



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:46 AM UTC

PDB ID : 3PXJ / pdb_00003pxj
Title : Tandem Ig repeats of Dlar
Authors : Biersmith, B.H.; Bouyain, S.
Deposited on : 2010-12-10
Resolution : 2.30 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

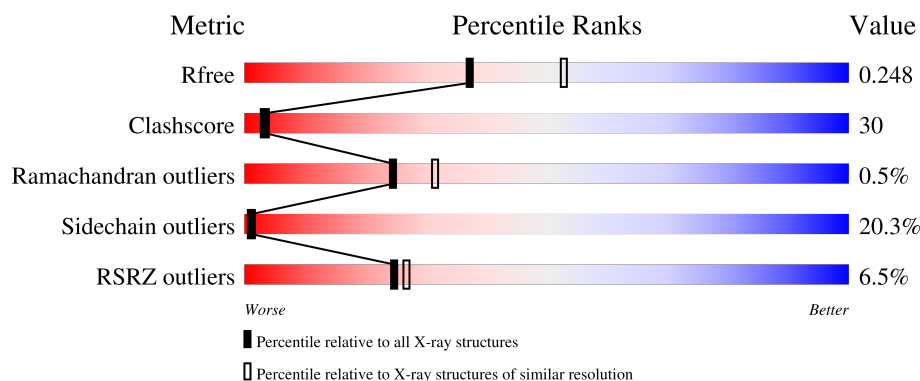
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	210	6% 48% 36% 10% 5%
1	B	210	5% 50% 34% 10% 5%
1	C	210	6% 44% 35% 10% 11%
1	D	210	8% 46% 36% 10% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6103 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 Tyrosine-protein phosphatase Lar.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	199	Total	C	N	O	S	0	0	0
			1519	946	273	293	7			
1	B	199	Total	C	N	O	S	0	0	0
			1519	946	273	293	7			
1	C	187	Total	C	N	O	S	0	0	0
			1441	904	259	271	7			
1	D	194	Total	C	N	O	S	0	0	0
			1486	928	267	284	7			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	GLY	-	expression tag	UNP P16621
A	29	PRO	-	expression tag	UNP P16621
A	30	GLY	-	expression tag	UNP P16621
A	31	SER	-	expression tag	UNP P16621
B	28	GLY	-	expression tag	UNP P16621
B	29	PRO	-	expression tag	UNP P16621
B	30	GLY	-	expression tag	UNP P16621
B	31	SER	-	expression tag	UNP P16621
C	28	GLY	-	expression tag	UNP P16621
C	29	PRO	-	expression tag	UNP P16621
C	30	GLY	-	expression tag	UNP P16621
C	31	SER	-	expression tag	UNP P16621
D	28	GLY	-	expression tag	UNP P16621
D	29	PRO	-	expression tag	UNP P16621
D	30	GLY	-	expression tag	UNP P16621
D	31	SER	-	expression tag	UNP P16621

- Molecule 2 is water.

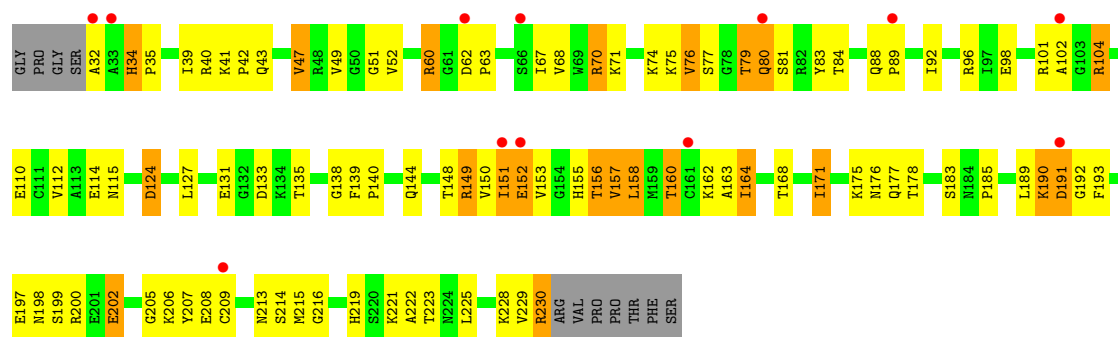
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	27	Total 27	O 27	0	0
2	B	43	Total 43	O 43	0	0
2	C	40	Total 40	O 40	0	0
2	D	28	Total 28	O 28	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

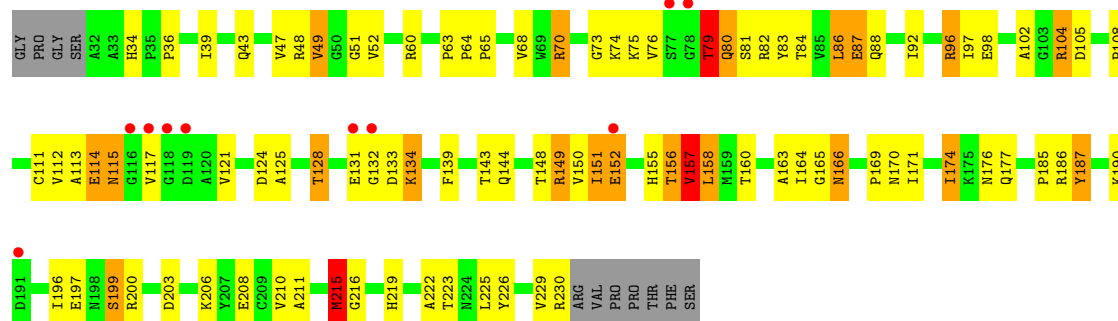
• Molecule 1: Tyrosine-protein phosphatase Lar

Chain A: 



• Molecule 1: Tyrosine-protein phosphatase Lar

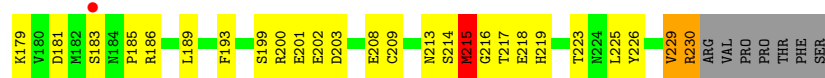
Chain B: 



• Molecule 1: Tyrosine-protein phosphatase Lar

Chain C: 





● Molecule 1: Tyrosine-protein phosphatase Lar



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.96Å 77.51Å 81.73Å 90.00° 101.08° 90.00°	Depositor
Resolution (Å)	41.38 – 2.30 41.38 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.0 (41.38-2.30) 92.8 (41.38-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
R, R_{free}	0.223 , 0.261 0.214 , 0.248	Depositor DCC
R_{free} test set	1964 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	45.5	Xtriage
Anisotropy	0.465	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 63.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6103	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.13 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8634e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.69	3/1551 (0.2%)	1.12	7/2106 (0.3%)
1	B	1.81	17/1551 (1.1%)	1.20	16/2106 (0.8%)
1	C	1.66	4/1470 (0.3%)	1.13	4/1992 (0.2%)
1	D	1.64	4/1516 (0.3%)	1.12	7/2055 (0.3%)
All	All	1.70	28/6088 (0.5%)	1.14	34/8259 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	208	GLU	C-N	15.43	1.55	1.33
1	B	114	GLU	C-N	10.62	1.50	1.33
1	B	115	ASN	C-N	8.30	1.44	1.33
1	B	144	GLN	C-N	-7.87	1.21	1.33
1	B	166	ASN	C-O	-7.20	1.19	1.25
1	B	143	THR	C-N	-6.83	1.24	1.33
1	D	91	GLY	C-N	-6.34	1.21	1.33
1	C	148	THR	C-N	-6.33	1.25	1.33
1	B	51	GLY	C-O	-5.99	1.18	1.23
1	A	51	GLY	C-O	-5.90	1.18	1.23
1	C	199	SER	C-O	-5.87	1.16	1.23
1	A	76	VAL	CA-CB	-5.86	1.46	1.54
1	D	92	ILE	C-N	-5.72	1.25	1.33
1	D	166	ASN	C-O	-5.68	1.20	1.25
1	B	76	VAL	CA-CB	-5.61	1.47	1.54
1	B	125	ALA	C-O	-5.59	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	111	CYS	C-O	-5.35	1.17	1.24
1	B	210	VAL	C-O	-5.34	1.18	1.24
1	B	211	ALA	CA-CB	-5.32	1.46	1.53
1	B	196	ILE	C-O	-5.21	1.18	1.24
1	B	174	ILE	C-O	-5.20	1.18	1.24
1	B	187	TYR	C-O	-5.17	1.17	1.24
1	B	43	GLN	C-O	-5.17	1.17	1.23
1	C	112	VAL	C-O	-5.11	1.18	1.24
1	B	68	VAL	C-O	-5.06	1.18	1.24
1	B	112	VAL	C-O	-5.06	1.18	1.24
1	A	68	VAL	C-O	-5.04	1.19	1.24
1	C	157	VAL	C-O	-5.02	1.18	1.24

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	114	GLU	CA-C-N	-11.79	103.73	122.24
1	B	114	GLU	C-N-CA	-11.79	103.73	122.24
1	B	79	THR	N-CA-C	9.22	121.10	111.14
1	B	115	ASN	O-C-N	8.79	132.01	122.34
1	A	215	MET	N-CA-C	8.11	120.20	111.36
1	A	34	HIS	CA-C-N	7.70	125.20	119.66
1	A	34	HIS	C-N-CA	7.70	125.20	119.66
1	B	114	GLU	O-C-N	7.57	132.55	123.24
1	D	139	PHE	N-CA-C	-6.54	101.06	110.08
1	D	82	ARG	N-CA-C	-6.49	105.36	113.28
1	D	209	CYS	O-C-N	6.36	130.63	123.19
1	C	171	ILE	N-CA-C	6.29	118.08	108.71
1	C	215	MET	N-CA-C	6.28	118.20	111.36
1	D	124	ASP	N-CA-C	6.16	118.31	109.14
1	A	192	GLY	N-CA-C	-6.16	107.41	115.47
1	B	143	THR	O-C-N	-6.00	115.31	122.15
1	D	212	GLU	N-CA-C	6.00	118.83	109.52
1	B	81	SER	N-CA-C	5.69	120.07	113.18
1	A	102	ALA	N-CA-C	5.67	117.14	111.07
1	A	124	ASP	N-CA-C	5.66	118.28	109.52
1	C	82	ARG	N-CA-C	-5.59	106.12	113.17
1	D	102	ALA	N-CA-C	5.57	117.03	111.07
1	B	102	ALA	N-CA-C	5.55	117.33	111.28
1	B	157	VAL	CB-CA-C	-5.41	102.59	110.81
1	B	144	GLN	CA-C-N	5.33	130.24	121.87
1	B	144	GLN	C-N-CA	5.33	130.24	121.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	34	HIS	N-CA-C	5.25	118.14	109.58
1	D	62	ASP	CB-CA-C	-5.12	105.10	110.33
1	B	144	GLN	O-C-N	-5.11	117.33	123.31
1	B	143	THR	CA-C-N	5.10	130.59	122.74
1	B	143	THR	C-N-CA	5.10	130.59	122.74
1	B	215	MET	N-CA-C	5.08	116.89	111.36
1	B	124	ASP	N-CA-C	5.06	116.68	109.14
1	C	143	THR	N-CA-C	5.03	116.58	111.14

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	96	ARG	Sidechain

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	1519	0	1500	105	0
1	B	1519	0	1500	78	0
1	C	1441	0	1433	98	0
1	D	1486	0	1471	87	0
2	A	27	0	0	2	0
2	B	43	0	0	6	0
2	C	40	0	0	7	0
2	D	28	0	0	4	0
All	All	6103	0	5904	351	0

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 30.

All (351) 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:130:TYR:CE1	1:C:136:PRO:HD3	1.80	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:ARG:HH11	1:B:70:ARG:CG	1.61	1.13
1:A:77:SER:OG	1:A:80:GLN:HB2	1.48	1.10
1:B:36:PRO:HG3	1:B:115:ASN:HB2	1.27	1.10
1:A:70:ARG:NH1	1:A:70:ARG:HG2	1.44	1.09
1:B:70:ARG:HH11	1:B:70:ARG:HG2	1.10	1.08
1:A:70:ARG:HH11	1:A:70:ARG:CG	1.61	1.06
1:C:130:TYR:CE1	1:C:136:PRO:CD	2.38	1.05
1:B:70:ARG:HG2	1:B:70:ARG:NH1	1.63	1.04
1:A:208:GLU:OE1	1:A:219:HIS:HD2	1.42	1.01
1:D:34:HIS:H	1:D:62:ASP:HB2	1.26	1.00
1:B:208:GLU:OE1	1:B:219:HIS:HD2	1.42	0.99
1:C:88:GLN:NE2	1:C:94:ILE:CD1	2.28	0.96
1:C:88:GLN:HE22	1:C:94:ILE:CD1	1.78	0.96
1:B:117:VAL:HG12	1:B:117:VAL:O	1.67	0.94
1:A:152:GLU:HG3	1:A:155:HIS:CE1	2.01	0.94
1:C:229:VAL:CG1	1:C:230:ARG:N	2.31	0.93
1:D:117:VAL:HG12	1:D:117:VAL:O	1.68	0.93
1:C:88:GLN:HE22	1:C:94:ILE:HD11	1.34	0.92
1:A:229:VAL:CG1	1:A:230:ARG:N	2.30	0.92
1:B:52:VAL:HG22	1:B:98:GLU:HG3	1.51	0.91
1:B:229:VAL:CG1	1:B:230:ARG:N	2.29	0.91
1:C:146:PRO:O	1:C:223:THR:HG21	1.72	0.90
1:C:52:VAL:HG22	1:C:98:GLU:HG3	1.53	0.89
1:B:80:GLN:NE2	1:B:104:ARG:NH2	2.20	0.89
1:C:88:GLN:NE2	1:C:94:ILE:HD12	1.85	0.89
1:A:70:ARG:HG2	1:A:70:ARG:HH11	0.75	0.88
1:A:88:GLN:HE22	1:A:178:THR:HG23	1.37	0.87
1:A:140:PRO:HG3	1:A:213:ASN:HB2	1.57	0.87
1:D:114:GLU:HG2	1:D:120:ALA:HB2	1.55	0.87
1:D:131:GLU:C	1:D:134:LYS:N	2.33	0.87
1:A:77:SER:HB2	1:A:79:THR:HG23	1.57	0.86
1:A:229:VAL:HG12	1:A:230:ARG:N	1.87	0.86
1:A:152:GLU:HG3	1:A:155:HIS:ND1	1.90	0.86
1:B:208:GLU:OE1	1:B:219:HIS:CD2	2.29	0.86
1:C:229:VAL:HG13	1:C:230:ARG:N	1.91	0.86
1:D:52:VAL:HG22	1:D:98:GLU:HG3	1.58	0.85
1:A:139:PHE:CE2	1:A:216:GLY:HA3	2.09	0.85
1:B:80:GLN:HE21	1:B:104:ARG:NH2	1.75	0.84
1:B:156:THR:HG22	2:B:255:HOH:O	1.76	0.84
1:B:70:ARG:HH11	1:B:70:ARG:CB	1.89	0.84
1:B:229:VAL:HG12	1:B:230:ARG:N	1.91	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:VAL:HG12	1:A:230:ARG:HB2	1.61	0.82
1:B:96:ARG:HH11	1:B:177:GLN:HE21	1.28	0.82
1:D:88:GLN:HG3	1:D:89:PRO:HD2	1.62	0.82
1:A:208:GLU:OE1	1:A:219:HIS:CD2	2.32	0.81
1:C:166:ASN:HA	1:C:167:PRO:C	2.04	0.81
1:C:84:THR:HG23	2:C:247:HOH:O	1.80	0.81
1:D:117:VAL:O	1:D:117:VAL:CG1	2.30	0.80
1:D:230:ARG:HD2	1:D:230:ARG:C	2.05	0.80
1:B:86:LEU:C	1:B:87:GLU:OE2	2.24	0.80
1:A:77:SER:HG	1:A:80:GLN:HB2	1.45	0.79
1:A:151:ILE:HG13	1:A:157:VAL:HG13	1.63	0.79
1:C:181:ASP:OD1	1:C:183:SER:CB	2.29	0.79
1:D:114:GLU:HG2	1:D:120:ALA:CB	2.11	0.79
1:A:52:VAL:HG22	1:A:98:GLU:HG3	1.65	0.79
1:C:161:CYS:CB	1:C:209:CYS:SG	2.72	0.78
1:D:200:ARG:HB2	1:D:202:GLU:HG2	1.65	0.78
1:B:117:VAL:O	1:B:117:VAL:CG1	2.30	0.78
1:B:229:VAL:HG12	1:B:230:ARG:O	1.84	0.78
1:C:140:PRO:HG3	1:C:213:ASN:HB2	1.63	0.78
1:B:49:VAL:HG12	1:B:131:GLU:HA	1.65	0.78
1:C:164:ILE:HG13	1:C:164:ILE:O	1.84	0.78
1:C:166:ASN:CA	1:C:167:PRO:C	2.58	0.77
1:D:110:GLU:HB2	1:D:124:ASP:HB3	1.64	0.77
1:D:143:THR:HG22	1:D:144:GLN:N	1.99	0.77
1:C:161:CYS:HB2	1:C:209:CYS:SG	2.25	0.77
1:C:165:GLY:H	1:C:169:PRO:HD3	1.49	0.76
1:B:229:VAL:HG13	1:B:230:ARG:H	1.49	0.76
1:C:38:ILE:HD13	1:C:121:VAL:HG23	1.66	0.76
1:C:88:GLN:HE21	1:C:94:ILE:HD12	1.48	0.76
1:C:181:ASP:OD1	1:C:183:SER:HB3	1.85	0.76
1:C:38:ILE:HD13	1:C:121:VAL:CG2	2.16	0.75
1:D:230:ARG:C	1:D:230:ARG:CD	2.59	0.75
1:D:83:TYR:CD1	1:D:97:ILE:HG12	2.22	0.75
1:C:229:VAL:HG13	1:C:230:ARG:H	1.54	0.73
1:D:131:GLU:C	1:D:134:LYS:HA	2.13	0.73
1:C:165:GLY:N	1:C:169:PRO:HD3	2.04	0.72
1:C:130:TYR:CD1	1:C:136:PRO:HD3	2.24	0.72
1:C:229:VAL:HG13	1:C:230:ARG:HG3	1.72	0.72
1:D:142:ILE:HG22	1:D:142:ILE:O	1.89	0.71
1:D:83:TYR:HD1	1:D:97:ILE:HG12	1.52	0.71
1:C:104:ARG:HG2	2:C:253:HOH:O	1.88	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:131:GLU:C	1:D:134:LYS:CA	2.64	0.71
1:A:229:VAL:HG13	1:A:230:ARG:H	1.56	0.70
1:A:71:LYS:NZ	1:A:83:TYR:OH	2.25	0.70
1:B:96:ARG:NH1	1:B:177:GLN:HE21	1.88	0.70
1:C:128:THR:HG22	2:C:267:HOH:O	1.90	0.70
1:B:49:VAL:CG1	1:B:131:GLU:HA	2.22	0.70
1:B:186:ARG:NH2	1:B:203:ASP:OD2	2.25	0.69
1:A:152:GLU:CG	1:A:155:HIS:ND1	2.54	0.69
1:B:83:TYR:CD1	1:B:97:ILE:HG12	2.27	0.69
1:B:128:THR:HG22	2:B:10:HOH:O	1.91	0.69
1:A:88:GLN:HE22	1:A:178:THR:CG2	2.05	0.69
1:A:88:GLN:NE2	1:A:178:THR:HG23	2.08	0.68
1:C:37:GLU:OE1	1:C:60:ARG:NH2	2.26	0.68
1:C:88:GLN:NE2	1:C:94:ILE:HD11	2.00	0.68
1:A:229:VAL:CG1	1:A:230:ARG:H	2.07	0.68
1:B:80:GLN:HE21	1:B:104:ARG:HH21	1.42	0.68
1:C:176:ASN:O	1:C:177:GLN:HB2	1.94	0.68
1:A:158:LEU:HD12	1:A:193:PHE:HB3	1.77	0.67
1:B:36:PRO:CG	1:B:115:ASN:HB2	2.14	0.67
1:C:229:VAL:HG12	1:C:230:ARG:N	2.10	0.67
1:A:34:HIS:H	1:A:62:ASP:HB2	1.60	0.67
1:B:185:PRO:O	1:C:40:ARG:HD2	1.96	0.66
1:B:96:ARG:HH11	1:B:177:GLN:NE2	1.91	0.66
1:B:34:HIS:O	1:B:115:ASN:ND2	2.28	0.66
1:D:34:HIS:N	1:D:62:ASP:HB2	2.08	0.66
1:A:151:ILE:HG13	1:A:157:VAL:CG1	2.26	0.66
1:B:87:GLU:OE2	1:B:87:GLU:N	2.29	0.65
1:D:201:GLU:OE2	1:D:229:VAL:CG1	2.45	0.65
1:D:201:GLU:OE2	1:D:229:VAL:HG11	1.96	0.65
1:B:128:THR:CG2	2:B:10:HOH:O	2.44	0.65
1:D:174:ILE:HD12	1:D:210:VAL:HG21	1.79	0.65
1:A:190:LYS:NZ	1:D:43:GLN:HE22	1.96	0.64
1:C:100:VAL:CG1	1:C:129:ILE:HG12	2.27	0.64
1:C:130:TYR:HE1	1:C:136:PRO:HD3	1.59	0.64
1:A:96:ARG:NH1	1:A:98:GLU:OE1	2.30	0.64
1:A:198:ASN:HD21	1:D:60:ARG:NH1	1.94	0.64
1:D:224:ASN:ND2	1:D:226:TYR:OH	2.29	0.64
1:A:70:ARG:CD	1:A:75:LYS:HA	2.27	0.64
1:D:114:GLU:CG	1:D:120:ALA:HB2	2.27	0.64
1:A:101:ARG:HB3	1:A:104:ARG:HD3	1.80	0.64
1:C:136:PRO:HB2	1:C:215:MET:SD	2.37	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:VAL:HG13	1:A:230:ARG:N	2.13	0.64
1:C:217:THR:HB	2:C:260:HOH:O	1.97	0.64
1:D:88:GLN:HG3	1:D:89:PRO:CD	2.27	0.63
1:D:115:ASN:OD1	1:D:117:VAL:N	2.29	0.63
1:A:80:GLN:HG2	2:A:243:HOH:O	1.99	0.62
1:B:88:GLN:HB2	1:B:92:ILE:HB	1.81	0.62
1:B:149:ARG:HD2	2:B:254:HOH:O	2.00	0.62
1:A:229:VAL:HG12	1:A:230:ARG:CB	2.29	0.62
1:D:146:PRO:O	1:D:223:THR:HG21	2.00	0.62
1:A:32:ALA:HB1	1:A:62:ASP:HB3	1.82	0.61
1:A:70:ARG:HD2	1:A:75:LYS:HA	1.80	0.61
1:C:83:TYR:CD2	1:C:97:ILE:HG12	2.35	0.61
1:A:175:LYS:HD2	1:A:205:GLY:HA3	1.83	0.61
1:A:190:LYS:HZ3	1:D:43:GLN:HE22	1.48	0.61
1:C:83:TYR:HD2	1:C:97:ILE:HG12	1.65	0.61
1:A:43:GLN:HE22	1:D:190:LYS:NZ	1.98	0.60
1:A:43:GLN:NE2	1:D:190:LYS:NZ	2.49	0.60
1:A:70:ARG:HD2	1:A:75:LYS:CA	2.32	0.59
1:B:52:VAL:HG11	1:B:96:ARG:HH21	1.66	0.59
1:B:215:MET:N	1:B:215:MET:HE2	2.17	0.59
1:C:88:GLN:HB3	1:C:89:PRO:CD	2.33	0.59
1:C:181:ASP:OD1	1:C:183:SER:HB2	2.03	0.59
1:A:206:LYS:HG2	1:A:222:ALA:HB1	1.85	0.58
1:B:139:PHE:CE2	1:B:216:GLY:HA3	2.38	0.58
1:A:70:ARG:NH1	1:A:70:ARG:CG	2.30	0.58
1:A:221:LYS:HE3	2:A:251:HOH:O	2.04	0.58
1:D:119:ASP:OD2	1:D:119:ASP:N	2.30	0.58
1:C:130:TYR:CD1	1:C:136:PRO:CD	2.84	0.57
1:A:139:PHE:CD2	1:A:216:GLY:HA3	2.38	0.57
1:C:161:CYS:HB2	1:C:209:CYS:CB	2.35	0.57
1:C:166:ASN:N	1:C:167:PRO:O	2.37	0.57
1:A:222:ALA:O	1:A:223:THR:HG22	2.04	0.57
1:D:212:GLU:CG	1:D:213:ASN:N	2.68	0.57
1:A:140:PRO:CG	1:A:213:ASN:HB2	2.32	0.57
1:D:182:MET:HE2	1:D:189:LEU:CD1	2.35	0.57
1:A:222:ALA:C	1:A:223:THR:CG2	2.78	0.57
1:C:161:CYS:SG	1:C:209:CYS:CB	2.93	0.57
1:A:43:GLN:HE22	1:D:190:LYS:CE	2.19	0.56
1:D:212:GLU:HG2	1:D:213:ASN:N	2.20	0.56
1:C:186:ARG:NH2	1:C:203:ASP:OD2	2.38	0.56
1:C:229:VAL:HG13	1:C:230:ARG:CG	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:PRO:O	1:D:40:ARG:HD2	2.06	0.56
1:C:100:VAL:HG11	1:C:129:ILE:HG12	1.88	0.55
1:A:158:LEU:CD1	1:A:193:PHE:HB3	2.35	0.55
1:B:80:GLN:NE2	1:B:104:ARG:CZ	2.69	0.55
1:A:34:HIS:N	1:A:62:ASP:O	2.39	0.55
1:B:83:TYR:CE1	1:B:97:ILE:HG12	2.42	0.55
1:A:124:ASP:OD1	1:A:124:ASP:N	2.39	0.55
1:D:149:ARG:HD2	2:D:248:HOH:O	2.06	0.55
1:C:38:ILE:CD1	1:C:121:VAL:HG23	2.37	0.54
1:A:34:HIS:N	1:A:62:ASP:HB2	2.23	0.54
1:B:70:ARG:HG3	1:B:75:LYS:HA	1.87	0.54
1:A:153:VAL:HG23	1:A:228:LYS:O	2.08	0.54
1:B:96:ARG:NH1	1:B:177:GLN:HG2	2.23	0.54
1:A:190:LYS:CE	1:D:43:GLN:OE1	2.55	0.53
1:C:130:TYR:CE1	1:C:136:PRO:HD2	2.37	0.53
1:B:200:ARG:NH1	2:B:261:HOH:O	2.40	0.53
1:C:57:CYS:CB	1:C:111:CYS:SG	2.97	0.53
1:A:60:ARG:CG	1:A:60:ARG:HH11	2.21	0.53
1:A:40:ARG:HD2	1:D:185:PRO:O	2.08	0.53
1:A:75:LYS:HG3	1:A:76:VAL:N	2.23	0.53
1:C:165:GLY:C	1:C:167:PRO:O	2.52	0.52
1:A:138:GLY:O	1:A:213:ASN:ND2	2.41	0.52
1:C:138:GLY:HA3	1:C:166:ASN:O	2.08	0.52
1:C:88:GLN:CB	1:C:89:PRO:CD	2.87	0.52
1:C:130:TYR:CE1	1:C:136:PRO:CG	2.92	0.52
1:B:229:VAL:HG13	1:B:230:ARG:N	2.08	0.52
1:A:43:GLN:HE22	1:D:190:LYS:HE3	1.75	0.52
1:D:45:GLN:NE2	2:D:26:HOH:O	2.36	0.52
1:D:82:ARG:HG2	1:D:98:GLU:O	2.09	0.52
1:D:140:PRO:HG3	1:D:213:ASN:HB2	1.91	0.51
1:D:130:TYR:CD2	1:D:136:PRO:HD3	2.46	0.51
1:A:70:ARG:HD3	1:A:75:LYS:HA	1.93	0.51
1:B:70:ARG:HH11	1:B:70:ARG:HB3	1.73	0.51
1:B:80:GLN:HG2	1:B:83:TYR:CD2	2.45	0.51
1:D:115:ASN:OD1	1:D:117:VAL:HB	2.11	0.51
1:A:191:ASP:C	1:A:193:PHE:H	2.19	0.51
1:A:171:ILE:HG23	1:A:209:CYS:SG	2.51	0.50
1:B:82:ARG:O	1:B:97:ILE:HA	2.11	0.50
1:A:164:ILE:O	1:A:164:ILE:HG12	2.10	0.50
1:C:134:LYS:N	1:C:134:LYS:HD3	2.26	0.50
1:D:176:ASN:O	1:D:177:GLN:HB2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:LYS:HE2	1:C:43:GLN:OE1	2.12	0.50
1:C:130:TYR:HB3	1:C:135:THR:HG22	1.94	0.50
1:C:229:VAL:O	1:C:230:ARG:C	2.52	0.50
1:D:160:THR:HB	1:D:193:PHE:CZ	2.47	0.50
1:B:79:THR:HG22	1:B:80:GLN:N	2.26	0.50
1:B:82:ARG:NH1	1:B:105:ASP:OD2	2.41	0.50
1:C:49:VAL:HG12	1:C:130:TYR:O	2.11	0.50
1:D:144:GLN:OE1	1:D:162:LYS:HE3	2.12	0.50
1:B:80:GLN:O	1:B:80:GLN:HG3	2.03	0.50
1:C:130:TYR:CZ	1:C:136:PRO:CD	2.94	0.50
1:C:139:PHE:CE2	1:C:216:GLY:HA3	2.46	0.50
1:C:161:CYS:CB	1:C:209:CYS:CB	2.89	0.50
1:A:150:VAL:HG12	1:A:228:LYS:HE2	1.92	0.49
1:B:83:TYR:HD1	1:B:97:ILE:HG12	1.76	0.49
1:B:151:ILE:HG13	1:B:152:GLU:N	2.27	0.49
1:D:77:SER:O	1:D:78:GLY:C	2.54	0.49
1:D:187:TYR:OH	1:D:203:ASP:OD1	2.23	0.49
1:C:107:ALA:HB1	1:C:108:PRO:HD2	1.94	0.49
1:B:163:ALA:HB3	1:B:171:ILE:HD11	1.94	0.49
1:A:101:ARG:CB	1:A:104:ARG:HD3	2.42	0.49
1:D:200:ARG:CB	1:D:202:GLU:HG2	2.40	0.49
1:B:70:ARG:CG	1:B:70:ARG:NH1	2.30	0.49
1:C:142:ILE:HD12	1:C:218:GLU:HG3	1.94	0.49
1:B:206:LYS:HG2	1:B:222:ALA:HB1	1.93	0.49
1:C:166:ASN:OD1	1:C:167:PRO:HA	2.11	0.49
1:A:92:ILE:N	1:A:92:ILE:HD12	2.28	0.49
1:A:34:HIS:O	1:A:115:ASN:ND2	2.44	0.48
1:C:62:ASP:HA	1:C:63:PRO:C	2.38	0.48
1:C:38:ILE:O	1:C:38:ILE:HG22	2.12	0.48
1:D:173:TRP:CE2	1:D:194:LEU:HB2	2.48	0.48
1:D:128:THR:HG22	2:D:25:HOH:O	2.12	0.48
1:A:176:ASN:O	1:A:177:GLN:HB2	2.13	0.48
1:A:222:ALA:C	1:A:223:THR:HG23	2.38	0.48
1:C:161:CYS:SG	1:C:209:CYS:HB2	2.54	0.48
1:A:200:ARG:HB2	1:A:202:GLU:HG2	1.96	0.48
1:D:114:GLU:HG2	1:D:120:ALA:HB1	1.94	0.48
1:C:38:ILE:HD13	1:C:121:VAL:HG21	1.93	0.48
1:A:34:HIS:ND1	1:A:62:ASP:OD2	2.34	0.48
1:A:70:ARG:HD2	1:A:75:LYS:N	2.28	0.48
1:A:88:GLN:NE2	1:A:178:THR:CG2	2.73	0.48
1:D:212:GLU:CG	1:D:213:ASN:H	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:THR:O	1:C:128:THR:OG1	2.30	0.47
1:D:33:ALA:HA	1:D:62:ASP:HB2	1.96	0.47
1:D:142:ILE:O	1:D:143:THR:C	2.57	0.47
1:B:64:PRO:HA	1:B:65:PRO:HD3	1.72	0.47
1:C:75:LYS:HE2	1:C:75:LYS:HB2	1.15	0.47
1:C:166:ASN:CA	1:C:167:PRO:O	2.62	0.47
1:C:167:PRO:HD2	1:C:213:ASN:OD1	2.14	0.47
1:D:114:GLU:CD	1:D:120:ALA:HB2	2.40	0.47
1:D:115:ASN:OD1	1:D:115:ASN:C	2.58	0.46
1:D:187:TYR:CE1	1:D:196:ILE:HG23	2.50	0.46
1:A:43:GLN:NE2	1:D:190:LYS:HZ2	2.11	0.46
1:C:139:PHE:CD1	1:C:140:PRO:HD2	2.50	0.46
1:B:215:MET:HE3	1:B:215:MET:HB2	1.81	0.46
1:C:40:ARG:HB3	1:C:58:ALA:HB3	1.97	0.46
1:B:70:ARG:HH12	1:B:73:GLY:HA2	1.81	0.46
1:B:158:LEU:HD22	1:B:158:LEU:HA	1.74	0.46
1:D:190:LYS:HB3	1:D:190:LYS:HE2	1.78	0.46
1:B:170:ASN:HB2	2:B:249:HOH:O	2.14	0.46
1:C:130:TYR:CZ	1:C:136:PRO:CG	2.98	0.46
1:D:150:VAL:HB	1:D:226:TYR:HB2	1.98	0.45
1:D:182:MET:HE2	1:D:189:LEU:HD12	1.98	0.45
1:A:150:VAL:CG1	1:A:228:LYS:HE2	2.47	0.45
1:A:60:ARG:CG	1:A:60:ARG:NH1	2.76	0.45
1:A:163:ALA:HB3	1:A:171:ILE:HD11	1.98	0.45
1:B:151:ILE:HD11	1:B:155:HIS:HB3	1.99	0.45
1:B:151:ILE:HD13	1:B:157:VAL:HG22	1.98	0.45
1:D:54:SER:HA	1:D:95:LEU:O	2.17	0.45
1:A:47:VAL:CG1	1:A:127:LEU:HD11	2.46	0.45
1:A:67:ILE:HA	1:A:112:VAL:O	2.17	0.45
1:A:190:LYS:NZ	1:D:43:GLN:NE2	2.62	0.45
1:A:222:ALA:O	1:A:223:THR:CG2	2.64	0.45
1:B:63:PRO:O	1:B:115:ASN:OD1	2.35	0.45
1:B:131:GLU:HG3	1:B:132:GLY:N	2.32	0.45
1:C:96:ARG:HG2	2:C:247:HOH:O	2.17	0.45
1:D:160:THR:HB	1:D:193:PHE:CE2	2.51	0.45
1:D:76:VAL:O	1:D:76:VAL:CG2	2.62	0.44
1:A:190:LYS:NZ	1:D:43:GLN:OE1	2.51	0.44
1:B:222:ALA:C	1:B:223:THR:HG23	2.42	0.44
1:A:160:THR:HB	1:A:193:PHE:CE2	2.53	0.44
1:C:160:THR:HB	1:C:193:PHE:CZ	2.52	0.44
1:B:80:GLN:HE22	1:B:104:ARG:CZ	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:THR:HB	1:B:197:GLU:HA	1.99	0.44
1:C:151:ILE:HD13	1:C:157:VAL:HG22	2.00	0.44
1:A:189:LEU:HA	1:A:193:PHE:O	2.18	0.44
1:A:144:GLN:HB3	1:A:162:LYS:HG3	2.00	0.43
1:D:111:CYS:O	1:D:122:SER:HB2	2.18	0.43
1:B:176:ASN:O	1:B:177:GLN:HB2	2.18	0.43
1:D:64:PRO:HA	1:D:65:PRO:HD3	1.64	0.43
1:A:213:ASN:HB3	1:A:216:GLY:H	1.82	0.43
1:A:150:VAL:HG12	1:A:228:LYS:HG2	2.00	0.43
1:A:190:LYS:HZ3	1:D:43:GLN:NE2	2.16	0.43
1:D:48:ARG:HD3	1:D:218:GLU:OE1	2.19	0.43
1:C:229:VAL:CG1	1:C:230:ARG:HD3	2.48	0.43
1:D:130:TYR:CE2	1:D:136:PRO:HD3	2.53	0.43
1:C:130:TYR:CZ	1:C:136:PRO:HG2	2.54	0.43
1:B:47:VAL:HG22	1:B:48:ARG:N	2.34	0.42
1:C:57:CYS:HB2	1:C:111:CYS:SG	2.59	0.42
1:A:92:ILE:HD12	1:A:92:ILE:H	1.83	0.42
1:A:149:ARG:HE	1:A:149:ARG:HB3	1.75	0.42
1:A:193:PHE:N	1:A:193:PHE:CD1	2.87	0.42
1:B:113:ALA:HB3	1:B:121:VAL:HG23	2.00	0.42
1:B:134:LYS:HE2	1:B:134:LYS:HB2	1.67	0.42
1:C:200:ARG:HB2	1:C:202:GLU:HG2	2.01	0.42
1:C:57:CYS:SG	1:C:111:CYS:CB	3.05	0.42
1:C:88:GLN:HB3	1:C:89:PRO:HD2	2.00	0.42
1:B:222:ALA:C	1:B:223:THR:CG2	2.93	0.42
1:D:84:THR:HA	2:D:27:HOH:O	2.19	0.42
1:D:88:GLN:HB3	1:D:92:ILE:HB	2.01	0.42
1:B:96:ARG:HH11	1:B:177:GLN:HG2	1.85	0.42
1:C:150:VAL:HB	1:C:226:TYR:HB2	2.01	0.41
1:C:219:HIS:HE1	2:C:262:HOH:O	2.02	0.41
1:D:66:SER:O	1:D:113:ALA:HA	2.20	0.41
1:D:142:ILE:HG13	1:D:218:GLU:HG3	2.01	0.41
1:A:34:HIS:HA	1:A:35:PRO:HD3	1.75	0.41
1:B:206:LYS:HE3	1:B:222:ALA:CB	2.50	0.41
1:C:107:ALA:HB1	1:C:108:PRO:CD	2.51	0.41
1:C:181:ASP:CG	1:C:183:SER:HB3	2.45	0.41
1:D:144:GLN:HB3	1:D:162:LYS:HG3	2.02	0.41
1:D:197:GLU:O	1:D:198:ASN:C	2.63	0.41
1:A:62:ASP:HA	1:A:63:PRO:HA	1.87	0.41
1:C:156:THR:HG23	2:C:250:HOH:O	2.20	0.41
1:A:41:LYS:HA	1:A:42:PRO:HD3	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:168:THR:HA	1:C:169:PRO:HD3	1.75	0.41
1:D:33:ALA:HA	1:D:62:ASP:CB	2.50	0.41
1:D:193:PHE:HD1	1:D:193:PHE:N	2.19	0.41
1:A:88:GLN:HA	1:A:89:PRO:HD3	1.91	0.41
1:A:193:PHE:N	1:A:193:PHE:HD1	2.19	0.41
1:A:229:VAL:HG12	1:A:230:ARG:CA	2.49	0.41
1:B:92:ILE:HD11	1:C:185:PRO:HG3	2.02	0.41
1:B:155:HIS:O	1:B:199:SER:HB2	2.21	0.41
1:C:158:LEU:HA	1:C:158:LEU:HD22	1.50	0.41
1:C:186:ARG:HH22	1:C:203:ASP:CG	2.29	0.41
1:D:193:PHE:N	1:D:193:PHE:CD1	2.88	0.41
1:A:156:THR:HB	1:A:197:GLU:HA	2.02	0.41
1:B:165:GLY:HA3	1:B:169:PRO:HD3	2.03	0.40
1:C:64:PRO:HA	1:C:65:PRO:HD3	1.77	0.40
1:D:168:THR:HA	1:D:169:PRO:HD3	1.88	0.40
1:A:175:LYS:HD3	1:A:207:TYR:CZ	2.57	0.40
1:D:212:GLU:HG3	1:D:213:ASN:H	1.86	0.40
1:A:47:VAL:HG12	1:A:127:LEU:HD11	2.04	0.40
1:A:110:GLU:HB2	1:A:124:ASP:HB3	2.03	0.40
1:A:152:GLU:HG3	1:A:155:HIS:CG	2.54	0.40
1:B:150:VAL:HG22	1:B:226:TYR:HB2	2.03	0.40
1:C:98:GLU:HA	1:C:99:PRO:HA	1.89	0.40

There are no symmetry-related clashes.

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	197/210 (94%)	192 (98%)	5 (2%)	0	100	100
1	B	197/210 (94%)	193 (98%)	3 (2%)	1 (0%)	24	31
1	C	179/210 (85%)	175 (98%)	3 (2%)	1 (1%)	21	27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	188/210 (90%)	184 (98%)	2 (1%)	2 (1%)	11	13
All	All	761/840 (91%)	744 (98%)	13 (2%)	4 (0%)	24	31

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	143	THR
1	D	146	PRO
1	B	108	PRO
1	C	108	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	A	164/173 (95%)	130 (79%)	34 (21%)	1	1
1	B	164/173 (95%)	134 (82%)	30 (18%)	2	2
1	C	156/173 (90%)	120 (77%)	36 (23%)	1	1
1	D	161/173 (93%)	130 (81%)	31 (19%)	1	1
All	All	645/692 (93%)	514 (80%)	131 (20%)	1	1

All (131) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ILE
1	A	47	VAL
1	A	49	VAL
1	A	60	ARG
1	A	70	ARG
1	A	74	LYS
1	A	79	THR
1	A	80	GLN
1	A	81	SER
1	A	84	THR

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Mol	Chain	Res	Type
1	A	104	ARG
1	A	114	GLU
1	A	131	GLU
1	A	133	ASP
1	A	135	THR
1	A	148	THR
1	A	149	ARG
1	A	151	ILE
1	A	152	GLU
1	A	156	THR
1	A	157	VAL
1	A	158	LEU
1	A	160	THR
1	A	164	ILE
1	A	168	THR
1	A	171	ILE
1	A	183	SER
1	A	190	LYS
1	A	191	ASP
1	A	199	SER
1	A	202	GLU
1	A	214	SER
1	A	225	LEU
1	A	230	ARG
1	B	39	ILE
1	B	49	VAL
1	B	60	ARG
1	B	70	ARG
1	B	74	LYS
1	B	79	THR
1	B	80	GLN
1	B	84	THR
1	B	86	LEU
1	B	87	GLU
1	B	104	ARG
1	B	114	GLU
1	B	128	THR
1	B	133	ASP
1	B	134	LYS
1	B	148	THR
1	B	149	ARG
1	B	151	ILE

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Mol	Chain	Res	Type
1	B	152	GLU
1	B	156	THR
1	B	157	VAL
1	B	158	LEU
1	B	160	THR
1	B	164	ILE
1	B	166	ASN
1	B	174	ILE
1	B	187	TYR
1	B	199	SER
1	B	215	MET
1	B	225	LEU
1	C	37	GLU
1	C	38	ILE
1	C	49	VAL
1	C	60	ARG
1	C	62	ASP
1	C	74	LYS
1	C	75	LYS
1	C	76	VAL
1	C	84	THR
1	C	86	LEU
1	C	87	GLU
1	C	92	ILE
1	C	104	ARG
1	C	128	THR
1	C	131	GLU
1	C	134	LYS
1	C	148	THR
1	C	150	VAL
1	C	151	ILE
1	C	156	THR
1	C	157	VAL
1	C	158	LEU
1	C	159	MET
1	C	160	THR
1	C	161	CYS
1	C	170	ASN
1	C	174	ILE
1	C	179	LYS
1	C	189	LEU
1	C	201	GLU

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Mol	Chain	Res	Type
1	C	208	GLU
1	C	214	SER
1	C	215	MET
1	C	225	LEU
1	C	229	VAL
1	C	230	ARG
1	D	38	ILE
1	D	48	ARG
1	D	54	SER
1	D	70	ARG
1	D	76	VAL
1	D	82	ARG
1	D	84	THR
1	D	86	LEU
1	D	104	ARG
1	D	119	ASP
1	D	128	THR
1	D	131	GLU
1	D	135	THR
1	D	142	ILE
1	D	143	THR
1	D	148	THR
1	D	149	ARG
1	D	150	VAL
1	D	157	VAL
1	D	158	LEU
1	D	159	MET
1	D	160	THR
1	D	174	ILE
1	D	179	LYS
1	D	190	LYS
1	D	202	GLU
1	D	214	SER
1	D	215	MET
1	D	221	LYS
1	D	225	LEU
1	D	230	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN

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Mol	Chain	Res	Type
1	A	45	GLN
1	A	88	GLN
1	A	144	GLN
1	A	177	GLN
1	A	198	ASN
1	A	219	HIS
1	A	224	ASN
1	B	45	GLN
1	B	80	GLN
1	B	88	GLN
1	B	177	GLN
1	B	195	GLN
1	B	219	HIS
1	C	45	GLN
1	C	88	GLN
1	C	224	ASN
1	D	219	HIS
1	D	224	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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	199/210 (94%)	0.63	12 (6%) 27 29	40, 57, 100, 112	0
1	B	199/210 (94%)	0.37	10 (5%) 34 35	27, 47, 92, 157	0
1	C	187/210 (89%)	0.55	13 (6%) 22 24	32, 53, 88, 109	0
1	D	194/210 (92%)	0.78	16 (8%) 17 19	37, 58, 92, 124	0
All	All	779/840 (92%)	0.58	51 (6%) 25 27	27, 55, 93, 157	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	183	SER	5.2
1	A	32	ALA	4.8
1	D	117	VAL	4.1
1	C	124	ASP	4.1
1	D	78	GLY	4.0
1	C	161	CYS	3.8
1	C	88	GLN	3.7
1	C	62	ASP	3.6
1	D	92	ILE	3.5
1	A	33	ALA	3.3
1	B	77	SER	3.2
1	C	131	GLU	3.2
1	D	62	ASP	3.1
1	B	117	VAL	3.0
1	D	143	THR	3.0
1	B	118	GLY	2.9
1	D	91	GLY	2.9
1	B	116	GLY	2.8
1	D	116	GLY	2.7
1	B	131	GLU	2.7
1	B	132	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	191	ASP	2.6
1	D	165	GLY	2.6
1	C	137	ALA	2.6
1	D	77	SER	2.6
1	A	209	CYS	2.5
1	D	229	VAL	2.5
1	B	152	GLU	2.4
1	A	62	ASP	2.4
1	C	165	GLY	2.4
1	A	89	PRO	2.4
1	D	134	LYS	2.4
1	C	91	GLY	2.3
1	A	102	ALA	2.3
1	C	114	GLU	2.3
1	A	152	GLU	2.2
1	D	142	ILE	2.2
1	D	87	GLU	2.2
1	A	151	ILE	2.2
1	B	119	ASP	2.2
1	A	80	GLN	2.2
1	C	128	THR	2.2
1	C	89	PRO	2.1
1	A	161	CYS	2.1
1	C	168	THR	2.1
1	D	124	ASP	2.1
1	B	78	GLY	2.1
1	D	118	GLY	2.1
1	D	81	SER	2.1
1	B	191	ASP	2.0
1	A	66	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.