11	127.12	112.47
1	P	39	ARG	CA-C-N	7.11	134.49	121.70
1	P	39	ARG	C-N-CA	7.11	134.49	121.70
3	B	103	GLN	N-CA-C	6.97	121.26	111.92
2	A	26	SER	N-CA-C	6.89	125.47	110.80
4	M	25	VAL	N-CA-C	6.85	117.70	108.11
3	B	83	ILE	CB-CA-C	-6.75	100.77	110.83
2	C	33	ARG	CA-CB-CG	6.75	127.61	114.10
2	A	143	ASN	N-CA-C	6.52	120.52	110.42
2	C	143	ASN	N-CA-C	6.50	120.64	110.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	56	TRP	CB-CA-C	-6.49	97.50	110.42
2	A	25	SER	CA-C-N	6.43	133.82	121.54
2	A	25	SER	C-N-CA	6.43	133.82	121.54
3	D	148	LYS	N-CA-C	6.37	119.85	109.59
2	A	56	TRP	CA-C-O	6.35	129.59	120.51
1	P	37	LEU	CA-C-N	-6.31	111.95	119.84
1	P	37	LEU	C-N-CA	-6.31	111.95	119.84
4	L	18	LYS	N-CA-CB	-6.26	99.92	110.49
2	C	29	LEU	CA-C-N	6.25	133.47	121.54
2	C	29	LEU	C-N-CA	6.25	133.47	121.54
3	B	125	SER	N-CA-C	-6.17	99.64	109.76
4	L	14	PRO	CA-C-N	-6.14	112.79	122.79
4	L	14	PRO	C-N-CA	-6.14	112.79	122.79
4	M	71	GLY	N-CA-C	-6.08	105.22	115.80
2	A	121	VAL	N-CA-C	6.08	117.77	108.71
2	A	82	SER	N-CA-C	5.99	117.88	111.36
2	A	29	LEU	CB-CA-C	5.98	120.77	110.01
3	B	155	VAL	N-CA-C	-5.96	99.63	108.45
2	C	152	VAL	N-CA-C	5.88	116.40	108.17
4	L	14	PRO	N-CA-C	5.87	124.56	112.47
3	B	83	ILE	CA-CB-CG1	5.85	120.35	110.40
2	C	148	LYS	N-CA-C	5.85	119.67	112.54
4	L	17	PRO	O-C-N	5.75	130.40	122.64
3	B	135	ALA	N-CA-C	5.71	122.97	110.80
3	B	109	GLY	N-CA-C	-5.68	103.91	112.89
3	B	136	GLN	CA-C-O	5.68	127.98	121.28
4	L	18	LYS	CA-C-N	5.64	132.32	121.54
4	L	18	LYS	C-N-CA	5.64	132.32	121.54
2	A	56	TRP	O-C-N	-5.58	115.16	122.59
2	A	84	VAL	N-CA-C	5.55	116.58	109.30
2	C	64	VAL	N-CA-C	5.55	114.39	108.95
4	L	15	GLU	N-CA-CB	5.44	119.88	111.39
3	B	171	PHE	N-CA-C	5.44	117.79	110.29
2	C	144	ASN	N-CA-C	5.41	117.78	111.02
2	A	28	SER	CA-C-O	-5.38	114.54	120.30
2	C	53	LEU	N-CA-C	-5.31	105.86	112.93
2	C	151	ASN	N-CA-C	5.31	118.08	108.69
2	C	171	ASP	N-CA-C	-5.31	103.68	110.53
2	A	105	GLY	N-CA-C	-5.29	104.34	112.85
4	L	25	VAL	N-CA-C	5.28	116.33	108.46
2	A	58	SER	CA-C-N	-5.17	115.55	122.42
2	A	58	SER	C-N-CA	-5.17	115.55	122.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	125	SER	N-CA-C	-5.16	100.82	109.24
2	A	120	THR	N-CA-C	-5.09	99.49	108.20
3	B	134	ALA	CA-C-N	5.09	131.25	121.54
3	B	134	ALA	C-N-CA	5.09	131.25	121.54
3	B	43	LYS	N-CA-C	5.07	121.21	113.61

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	26	SER	Mainchain,Peptide
2	A	27	GLN	Mainchain,Sidechain,Peptide
2	A	28	SER	Mainchain
2	A	29	LEU	Mainchain
2	A	56	TRP	Mainchain
2	A	57	ALA	Mainchain
3	B	136	GLN	Mainchain

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	P	115	0	123	17	0
1	Q	59	0	62	14	0
2	A	1708	0	1655	87	0
2	C	1708	0	1658	125	0
3	B	1660	0	1615	82	0
3	D	1660	0	1615	99	0
4	L	547	0	520	68	0
4	M	517	0	499	34	0
5	A	141	0	0	8	0
5	B	115	0	0	5	0
5	C	68	0	0	3	1
5	D	75	0	0	0	1
5	L	15	0	0	3	0
5	M	23	0	0	3	0
5	P	8	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	Q	1	0	0	0	0
All	All	8420	0	7747	496	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 32.

All (496) 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)
3:D:68:PHE:CD1	3:D:83:ILE:HD12	1.90	1.07
3:D:68:PHE:HD1	3:D:83:ILE:HD12	1.07	1.04
2:A:150:ILE:HG23	2:A:181:MET:HE3	1.44	0.99
4:L:18:LYS:HD2	4:L:18:LYS:H	1.25	0.98
2:C:28:SER:OG	2:C:30:LEU:HB3	1.64	0.98
2:C:96:GLN:HE21	2:C:98:TYR:N	1.62	0.98
4:L:18:LYS:H	4:L:18:LYS:CD	1.75	0.96
2:A:169:TRP:HE1	2:A:181:MET:HE2	1.31	0.96
3:B:136:GLN:CD	3:B:137:THR:H	1.74	0.94
3:D:161:SER:H	3:D:201:ASN:HD21	1.16	0.93
3:D:82:GLN:C	3:D:83:ILE:HD13	1.93	0.93
2:C:87:GLU:H	2:C:87:GLU:CD	1.73	0.91
3:B:161:SER:H	3:B:201:ASN:HD21	1.17	0.90
2:C:30:LEU:HD12	2:C:31:ASN:HA	1.52	0.90
2:A:33:ARG:HG2	2:A:33:ARG:HH21	1.37	0.89
2:A:150:ILE:CG2	2:A:181:MET:HE3	2.03	0.89
3:B:135:ALA:O	3:B:136:GLN:NE2	2.05	0.88
2:A:199:THR:HB	2:A:214:SER:HB3	1.55	0.88
1:P:39:ARG:HD2	1:P:39:ARG:H	1.37	0.87
2:C:30:LEU:HD12	2:C:31:ASN:N	1.90	0.86
2:C:96:GLN:NE2	2:C:98:TYR:H	1.71	0.86
4:M:21:VAL:HG21	4:M:43:PHE:HB2	1.58	0.86
3:D:68:PHE:HD1	3:D:83:ILE:CD1	1.91	0.83
1:P:39:ARG:NH2	5:P:284:HOH:O	2.04	0.83
2:C:19:VAL:HG12	2:C:81:ILE:HB	1.58	0.82
3:D:124:PRO:HB3	3:D:150:TYR:HB3	1.62	0.82
2:C:19:VAL:HG13	2:C:81:ILE:HD13	1.61	0.82
2:C:60:ARG:CZ	2:C:66:ASP:HA	2.09	0.81
2:A:169:TRP:NE1	2:A:181:MET:HE2	1.94	0.81
3:D:101:LEU:HD22	3:D:101:LEU:H	1.45	0.80
4:L:18:LYS:HA	4:L:18:LYS:HE3	1.64	0.80
1:P:39:ARG:H	1:P:39:ARG:CD	1.91	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:30:LEU:HD12	2:C:31:ASN:CA	2.12	0.79
3:D:202:VAL:HG13	3:D:211:VAL:HG13	1.64	0.79
3:D:102:ARG:HG3	3:D:102:ARG:HH11	1.48	0.78
3:B:168:VAL:HG22	3:B:186:VAL:HG12	1.65	0.78
3:D:53:THR:HA	3:D:72:LEU:HD11	1.64	0.78
3:D:87:LYS:HZ2	3:D:89:GLU:HG2	1.49	0.77
2:C:31:ASN:HB2	2:C:98:TYR:HE1	1.49	0.76
3:D:20:ILE:HD12	3:D:112:THR:HG21	1.66	0.76
4:L:20:GLU:HB3	5:L:91:HOH:O	1.85	0.76
2:C:96:GLN:HE21	2:C:98:TYR:H	0.85	0.76
4:L:28:ILE:HG12	4:L:34:ILE:HD12	1.68	0.75
2:C:85:GLN:HB3	2:C:87:GLU:OE2	1.87	0.74
2:A:24:LYS:NZ	5:A:297:HOH:O	2.21	0.73
4:L:18:LYS:HE3	4:L:18:LYS:CA	2.18	0.73
3:B:136:GLN:NE2	3:B:136:GLN:HA	2.04	0.73
3:B:161:SER:H	3:B:201:ASN:ND2	1.87	0.73
4:L:17:PRO:HA	4:L:18:LYS:HD2	1.71	0.72
1:Q:29:GLN:CG	3:D:33:SER:HB3	2.19	0.72
1:P:37:LEU:O	1:P:38:PRO:C	2.32	0.72
3:D:2:ILE:HG12	3:D:98:ARG:NH2	2.03	0.72
1:P:39:ARG:CD	1:P:39:ARG:N	2.52	0.71
4:L:23:ILE:HG21	4:L:75:MET:HE3	1.71	0.71
1:P:39:ARG:HD2	1:P:39:ARG:N	2.05	0.71
3:D:140:MET:HE3	3:D:187:THR:HG22	1.72	0.71
3:B:168:VAL:HG22	3:B:186:VAL:CG1	2.21	0.70
2:C:2:ILE:CD1	2:C:2:ILE:H	2.04	0.70
3:D:197:THR:HB	3:D:214:LYS:HE3	1.74	0.69
2:A:206:THR:HG22	5:A:256:HOH:O	1.91	0.69
2:C:166:LEU:HD21	3:D:174:VAL:HB	1.75	0.69
3:D:82:GLN:O	3:D:83:ILE:HD13	1.91	0.69
2:A:118:ALA:HA	2:A:206:THR:HG21	1.75	0.69
2:A:118:ALA:CA	2:A:206:THR:HG21	2.23	0.68
1:Q:29:GLN:HG2	3:D:33:SER:HB3	1.75	0.68
2:C:30:LEU:CD1	2:C:32:SER:H	2.06	0.67
2:C:41:TRP:HD1	2:C:54:ILE:HD11	1.59	0.67
4:M:42:THR:O	4:M:43:PHE:HB3	1.94	0.67
3:B:126:VAL:HG21	3:B:211:VAL:HG11	1.77	0.67
4:L:18:LYS:CA	4:L:18:LYS:CE	2.73	0.67
2:A:150:ILE:CG2	2:A:181:MET:CE	2.73	0.67
2:A:55:TYR:O	2:A:56:TRP:C	2.38	0.67
2:A:118:ALA:HB2	2:A:206:THR:CG2	2.25	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:142:LEU:HD11	2:A:152:VAL:HG22	1.77	0.66
1:P:39:ARG:NH2	5:P:325:HOH:O	2.27	0.66
2:C:81:ILE:HD12	2:C:81:ILE:N	2.10	0.66
4:L:21:VAL:HG22	4:L:22:THR:H	1.60	0.66
4:L:48:ALA:O	4:L:52:ARG:HG3	1.95	0.66
2:C:2:ILE:CD1	2:C:2:ILE:N	2.58	0.66
3:D:182:LEU:C	3:D:182:LEU:HD12	2.21	0.66
4:L:16:GLU:HB3	4:L:17:PRO:HD3	1.78	0.66
4:L:66:ALA:HB2	4:L:77:ILE:HD13	1.77	0.65
2:C:80:THR:C	2:C:81:ILE:HD12	2.21	0.65
2:C:19:VAL:CG1	2:C:81:ILE:HB	2.26	0.65
3:D:38:ASN:OD1	3:D:46:ASN:HB2	1.95	0.65
3:D:20:ILE:CD1	3:D:112:THR:HG21	2.26	0.65
4:L:21:VAL:HG22	4:L:22:THR:N	2.10	0.65
2:C:218:ASN:HD22	2:C:219:GLU:N	1.93	0.65
2:C:56:TRP:CZ2	3:D:102:ARG:HG2	2.32	0.65
2:A:33:ARG:HH21	2:A:33:ARG:CG	2.08	0.64
2:A:36:LYS:HD2	2:A:56:TRP:CG	2.33	0.64
2:C:199:THR:HB	2:C:214:SER:HB3	1.80	0.64
4:M:19:GLU:HB2	4:M:40:LYS:HB3	1.80	0.64
3:B:215:ILE:N	3:B:215:ILE:HD12	2.13	0.64
3:B:99:PHE:CZ	3:B:103:GLN:HA	2.33	0.64
4:M:26:ASN:ND2	4:M:36:THR:HG22	2.11	0.64
3:D:101:LEU:HD22	3:D:101:LEU:N	2.13	0.63
4:M:59:LYS:CE	5:M:103:HOH:O	2.46	0.63
2:A:3:VAL:H	2:A:26:SER:HG	1.45	0.63
3:B:93:THR:HA	3:B:113:THR:HA	1.80	0.63
3:B:138:ASN:HB2	3:B:140:MET:O	1.98	0.63
2:C:141:PHE:C	2:C:142:LEU:HD23	2.24	0.63
2:C:99:ILE:HG13	2:C:99:ILE:O	1.98	0.62
3:D:81:LEU:HD11	3:D:83:ILE:HD11	1.80	0.62
3:D:99:PHE:CZ	3:D:103:GLN:HA	2.33	0.62
2:A:11:LEU:HD13	2:A:19:VAL:CG2	2.30	0.62
4:M:19:GLU:HB2	4:M:40:LYS:HD3	1.82	0.62
4:L:42:THR:HG22	4:L:45:GLU:CG	2.29	0.62
4:M:18:LYS:HE2	4:M:20:GLU:O	2.00	0.62
3:D:70:PHE:CE2	3:D:81:LEU:HD23	2.35	0.62
1:Q:30:ILE:HG22	1:Q:31:VAL:H	1.65	0.62
2:A:3:VAL:N	2:A:26:SER:OG	2.27	0.61
2:C:199:THR:CB	2:C:214:SER:HB3	2.30	0.61
3:D:152:PRO:O	3:D:204:HIS:HE1	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:2:ILE:H	2:C:2:ILE:HD13	1.66	0.61
2:C:152:VAL:HG23	2:C:181:MET:HE1	1.82	0.61
3:D:87:LYS:NZ	3:D:89:GLU:HG2	2.14	0.61
3:D:1:GLN:O	3:D:1:GLN:HG2	2.01	0.61
1:Q:30:ILE:HG22	1:Q:31:VAL:N	2.14	0.61
3:B:1:GLN:OE1	3:B:1:GLN:N	2.25	0.61
4:L:42:THR:HG23	4:L:45:GLU:H	1.64	0.61
2:C:2:ILE:N	2:C:2:ILE:HD12	2.15	0.60
3:B:202:VAL:HG13	3:B:211:VAL:HG13	1.83	0.60
4:M:26:ASN:HD22	4:M:36:THR:HG22	1.66	0.60
3:D:102:ARG:HG3	3:D:102:ARG:NH1	2.17	0.60
2:C:121:VAL:HG22	2:C:142:LEU:HD22	1.84	0.60
2:A:118:ALA:HB2	2:A:206:THR:HG21	1.84	0.60
2:C:98:TYR:CD2	2:C:99:ILE:HG23	2.36	0.60
3:B:93:THR:HG22	3:B:113:THR:HB	1.83	0.60
4:L:51:TYR:OH	4:M:68:LEU:HD21	2.01	0.59
2:C:205:LYS:O	2:C:207:SER:N	2.35	0.59
4:L:18:LYS:CE	4:L:18:LYS:N	2.64	0.59
2:C:188:THR:HG23	5:C:233:HOH:O	2.02	0.59
3:D:126:VAL:O	3:D:213:LYS:HE3	2.02	0.59
2:A:97:ALA:HA	2:A:102:LEU:HD22	1.84	0.59
2:A:170:THR:HG21	5:A:277:HOH:O	2.02	0.59
3:B:124:PRO:HB3	3:B:150:TYR:HB3	1.84	0.59
4:L:16:GLU:CB	4:L:17:PRO:HD3	2.33	0.59
2:A:67:ARG:NH1	2:A:88:ASP:OD1	2.36	0.59
2:A:162:GLN:O	2:A:165:VAL:HG12	2.03	0.59
3:B:113:THR:HG21	5:B:329:HOH:O	2.01	0.59
2:A:141:PHE:C	2:A:142:LEU:HD23	2.27	0.59
2:A:37:ASN:ND2	2:A:74:GLY:H	2.01	0.59
3:D:101:LEU:H	3:D:101:LEU:CD2	2.15	0.58
2:A:199:THR:CB	2:A:214:SER:HB3	2.31	0.58
3:B:215:ILE:HD12	3:B:215:ILE:H	1.69	0.58
3:D:70:PHE:CD2	3:D:81:LEU:HD23	2.38	0.58
2:C:199:THR:HG22	2:C:214:SER:CB	2.33	0.58
2:A:56:TRP:HZ2	5:A:310:HOH:O	1.86	0.58
2:A:118:ALA:CB	2:A:206:THR:HG21	2.33	0.58
2:C:97:ALA:CB	3:D:103:GLN:HB3	2.34	0.58
3:D:216:VAL:HG13	3:D:217:PRO:HD2	1.85	0.58
2:A:54:ILE:HD13	2:A:60:ARG:HA	1.85	0.57
3:D:10:GLU:HG2	3:D:18:VAL:CG2	2.34	0.57
1:P:37:LEU:O	1:P:38:PRO:O	2.22	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:28:ILE:HG23	4:L:34:ILE:HD11	1.87	0.57
4:L:42:THR:HG22	4:L:45:GLU:CD	2.30	0.57
2:C:10:SER:HB3	4:M:38:GLU:HB2	1.86	0.57
3:D:218:ARG:HG3	3:D:218:ARG:O	2.05	0.57
3:B:136:GLN:OE1	3:B:137:THR:HA	2.03	0.57
3:B:132:GLY:HA2	3:B:218:ARG:HD2	1.86	0.56
4:L:17:PRO:HA	4:L:18:LYS:HZ3	1.69	0.56
2:C:169:TRP:CE2	2:C:181:MET:HG3	2.40	0.56
3:D:159:TRP:CZ3	3:D:200:CYS:HB3	2.40	0.56
4:M:20:GLU:HG3	4:M:21:VAL:H	1.70	0.56
4:L:45:GLU:O	4:L:49:GLU:HG2	2.06	0.56
3:B:21:SER:HB3	3:B:80:TYR:CE1	2.40	0.56
3:B:198:VAL:HG12	3:B:215:ILE:HD13	1.87	0.56
2:C:99:ILE:HD12	2:C:100:PRO:O	2.06	0.56
3:D:126:VAL:HG21	3:D:202:VAL:HG11	1.87	0.56
2:A:36:LYS:HD2	2:A:56:TRP:CD1	2.41	0.55
2:C:165:VAL:HA	2:C:184:THR:O	2.07	0.55
3:D:10:GLU:HG2	3:D:18:VAL:HG23	1.89	0.55
3:D:161:SER:H	3:D:201:ASN:ND2	1.95	0.55
3:B:20:ILE:HD11	3:B:81:LEU:HD23	1.87	0.55
3:B:93:THR:HG23	5:B:254:HOH:O	2.06	0.55
2:C:8:PRO:CG	2:C:11:LEU:HG	2.37	0.55
2:C:111:GLU:HG2	5:C:234:HOH:O	2.06	0.55
4:M:28:ILE:HA	4:M:34:ILE:HD13	1.88	0.55
2:A:170:THR:CG2	5:A:277:HOH:O	2.55	0.55
4:L:18:LYS:HA	4:L:18:LYS:CE	2.32	0.55
2:C:69:THR:CG2	2:C:71:ARG:HH11	2.20	0.55
2:C:218:ASN:HD22	2:C:219:GLU:H	1.53	0.55
4:L:43:PHE:HE1	4:M:51:TYR:CD2	2.25	0.55
2:A:67:ARG:HH12	2:A:88:ASP:CG	2.14	0.54
2:A:211:ILE:HD11	5:A:256:HOH:O	2.07	0.54
4:L:51:TYR:HD1	4:M:43:PHE:HZ	1.54	0.54
1:P:25:PRO:HD2	1:P:37:LEU:CD1	2.38	0.54
2:C:69:THR:HG21	2:C:71:ARG:HH11	1.71	0.54
4:M:31:ASP:OD1	4:M:33:LYS:HG3	2.06	0.54
4:L:68:LEU:HD13	4:L:75:MET:HG2	1.89	0.54
3:B:113:THR:HG23	5:B:236:HOH:O	2.06	0.54
2:A:7:SER:HA	2:A:8:PRO:C	2.32	0.54
2:A:142:LEU:HD23	2:A:142:LEU:N	2.23	0.54
1:Q:31:VAL:HG12	3:D:99:PHE:CE1	2.43	0.54
2:A:36:LYS:HE3	2:A:56:TRP:CE3	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:122:THR:HG21	3:D:179:LEU:HD11	1.90	0.54
2:C:205:LYS:C	2:C:207:SER:H	2.15	0.54
3:D:2:ILE:HG21	3:D:98:ARG:HH21	1.73	0.53
4:L:66:ALA:CB	4:L:77:ILE:HD13	2.37	0.53
4:M:52:ARG:HG2	5:M:94:HOH:O	2.07	0.53
1:Q:33:GLY:N	2:C:98:TYR:O	2.32	0.53
3:B:67:ARG:HD2	3:B:85:SER:HB2	1.89	0.53
3:D:83:ILE:HD13	3:D:83:ILE:N	2.23	0.53
4:L:17:PRO:HA	4:L:18:LYS:NZ	2.24	0.53
2:A:199:THR:HG22	2:A:214:SER:HB2	1.91	0.53
2:C:30:LEU:HD12	2:C:30:LEU:C	2.33	0.53
2:C:125:PRO:HB3	2:C:215:PHE:CZ	2.44	0.53
3:B:136:GLN:CG	3:B:137:THR:H	2.19	0.52
4:L:28:ILE:HG23	4:L:34:ILE:CD1	2.39	0.52
2:C:12:ALA:HB3	4:M:36:THR:OG1	2.09	0.52
2:C:153:LYS:HD3	2:C:155:LYS:HE3	1.91	0.52
3:D:48:MET:HG2	3:D:64:PHE:CZ	2.44	0.52
1:P:37:LEU:C	1:P:38:PRO:O	2.52	0.52
4:L:42:THR:CG2	4:L:45:GLU:H	2.22	0.52
2:C:148:LYS:HB3	2:C:179:TYR:CD1	2.44	0.52
3:B:73:GLU:HG2	3:B:76:ALA:HB3	1.92	0.52
2:A:44:GLN:O	2:A:90:ALA:HB1	2.10	0.52
3:B:161:SER:N	3:B:201:ASN:HD21	1.98	0.52
2:C:60:ARG:NH2	2:C:66:ASP:HB3	2.25	0.52
2:A:39:LEU:HD13	2:A:39:LEU:C	2.35	0.52
2:C:44:GLN:HE22	3:D:39:GLN:HE22	1.58	0.52
2:C:96:GLN:NE2	2:C:99:ILE:H	2.08	0.52
2:A:98:TYR:CD2	2:A:99:ILE:HG13	2.44	0.52
2:C:199:THR:HG22	2:C:214:SER:HB2	1.91	0.52
3:B:11:LEU:HG	3:B:152:PRO:HG3	1.91	0.52
2:C:28:SER:CB	2:C:30:LEU:HB3	2.40	0.52
2:C:87:GLU:CD	2:C:87:GLU:N	2.53	0.52
2:C:148:LYS:HG2	5:C:260:HOH:O	2.08	0.52
4:L:18:LYS:H	4:L:18:LYS:CE	2.23	0.51
2:C:28:SER:C	2:C:29:LEU:HD12	2.35	0.51
1:Q:29:GLN:OE1	3:D:33:SER:HA	2.10	0.51
4:L:42:THR:HG22	4:L:45:GLU:OE2	2.10	0.51
2:C:24:LYS:O	2:C:24:LYS:HG3	2.09	0.51
2:C:37:ASN:ND2	2:C:74:GLY:H	2.09	0.51
2:A:37:ASN:HD21	2:A:74:GLY:H	1.57	0.51
1:Q:36:LEU:HD12	1:Q:36:LEU:N	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:6:GLN:HG3	2:C:94:CYS:SG	2.50	0.51
2:C:217:ARG:HH11	2:C:217:ARG:HG2	1.74	0.51
2:A:36:LYS:HD2	2:A:56:TRP:CD2	2.46	0.51
3:B:98:ARG:NH1	3:B:106:ASP:OD2	2.44	0.51
3:D:2:ILE:HG12	3:D:98:ARG:HH22	1.76	0.51
3:B:189:PRO:HG2	3:B:192:THR:CG2	2.40	0.51
3:D:91:THR:O	3:D:92:ALA:HB2	2.10	0.51
2:A:205:LYS:C	2:A:207:SER:H	2.19	0.51
2:A:43:GLN:HB2	2:A:53:LEU:HD11	1.92	0.50
4:L:25:VAL:HG22	4:L:75:MET:HB2	1.93	0.50
2:C:36:LYS:HG2	2:C:56:TRP:CD1	2.46	0.50
3:D:126:VAL:HG21	3:D:211:VAL:CG1	2.42	0.50
3:B:159:TRP:HZ3	3:B:215:ILE:HD11	1.76	0.50
2:C:89:GLN:HE21	2:C:112:LEU:HG	1.75	0.50
2:C:218:ASN:ND2	2:C:219:GLU:N	2.58	0.50
2:C:201:GLU:HG2	2:C:212:VAL:HG12	1.92	0.50
3:B:6:GLN:HE21	3:B:109:GLY:HA3	1.75	0.50
3:D:81:LEU:CD1	3:D:83:ILE:HD11	2.41	0.50
1:P:38:PRO:HD3	2:A:33:ARG:CB	2.42	0.50
4:L:31:ASP:OD1	4:L:33:LYS:HG3	2.12	0.50
2:A:2:ILE:N	2:A:2:ILE:HD12	2.26	0.50
2:C:56:TRP:O	2:C:57:ALA:HB3	2.12	0.50
2:C:97:ALA:HB1	3:D:103:GLN:HB3	1.94	0.50
3:B:132:GLY:O	3:B:134:ALA:N	2.44	0.50
3:B:136:GLN:HA	3:B:136:GLN:HE21	1.70	0.50
4:M:28:ILE:HG23	4:M:34:ILE:HD11	1.93	0.50
3:D:68:PHE:CD1	3:D:83:ILE:CD1	2.77	0.50
1:Q:34:VAL:O	1:Q:34:VAL:HG22	2.12	0.49
2:C:19:VAL:CG1	2:C:81:ILE:HD13	2.36	0.49
2:C:28:SER:OG	2:C:30:LEU:CB	2.47	0.49
3:D:11:LEU:HD12	3:D:152:PRO:HD3	1.94	0.49
3:B:133:SER:O	3:B:134:ALA:HB2	2.12	0.49
2:A:7:SER:HA	2:A:8:PRO:O	2.13	0.49
2:A:24:LYS:HA	2:A:75:THR:O	2.12	0.49
2:C:7:SER:HA	2:C:8:PRO:C	2.37	0.49
4:L:21:VAL:N	4:L:41:GLY:O	2.44	0.49
3:D:9:PRO:HG3	3:D:205:PRO:HB2	1.94	0.49
1:Q:31:VAL:HG11	3:D:35:HIS:CE1	2.48	0.49
2:C:8:PRO:HG3	2:C:11:LEU:HG	1.94	0.49
2:C:195:HIS:O	2:C:217:ARG:NH2	2.44	0.49
4:M:42:THR:O	4:M:43:PHE:CB	2.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:170:THR:HG23	2:A:171:ASP:O	2.12	0.49
2:A:27:GLN:C	2:A:75:THR:HG22	2.37	0.48
2:A:150:ILE:HG23	2:A:181:MET:CE	2.29	0.48
4:L:42:THR:HG21	5:L:90:HOH:O	2.12	0.48
2:C:169:TRP:CZ2	2:C:181:MET:HE2	2.48	0.48
3:D:2:ILE:HD12	3:D:2:ILE:H	1.78	0.48
1:P:25:PRO:HD2	1:P:37:LEU:HD11	1.94	0.48
1:P:38:PRO:HA	1:P:39:ARG:HD2	1.95	0.48
3:B:152:PRO:HD2	3:B:206:ALA:CB	2.43	0.48
2:A:135:GLY:HA3	5:A:266:HOH:O	2.13	0.48
3:B:93:THR:HG22	3:B:113:THR:CG2	2.44	0.48
2:C:4:MET:HE3	2:C:23:CYS:SG	2.54	0.48
2:C:142:LEU:HD11	2:C:152:VAL:HG22	1.94	0.48
3:D:30:THR:HG22	3:D:53:THR:O	2.13	0.48
4:M:58:ALA:O	4:M:59:LYS:C	2.56	0.48
4:M:68:LEU:HD11	4:M:75:MET:HE2	1.96	0.48
3:B:126:VAL:CG2	3:B:211:VAL:HG11	2.43	0.48
3:B:198:VAL:O	3:B:215:ILE:HD13	2.13	0.48
2:A:13:VAL:HG12	4:L:35:GLN:HG2	1.96	0.48
2:A:211:ILE:HD12	2:A:211:ILE:N	2.29	0.48
3:B:13:LYS:HB2	3:B:16:GLU:CD	2.38	0.48
4:L:81:GLY:O	4:L:82:LYS:HB3	2.14	0.48
4:L:82:LYS:OXT	4:L:82:LYS:HD3	2.13	0.48
2:C:166:LEU:HD23	2:C:166:LEU:C	2.38	0.48
3:D:73:GLU:HG2	3:D:76:ALA:HB3	1.96	0.48
3:B:93:THR:HG22	3:B:113:THR:CB	2.43	0.48
2:C:30:LEU:HD12	2:C:32:SER:H	1.77	0.48
2:C:148:LYS:HB3	2:C:179:TYR:CG	2.49	0.48
1:Q:33:GLY:C	1:Q:35:TYR:H	2.22	0.48
2:C:31:ASN:HB2	2:C:98:TYR:CE1	2.40	0.48
3:B:42:GLY:HA2	5:B:298:HOH:O	2.14	0.47
3:D:161:SER:N	3:D:201:ASN:HD21	1.98	0.47
2:A:67:ARG:NH1	2:A:88:ASP:CG	2.71	0.47
2:C:60:ARG:HH22	2:C:66:ASP:HB3	1.79	0.47
4:M:28:ILE:HD12	4:M:78:LYS:HG2	1.96	0.47
3:B:6:GLN:HE22	3:B:95:PHE:HA	1.79	0.47
4:L:18:LYS:N	4:L:18:LYS:NZ	2.63	0.47
2:A:31:ASN:HB3	2:A:34:THR:OG1	2.14	0.47
2:C:92:TYR:N	2:C:92:TYR:CD1	2.82	0.47
2:C:153:LYS:CD	2:C:155:LYS:HE3	2.44	0.47
3:D:182:LEU:C	3:D:182:LEU:CD1	2.88	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:137:SER:HA	2:A:185:LEU:O	2.15	0.47
2:A:118:ALA:CB	2:A:206:THR:CG2	2.93	0.47
3:B:126:VAL:HG21	3:B:211:VAL:CG1	2.44	0.47
4:L:18:LYS:CD	4:L:18:LYS:N	2.51	0.47
2:C:156:ILE:HD12	2:C:161:ARG:NH2	2.30	0.47
3:B:11:LEU:HB2	3:B:152:PRO:HG3	1.98	0.46
2:C:30:LEU:HD13	2:C:32:SER:H	1.78	0.46
3:D:11:LEU:C	3:D:11:LEU:CD2	2.88	0.46
3:D:24:ALA:O	3:D:25:SER:HB3	2.14	0.46
4:M:20:GLU:OE2	4:M:20:GLU:HA	2.15	0.46
4:M:66:ALA:HA	4:M:76:ASN:O	2.14	0.46
2:C:45:LYS:HB3	2:C:46:PRO:HD2	1.97	0.46
2:C:52:VAL:HG11	3:D:104:TYR:CD1	2.50	0.46
3:D:126:VAL:HG21	3:D:211:VAL:HG11	1.97	0.46
4:L:61:ASN:O	4:L:81:GLY:HA2	2.15	0.46
3:D:11:LEU:C	3:D:11:LEU:HD23	2.39	0.46
2:A:143:ASN:HB3	2:A:144:ASN:ND2	2.30	0.46
2:A:12:ALA:O	2:A:113:LYS:HD2	2.15	0.46
3:B:91:THR:O	3:B:92:ALA:HB2	2.16	0.46
2:C:152:VAL:CG2	2:C:181:MET:HE1	2.46	0.46
1:P:38:PRO:HD3	2:A:33:ARG:HB2	1.97	0.46
2:C:96:GLN:NE2	2:C:98:TYR:N	2.43	0.46
3:B:27:TYR:CE2	3:B:98:ARG:HD2	2.50	0.46
4:L:12:LYS:N	5:L:97:HOH:O	2.48	0.46
4:L:42:THR:HG22	4:L:45:GLU:CB	2.46	0.46
4:L:67:ASP:HB2	4:L:76:ASN:HB2	1.97	0.46
3:D:218:ARG:NH2	3:D:218:ARG:HB2	2.31	0.46
4:M:81:GLY:O	4:M:82:LYS:HB3	2.16	0.46
4:L:82:LYS:HD3	4:L:82:LYS:C	2.40	0.46
2:A:56:TRP:CH2	3:B:102:ARG:HG2	2.50	0.45
3:D:2:ILE:HD12	3:D:2:ILE:N	2.30	0.45
3:D:100:LEU:HD12	3:D:104:TYR:CZ	2.51	0.45
3:D:175:LEU:HD22	3:D:180:TYR:CE2	2.51	0.45
2:A:29:LEU:O	2:A:37:ASN:HA	2.16	0.45
2:A:33:ARG:CG	2:A:33:ARG:NH2	2.74	0.45
2:A:43:GLN:NE2	2:A:45:LYS:HE3	2.32	0.45
3:D:50:TRP:CE2	3:D:59:THR:HG21	2.52	0.45
2:C:8:PRO:HG2	2:C:11:LEU:HG	1.97	0.45
2:C:31:ASN:OD1	2:C:31:ASN:C	2.58	0.45
2:C:42:TYR:CE2	2:C:52:VAL:HG12	2.51	0.45
2:C:55:TYR:C	2:C:57:ALA:N	2.72	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:60:ARG:O	2:C:61:GLU:C	2.60	0.45
3:D:2:ILE:HG12	3:D:98:ARG:HH21	1.78	0.45
3:B:51:VAL:HG13	3:B:51:VAL:O	2.15	0.45
3:B:83:ILE:N	3:B:83:ILE:HD13	2.31	0.45
2:C:190:ASP:O	2:C:194:ARG:HG3	2.16	0.45
4:L:42:THR:HG22	4:L:45:GLU:HB2	1.98	0.45
2:C:30:LEU:CD1	2:C:30:LEU:C	2.90	0.45
2:C:42:TYR:OH	2:C:95:LYS:NZ	2.36	0.45
4:L:16:GLU:HB3	4:L:17:PRO:CD	2.44	0.45
2:C:27:GLN:O	2:C:29:LEU:HD12	2.16	0.45
3:D:37:VAL:HG13	3:D:46:ASN:O	2.16	0.45
2:A:67:ARG:NH1	2:A:88:ASP:OD2	2.44	0.45
3:B:189:PRO:HG2	3:B:192:THR:HG23	1.98	0.45
2:C:42:TYR:CE2	2:C:52:VAL:CG1	3.00	0.45
2:A:206:THR:O	2:A:206:THR:HG23	2.17	0.44
3:B:198:VAL:O	3:B:215:ILE:CD1	2.65	0.44
4:L:51:TYR:HD1	4:M:43:PHE:CZ	2.34	0.44
4:L:69:GLU:HG3	4:L:74:CYS:HB3	1.99	0.44
3:B:129:LEU:HB2	3:B:144:GLY:C	2.42	0.44
2:A:89:GLN:NE2	2:A:172:GLN:HG2	2.32	0.44
2:A:95:LYS:HB2	2:A:104:PHE:CD2	2.53	0.44
4:L:43:PHE:CE1	4:M:51:TYR:HD2	2.35	0.44
3:D:48:MET:HG2	3:D:64:PHE:CE2	2.53	0.44
3:D:70:PHE:HE2	3:D:81:LEU:HD23	1.82	0.44
3:B:56:GLY:HA2	3:B:72:LEU:HD11	1.99	0.44
4:L:21:VAL:CG2	4:L:22:THR:H	2.29	0.44
2:C:199:THR:HG22	2:C:214:SER:HB3	2.00	0.44
2:C:166:LEU:HD21	3:D:174:VAL:CB	2.43	0.44
3:B:134:ALA:HB1	3:B:135:ALA:H	1.43	0.44
4:L:22:THR:OG1	4:L:40:LYS:HD3	2.17	0.44
1:P:31:VAL:HG21	3:B:50:TRP:CE2	2.53	0.44
1:Q:30:ILE:CG2	1:Q:31:VAL:H	2.31	0.44
2:A:211:ILE:HD12	2:A:211:ILE:H	1.83	0.44
2:C:199:THR:CG2	2:C:214:SER:HB3	2.48	0.43
2:A:48:GLN:HB3	5:A:252:HOH:O	2.19	0.43
3:B:151:PHE:CE1	3:B:152:PRO:HB3	2.53	0.43
2:C:166:LEU:HD23	2:C:167:ASN:N	2.33	0.43
3:D:168:VAL:HG22	3:D:186:VAL:HG23	1.99	0.43
3:B:30:THR:HG23	3:B:53:THR:O	2.19	0.43
3:B:30:THR:CG2	3:B:53:THR:O	2.66	0.43
3:B:159:TRP:CZ2	3:B:186:VAL:HG22	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:21:VAL:CG2	4:L:22:THR:N	2.78	0.43
2:C:30:LEU:O	2:C:31:ASN:C	2.60	0.43
2:C:60:ARG:HD3	2:C:64:VAL:O	2.18	0.43
3:D:128:PRO:CB	3:D:215:ILE:HD13	2.48	0.43
2:A:24:LYS:HG2	2:A:76:ASP:OD2	2.18	0.43
2:C:55:TYR:C	2:C:57:ALA:H	2.27	0.43
2:C:95:LYS:HE2	2:C:95:LYS:HB3	1.75	0.43
3:D:10:GLU:HG2	3:D:18:VAL:HG21	2.01	0.43
3:D:32:PHE:CG	3:D:98:ARG:HD3	2.54	0.43
2:A:28:SER:HA	2:A:74:GLY:O	2.19	0.43
2:A:116:ASP:OD1	2:A:205:LYS:HD2	2.19	0.43
3:B:71:SER:HB2	5:B:303:HOH:O	2.18	0.43
2:C:141:PHE:O	2:C:142:LEU:HD23	2.18	0.43
3:D:102:ARG:NH1	3:D:102:ARG:CG	2.80	0.43
4:M:60:VAL:O	4:M:60:VAL:CG2	2.65	0.43
3:D:64:PHE:HB3	3:D:68:PHE:CD2	2.53	0.43
2:A:118:ALA:HB2	2:A:206:THR:HG23	2.00	0.43
2:C:97:ALA:HB3	3:D:103:GLN:HB3	2.01	0.43
3:B:93:THR:CG2	3:B:113:THR:HG22	2.49	0.43
3:B:122:THR:HG21	3:B:179:LEU:HD21	2.01	0.43
4:L:68:LEU:CD1	4:L:75:MET:HG2	2.49	0.43
3:D:18:VAL:HG12	3:D:86:LEU:HD11	2.01	0.43
3:D:32:PHE:HE1	3:D:101:LEU:HD21	1.84	0.43
3:D:126:VAL:CG2	3:D:211:VAL:HG11	2.49	0.43
2:A:56:TRP:C	2:A:58:SER:H	2.27	0.43
2:C:54:ILE:HD12	2:C:54:ILE:C	2.43	0.43
3:D:64:PHE:O	3:D:65:LYS:HB2	2.19	0.43
2:A:146:TYR:HA	2:A:147:PRO:C	2.43	0.42
3:B:161:SER:N	3:B:201:ASN:ND2	2.60	0.42
3:D:11:LEU:HD23	3:D:12:LYS:N	2.33	0.42
2:A:216:ASN:HB2	2:A:219:GLU:OE2	2.19	0.42
3:B:124:PRO:CB	3:B:150:TYR:HB3	2.49	0.42
4:L:18:LYS:N	4:L:18:LYS:HZ2	2.17	0.42
2:C:41:TRP:CD1	2:C:54:ILE:HD11	2.48	0.42
2:A:6:GLN:HG3	2:A:94:CYS:SG	2.59	0.42
4:L:76:ASN:HD22	4:L:76:ASN:HA	1.68	0.42
2:A:173:ASP:HB3	2:A:176:ASP:OD2	2.19	0.42
4:L:43:PHE:CE1	4:M:51:TYR:CD2	3.06	0.42
1:Q:30:ILE:CG2	1:Q:31:VAL:N	2.81	0.42
3:B:136:GLN:CD	3:B:137:THR:N	2.59	0.42
4:L:16:GLU:CB	4:L:17:PRO:CD	2.98	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:18:LYS:HB3	4:L:19:GLU:H	1.20	0.42
2:C:121:VAL:O	2:C:213:LYS:HE3	2.19	0.42
4:M:59:LYS:HE3	5:M:103:HOH:O	2.18	0.42
3:B:141:VAL:HG12	3:B:188:VAL:O	2.20	0.42
3:D:218:ARG:HB2	3:D:218:ARG:CZ	2.49	0.42
4:M:69:GLU:O	4:M:70:ASP:HB2	2.20	0.42
1:P:25:PRO:HD2	1:P:37:LEU:HD12	2.02	0.42
4:L:56:LEU:HD23	4:L:59:LYS:HD3	2.00	0.42
3:D:152:PRO:HD2	3:D:206:ALA:CB	2.50	0.42
3:D:171:PHE:CD1	3:D:171:PHE:N	2.88	0.42
2:A:96:GLN:CD	2:A:96:GLN:C	2.88	0.41
2:C:56:TRP:CH2	3:D:102:ARG:HG2	2.55	0.41
4:M:42:THR:HG23	4:M:45:GLU:OE2	2.20	0.41
2:A:176:ASP:OD2	2:A:176:ASP:C	2.64	0.41
4:L:70:ASP:O	4:L:73:ASN:ND2	2.53	0.41
3:B:175:LEU:HB2	3:B:180:TYR:CE1	2.55	0.41
4:L:60:VAL:HG23	4:L:61:ASN:OD1	2.20	0.41
2:A:56:TRP:CZ2	3:B:102:ARG:HG2	2.56	0.41
2:A:156:ILE:HD12	2:A:161:ARG:HB2	2.02	0.41
3:B:11:LEU:C	3:B:11:LEU:HD22	2.45	0.41
3:B:71:SER:OG	3:B:80:TYR:HB2	2.20	0.41
3:B:84:ASN:O	3:B:85:SER:C	2.64	0.41
3:D:67:ARG:NH1	3:D:87:LYS:HE3	2.35	0.41
2:C:81:ILE:N	2:C:81:ILE:CD1	2.79	0.41
2:C:123:ILE:HD12	2:C:200:CYS:SG	2.60	0.41
2:C:169:TRP:CH2	2:C:181:MET:HE2	2.55	0.41
4:M:19:GLU:OE2	4:M:40:LYS:HD2	2.21	0.41
3:B:213:LYS:HD3	3:B:213:LYS:HA	1.72	0.41
4:L:28:ILE:HA	4:L:34:ILE:HD13	2.02	0.41
4:L:13:THR:HA	4:L:14:PRO:HD2	1.79	0.41
2:C:132:THR:O	2:C:132:THR:HG22	2.20	0.41
2:C:218:ASN:HD22	2:C:218:ASN:N	2.18	0.41
3:D:74:THR:C	3:D:77:SER:H	2.29	0.41
3:D:167:GLY:O	3:D:186:VAL:HA	2.21	0.41
3:B:35:HIS:N	3:B:35:HIS:CD2	2.89	0.41
3:B:93:THR:HG22	3:B:113:THR:HG22	2.03	0.41
3:B:179:LEU:HD23	3:B:179:LEU:HA	1.87	0.41
4:L:66:ALA:HB2	4:L:77:ILE:CD1	2.49	0.41
2:C:43:GLN:HB2	2:C:53:LEU:HD11	2.02	0.41
3:D:29:PHE:CD2	3:D:77:SER:HA	2.56	0.41
3:D:73:GLU:O	3:D:74:THR:HB	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:124:PRO:CA	3:B:150:TYR:HB3	2.51	0.41
2:C:85:GLN:CB	2:C:87:GLU:OE2	2.65	0.40
2:C:206:THR:O	2:C:207:SER:HB2	2.22	0.40
1:Q:31:VAL:O	1:Q:31:VAL:HG23	2.21	0.40
3:B:30:THR:HG22	3:B:54:GLU:HA	2.02	0.40
4:M:25:VAL:HG22	4:M:75:MET:HB3	2.02	0.40
3:B:136:GLN:CD	3:B:137:THR:HA	2.46	0.40
4:L:21:VAL:HG11	4:L:43:PHE:HA	2.04	0.40
2:C:218:ASN:ND2	2:C:219:GLU:HG3	2.36	0.40
2:A:171:ASP:O	2:A:172:GLN:C	2.64	0.40
3:B:100:LEU:HD12	3:B:104:TYR:CZ	2.56	0.40
1:P:38:PRO:HD3	2:A:33:ARG:HB3	2.03	0.40
2:C:212:VAL:O	2:C:212:VAL:HG23	2.22	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)
5:C:285:HOH:O	5:D:278:HOH:O[1_554]	2.06	0.14

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	P	14/44 (32%)	10 (71%)	4 (29%)	0	100	100
1	Q	6/44 (14%)	2 (33%)	3 (50%)	1 (17%)	0	0
2	A	218/220 (99%)	203 (93%)	9 (4%)	6 (3%)	4	6
2	C	218/220 (99%)	194 (89%)	18 (8%)	6 (3%)	4	6
3	B	216/218 (99%)	197 (91%)	14 (6%)	5 (2%)	5	9
3	D	216/218 (99%)	193 (89%)	15 (7%)	8 (4%)	2	3

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	L	69/80 (86%)	61 (88%)	2 (3%)	6 (9%)	0	0
4	M	65/80 (81%)	55 (85%)	7 (11%)	3 (5%)	2	2
All	All	1022/1124 (91%)	915 (90%)	72 (7%)	35 (3%)	3	4

All (35) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	B	133	SER
4	L	16	GLU
4	L	17	PRO
4	L	18	LYS
4	L	19	GLU
2	C	31	ASN
2	C	46	PRO
2	C	206	THR
3	D	135	ALA
3	D	136	GLN
3	D	137	THR
2	A	27	GLN
2	A	56	TRP
3	B	135	ALA
4	L	14	PRO
2	C	82	SER
3	D	29	PHE
3	D	132	GLY
4	M	43	PHE
2	A	26	SER
2	A	57	ALA
2	A	58	SER
3	B	134	ALA
3	B	137	THR
4	M	18	LYS
3	B	161	SER
2	C	28	SER
3	D	16	GLU
3	D	25	SER
2	A	206	THR
4	L	13	THR
3	D	92	ALA
4	M	20	GLU
1	Q	34	VAL

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Mol	Chain	Res	Type
2	C	50	PRO

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	P	11/37 (30%)	10 (91%)	1 (9%)	9	19
1	Q	6/37 (16%)	6 (100%)	0	100	100
2	A	195/195 (100%)	181 (93%)	14 (7%)	13	30
2	C	195/195 (100%)	179 (92%)	16 (8%)	10	24
3	B	187/187 (100%)	167 (89%)	20 (11%)	6	14
3	D	187/187 (100%)	173 (92%)	14 (8%)	12	28
4	L	54/63 (86%)	48 (89%)	6 (11%)	6	12
4	M	51/63 (81%)	47 (92%)	4 (8%)	11	26
All	All	886/964 (92%)	811 (92%)	75 (8%)	10	22

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	P	39	ARG
2	A	33	ARG
2	A	46	PRO
2	A	95	LYS
2	A	112	LEU
2	A	120	THR
2	A	142	LEU
2	A	150	ILE
2	A	166	LEU
2	A	170	THR
2	A	181	MET
2	A	199	THR
2	A	205	LYS
2	A	206	THR

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Mol	Chain	Res	Type
2	A	208	THR
3	B	4	LEU
3	B	11	LEU
3	B	30	THR
3	B	54	GLU
3	B	83	ILE
3	B	101	LEU
3	B	113	THR
3	B	114	VAL
3	B	122	THR
3	B	136	GLN
3	B	137	THR
3	B	138	ASN
3	B	152	PRO
3	B	154	PRO
3	B	164	LEU
3	B	182	LEU
3	B	186	VAL
3	B	187	THR
3	B	202	VAL
3	B	211	VAL
4	L	15	GLU
4	L	18	LYS
4	L	19	GLU
4	L	36	THR
4	L	56	LEU
4	L	57	LEU
2	C	2	ILE
2	C	30	LEU
2	C	32	SER
2	C	33	ARG
2	C	36	LYS
2	C	46	PRO
2	C	79	LEU
2	C	81	ILE
2	C	87	GLU
2	C	99	ILE
2	C	112	LEU
2	C	165	VAL
2	C	175	LYS
2	C	199	THR
2	C	212	VAL

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Mol	Chain	Res	Type
2	C	218	ASN
3	D	11	LEU
3	D	31	ASP
3	D	59	THR
3	D	81	LEU
3	D	83	ILE
3	D	87	LYS
3	D	89	GLU
3	D	101	LEU
3	D	114	VAL
3	D	122	THR
3	D	175	LEU
3	D	182	LEU
3	D	202	VAL
3	D	211	VAL
4	M	56	LEU
4	M	60	VAL
4	M	68	LEU
4	M	82	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	37	ASN
2	A	43	GLN
2	A	48	GLN
2	A	85	GLN
2	A	89	GLN
2	A	144	ASN
2	A	163	ASN
2	A	167	ASN
2	A	196	ASN
3	B	6	GLN
3	B	39	GLN
3	B	88	ASN
3	B	169	HIS
3	B	201	ASN
4	L	35	GLN
4	L	73	ASN
4	L	76	ASN
2	C	37	ASN
2	C	43	GLN

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Mol	Chain	Res	Type
2	C	85	GLN
2	C	89	GLN
2	C	96	GLN
2	C	143	ASN
2	C	216	ASN
2	C	218	ASN
3	D	39	GLN
3	D	138	ASN
3	D	201	ASN
3	D	204	HIS
4	M	26	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	55:TYR	C	56:TRP	N	1.12

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	P	16/44 (36%)	0.67	2 (12%) 8 6	38, 57, 94, 96	0
1	Q	8/44 (18%)	2.11	5 (62%) 0 0	104, 109, 111, 111	0
2	A	220/220 (100%)	-0.37	1 (0%) 87 85	17, 32, 52, 75	0
2	C	220/220 (100%)	-0.01	2 (0%) 81 78	24, 45, 73, 90	1 (0%)
3	B	218/218 (100%)	-0.40	4 (1%) 67 63	15, 29, 48, 93	0
3	D	218/218 (100%)	0.01	2 (0%) 81 78	23, 46, 71, 97	0
4	L	71/80 (88%)	0.24	3 (4%) 40 35	29, 45, 124, 137	0
4	M	67/80 (83%)	0.08	4 (5%) 27 22	26, 37, 88, 113	0
All	All	1038/1124 (92%)	-0.11	23 (2%) 62 57	15, 38, 74, 137	1 (0%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	135	ALA	3.8
2	C	30	LEU	3.7
3	B	135	ALA	3.3
3	B	137	THR	3.2
4	L	13	THR	3.0
2	A	56	TRP	2.9
1	Q	29	GLN	2.9
4	M	18	LYS	2.7
2	C	38	TYR	2.6
1	Q	36	LEU	2.6
4	M	19	GLU	2.6
1	Q	31	VAL	2.5
3	B	136	GLN	2.4
1	Q	34	VAL	2.4
1	P	39	ARG	2.3
4	L	17	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
4	M	16	GLU	2.3
3	D	136	GLN	2.2
4	L	14	PRO	2.2
4	M	17	PRO	2.1
1	Q	30	ILE	2.1
3	B	138	ASN	2.1
1	P	26	GLY	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.