



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:06 PM UTC

PDB ID : 1IHM / pdb_00001ihm
Title : CRYSTAL STRUCTURE ANALYSIS OF NORWALK VIRUS CAPSID
Authors : Prasad, B.V.; Hardy, M.E.; Dokland, T.; Bella, J.; Rossmann, M.G.; Estes, M.K.
Deposited on : 2001-04-19
Resolution : 3.40 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

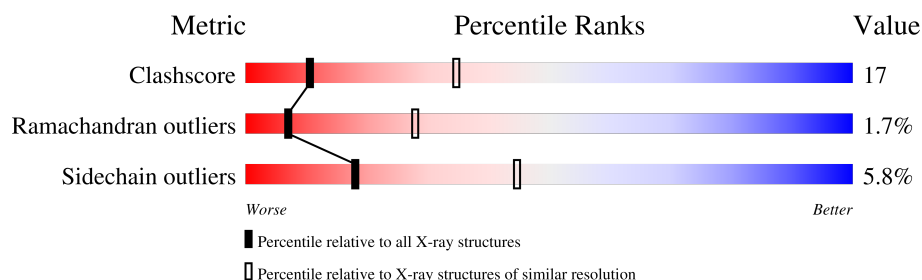
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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))
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	530	
1	B	530	
1	C	530	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11284 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	492	Total	C	N	O	S	0	0	0
			3721	2378	628	697	18			
1	B	511	Total	C	N	O	S	0	0	0
			3842	2449	649	726	18			
1	C	492	Total	C	N	O	S	0	0	0
			3721	2378	628	697	18			

There are 3 discrepancies between the modelled and reference sequences:

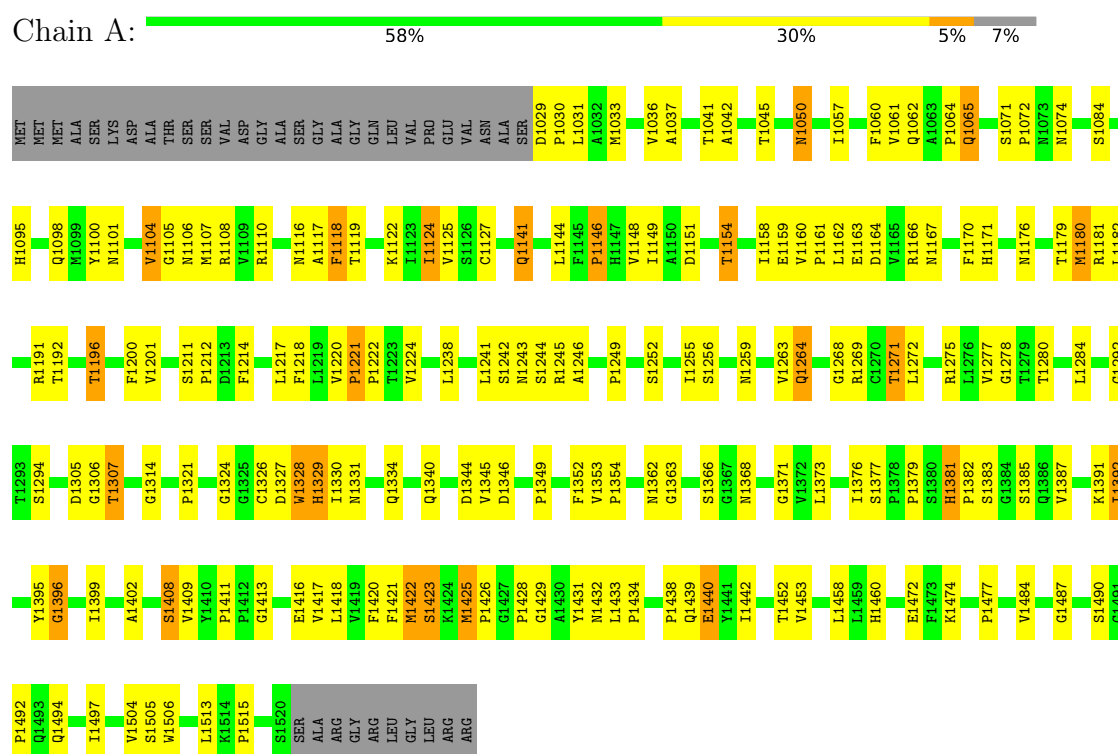
Chain	Residue	Modelled	Actual	Comment	Reference
A	1253	MET	ILE	SEE REMARK 999	UNP Q83884
B	2253	MET	ILE	SEE REMARK 999	UNP Q83884
C	3253	MET	ILE	SEE REMARK 999	UNP Q83884

3 Residue-property plots

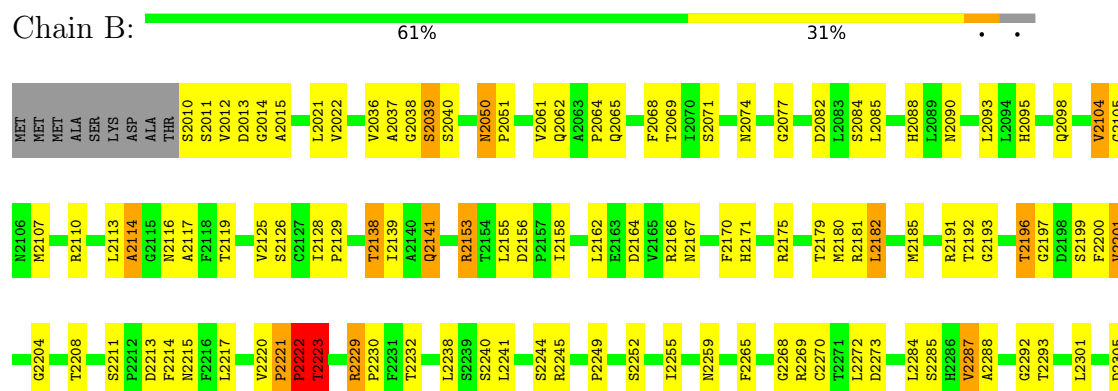
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

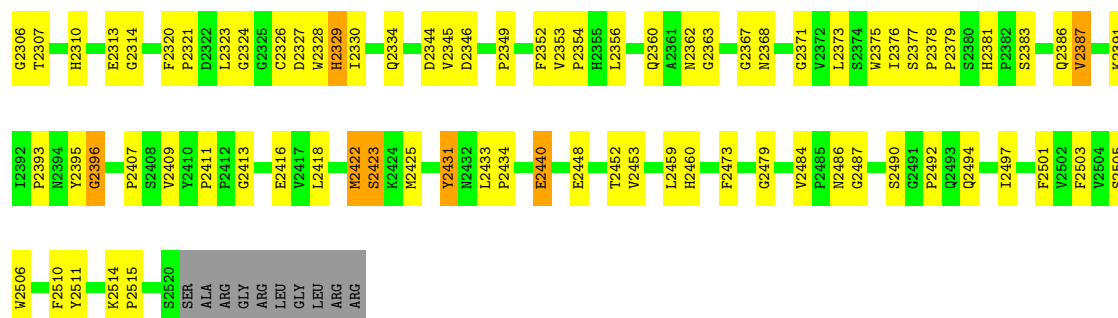
Note EDS was not executed.

- Molecule 1: capsid protein



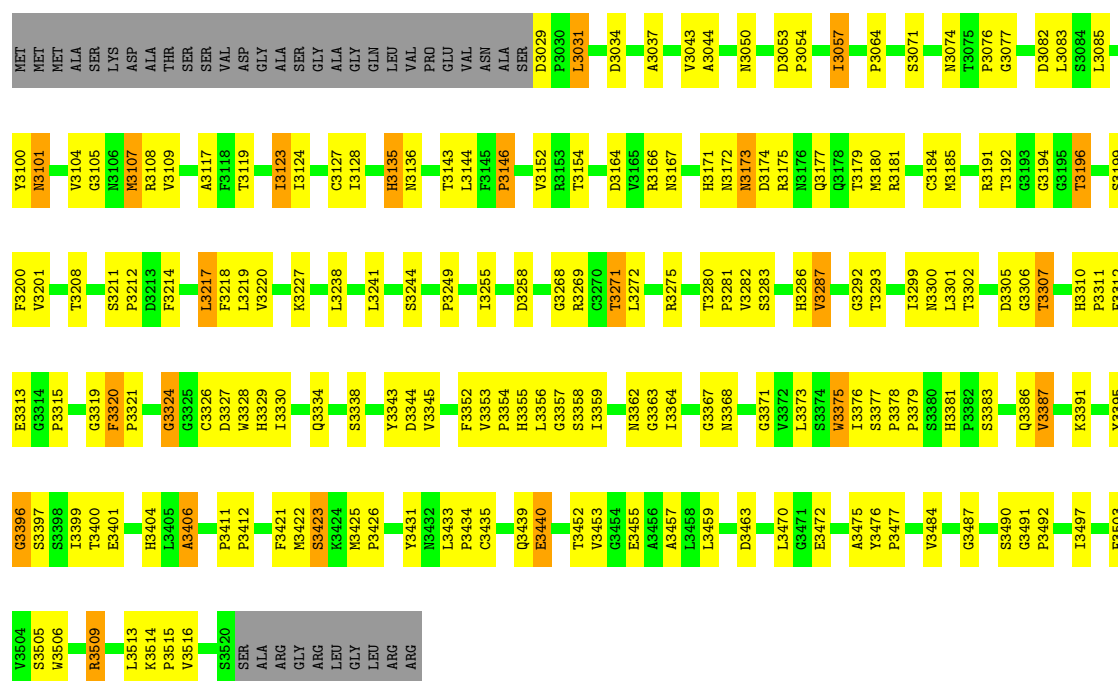
- Molecule 1: capsid protein





• Molecule 1: capsid protein

Chain C: 58% 31% 7%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	605.74Å 605.74Å 466.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.00 – 3.40	Depositor
% Data completeness (in resolution range)	(Not available) (35.00-3.40)	Depositor
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.260 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	11284	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

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	0.50	0/3835	1.04	21/5261 (0.4%)
1	B	0.49	0/3957	1.04	28/5428 (0.5%)
1	C	0.48	0/3835	1.05	24/5261 (0.5%)
All	All	0.49	0/11627	1.04	73/15950 (0.5%)

There are no bond length outliers.

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2221	PRO	CA-C-N	9.47	131.68	119.84
1	B	2221	PRO	C-N-CA	9.47	131.68	119.84
1	A	1306	GLY	N-CA-C	-7.90	104.18	114.85
1	A	1167	ASN	N-CA-C	-7.22	104.72	113.97
1	A	1217	LEU	N-CA-C	7.02	120.07	108.48
1	B	2386	GLN	N-CA-C	6.95	119.53	110.43
1	C	3175	ARG	N-CA-C	-6.80	105.01	113.38
1	B	2199	SER	N-CA-C	-6.80	104.85	113.01
1	C	3381	HIS	N-CA-C	6.68	122.79	113.57
1	B	2065	GLN	N-CA-C	-6.60	102.51	111.55
1	B	2510	PHE	N-CA-C	6.43	120.13	112.93
1	C	3117	ALA	N-CA-C	-6.40	105.47	113.28
1	C	3034	ASP	N-CA-C	-6.32	102.11	110.07
1	A	1117	ALA	N-CA-C	-6.25	105.66	113.28
1	C	3217	LEU	N-CA-C	6.14	123.89	110.80
1	C	3146	PRO	N-CA-C	-6.14	101.57	111.03
1	A	1345	VAL	N-CA-C	6.08	116.62	108.11
1	A	1221	PRO	CA-C-N	5.93	125.94	119.90
1	A	1221	PRO	C-N-CA	5.93	125.94	119.90
1	C	3301	LEU	N-CA-C	5.93	118.35	110.53
1	B	2381	HIS	N-CA-C	5.91	121.66	113.45
1	C	3476	TYR	CA-C-N	5.89	127.21	119.84
1	C	3476	TYR	C-N-CA	5.89	127.21	119.84
1	B	2193	GLY	N-CA-C	-5.83	104.26	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2314	GLY	CA-C-N	5.80	125.34	119.19
1	B	2314	GLY	C-N-CA	5.80	125.34	119.19
1	C	3167	ASN	N-CA-C	-5.80	106.36	113.50
1	C	3491	GLY	CA-C-N	5.77	125.45	119.56
1	C	3491	GLY	C-N-CA	5.77	125.45	119.56
1	A	1065	GLN	N-CA-C	-5.75	103.40	110.65
1	A	1176	ASN	N-CA-C	-5.72	106.34	113.38
1	C	3044	ALA	N-CA-C	5.72	117.51	111.28
1	B	2167	ASN	N-CA-C	-5.68	106.48	113.41
1	B	2306	GLY	N-CA-C	-5.66	107.86	115.21
1	B	2197	GLY	N-CA-C	-5.64	99.82	113.18
1	B	2287	VAL	N-CA-C	5.55	116.22	108.89
1	C	3306	GLY	N-CA-C	-5.55	107.60	115.43
1	A	1146	PRO	N-CA-C	-5.50	102.56	111.03
1	A	1328	TRP	N-CA-C	5.49	119.05	110.32
1	A	1314	GLY	CA-C-N	5.44	124.95	119.19
1	A	1314	GLY	C-N-CA	5.44	124.95	119.19
1	B	2201	VAL	N-CA-C	-5.39	100.08	108.81
1	C	3406	ALA	CA-C-N	5.39	125.39	119.89
1	C	3406	ALA	C-N-CA	5.39	125.39	119.89
1	B	2422	MET	N-CA-C	5.36	118.14	109.40
1	B	2345	VAL	N-CA-C	5.35	115.59	108.11
1	B	2222	PRO	N-CA-C	5.33	123.45	112.47
1	C	3386	GLN	N-CA-C	5.32	117.40	110.43
1	C	3199	SER	N-CA-C	-5.30	106.31	112.89
1	C	3128	ILE	CA-C-N	5.29	123.47	119.66
1	C	3128	ILE	C-N-CA	5.29	123.47	119.66
1	A	1422	MET	N-CA-C	5.27	118.00	109.40
1	B	2050	ASN	CA-C-N	5.21	125.12	119.85
1	B	2050	ASN	C-N-CA	5.21	125.12	119.85
1	A	1050	ASN	CA-C-N	5.21	125.42	120.31
1	A	1050	ASN	C-N-CA	5.21	125.42	120.31
1	C	3320	PHE	CA-C-N	5.21	125.21	119.90
1	C	3320	PHE	C-N-CA	5.21	125.21	119.90
1	B	2211	SER	N-CA-C	-5.20	102.64	110.24
1	B	2039	SER	N-CA-C	5.17	117.59	111.33
1	B	2217	LEU	N-CA-C	5.17	121.81	110.80
1	B	2301	LEU	N-CA-C	5.16	118.12	110.48
1	C	3287	VAL	N-CA-C	5.16	116.00	108.58
1	A	1381	HIS	N-CA-C	5.15	120.67	113.57
1	B	2229	ARG	N-CA-C	5.13	116.87	109.48
1	A	1402	ALA	N-CA-C	5.12	117.03	108.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1211	SER	N-CA-C	-5.10	103.04	110.08
1	A	1104	VAL	CB-CA-C	-5.08	103.97	111.80
1	C	3172	ASN	N-CA-C	-5.07	101.43	109.59
1	B	2156	ASP	CA-C-N	5.06	124.97	119.76
1	B	2156	ASP	C-N-CA	5.06	124.97	119.76
1	A	1154	THR	N-CA-C	5.04	117.78	110.42
1	B	2138	THR	N-CA-C	-5.00	103.12	110.52

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	3721	0	3624	136	0
1	B	3842	0	3737	129	0
1	C	3721	0	3624	129	0
All	All	11284	0	10985	373	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 17.

All (373) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2330:ILE:CG2	1:B:2373:LEU:HD23	1.44	1.46
1:A:1330:ILE:CG2	1:A:1373:LEU:HD23	1.47	1.42
1:B:2321:PRO:CB	1:B:2373:LEU:HD11	1.59	1.30
1:C:3321:PRO:CB	1:C:3373:LEU:HD11	1.72	1.20
1:B:2330:ILE:HG22	1:B:2373:LEU:CD2	1.72	1.20
1:B:2321:PRO:HB3	1:B:2373:LEU:CD1	1.72	1.19
1:C:3330:ILE:HG22	1:C:3373:LEU:HD23	1.21	1.17
1:C:3330:ILE:CG2	1:C:3373:LEU:HD23	1.77	1.13
1:A:1330:ILE:HG22	1:A:1373:LEU:CD2	1.78	1.12
1:B:2321:PRO:HB3	1:B:2373:LEU:HD11	1.20	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1321:PRO:CB	1:A:1373:LEU:HD11	1.84	1.07
1:C:3321:PRO:HB3	1:C:3373:LEU:CD1	1.86	1.04
1:A:1330:ILE:CG2	1:A:1373:LEU:CD2	2.35	1.00
1:B:2330:ILE:CG2	1:B:2373:LEU:CD2	2.33	1.00
1:A:1321:PRO:HB3	1:A:1373:LEU:HD11	1.41	0.99
1:C:3123:ILE:HD11	1:C:3184:CYS:SG	2.02	0.97
1:A:1321:PRO:HB3	1:A:1373:LEU:CD1	1.95	0.97
1:B:2329:HIS:HB3	1:B:2344:ASP:HA	1.50	0.93
1:C:3330:ILE:HD12	1:C:3371:GLY:HA3	1.51	0.91
1:A:1329:HIS:HB3	1:A:1344:ASP:HA	1.49	0.91
1:B:2321:PRO:CB	1:B:2373:LEU:CD1	2.41	0.91
1:C:3321:PRO:CB	1:C:3373:LEU:CD1	2.46	0.90
1:C:3387:VAL:HG21	1:C:3431:TYR:HE2	1.36	0.90
1:A:1326:CYS:SG	1:A:1376:ILE:HB	2.15	0.86
1:B:2138:THR:H	1:B:2141:GLN:HG3	1.39	0.86
1:C:3329:HIS:HB3	1:C:3344:ASP:HA	1.58	0.85
1:B:2422:MET:HE3	1:B:2434:PRO:HB3	1.59	0.84
1:C:3487:GLY:H	1:C:3490:SER:HB3	1.42	0.84
1:B:2321:PRO:HB3	1:B:2373:LEU:HD12	1.60	0.83
1:A:1330:ILE:HG22	1:A:1373:LEU:HD23	0.82	0.82
1:B:2119:THR:HG22	1:B:2192:THR:HG22	1.64	0.80
1:C:3321:PRO:HB3	1:C:3373:LEU:HD12	1.60	0.80
1:B:2272:LEU:HB2	1:B:2452:THR:HG21	1.64	0.80
1:B:2326:CYS:SG	1:B:2376:ILE:HB	2.22	0.79
1:A:1327:ASP:HB2	1:A:1379:PRO:HA	1.64	0.78
1:C:3321:PRO:CG	1:C:3373:LEU:HD11	2.14	0.78
1:B:2038:GLY:HA3	1:B:2213:ASP:HB3	1.64	0.78
1:C:3321:PRO:HB3	1:C:3373:LEU:HD11	1.45	0.76
1:C:3326:CYS:SG	1:C:3376:ILE:HB	2.26	0.76
1:C:3328:TRP:HB3	1:C:3376:ILE:HG22	1.67	0.75
1:A:1272:LEU:HB2	1:A:1452:THR:HG21	1.68	0.75
1:B:2321:PRO:CG	1:B:2373:LEU:HD11	2.16	0.75
1:B:2330:ILE:HG22	1:B:2373:LEU:HD23	0.76	0.75
1:C:3174:ASP:HB2	1:C:3177:GLN:HG2	1.68	0.74
1:A:1249:PRO:O	1:A:1423:SER:HB3	1.87	0.74
1:B:2061:VAL:HG11	1:C:3144:LEU:HD23	1.70	0.74
1:C:3387:VAL:HG21	1:C:3431:TYR:CE2	2.21	0.74
1:B:2321:PRO:HB2	1:B:2373:LEU:HD11	1.67	0.73
1:B:2327:ASP:HB2	1:B:2379:PRO:HA	1.71	0.72
1:B:2249:PRO:O	1:B:2423:SER:HB3	1.89	0.72
1:C:3064:PRO:HB2	1:C:3505:SER:HB3	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3127:CYS:HB2	1:C:3146:PRO:HG2	1.72	0.71
1:B:2129:PRO:HA	1:B:2180:MET:HE2	1.72	0.71
1:C:3472:GLU:OE2	1:C:3514:LYS:HG3	1.91	0.70
1:A:1141:GLN:HA	1:A:1144:LEU:HD12	1.74	0.69
1:A:1425:MET:HE3	1:A:1426:PRO:HD2	1.75	0.69
1:C:3127:CYS:HB3	1:C:3180:MET:HE3	1.75	0.68
1:C:3421:PHE:O	1:C:3435:CYS:HB3	1.94	0.68
1:C:3311:PRO:HB2	1:C:3401:GLU:HG3	1.76	0.68
1:B:2255:ILE:HD11	1:B:2492:PRO:HD2	1.76	0.68
1:B:2330:ILE:HG21	1:B:2373:LEU:HD23	1.70	0.68
1:C:3321:PRO:HB2	1:C:3373:LEU:HD11	1.70	0.68
1:A:1399:ILE:HB	1:C:3399:ILE:HG21	1.75	0.67
1:C:3315:PRO:HA	1:C:3355:HIS:HD2	1.58	0.67
1:C:3329:HIS:CB	1:C:3344:ASP:HA	2.24	0.67
1:A:1353:VAL:HG13	1:A:1396:GLY:HA3	1.76	0.67
1:C:3330:ILE:HG22	1:C:3373:LEU:CD2	2.14	0.67
1:C:3287:VAL:HG21	1:C:3439:GLN:OE1	1.95	0.67
1:B:2119:THR:HG21	1:B:2200:PHE:CE2	2.30	0.66
1:A:1387:VAL:HG21	1:A:1431:TYR:HE2	1.60	0.66
1:C:3330:ILE:CG2	1:C:3373:LEU:CD2	2.67	0.66
1:B:2330:ILE:HD12	1:B:2371:GLY:HA3	1.78	0.66
1:A:1116:ASN:ND2	1:A:1200:PHE:CE1	2.65	0.65
1:C:3310:HIS:O	1:C:3313:GLU:HG2	1.96	0.65
1:C:3173:ASN:O	1:C:3227:LYS:HD3	1.97	0.65
1:A:1255:ILE:HG23	1:A:1417:VAL:HG11	1.78	0.65
1:B:2116:ASN:HD22	1:B:2200:PHE:HE1	1.45	0.64
1:C:3241:LEU:HD22	1:C:3440:GLU:HG3	1.79	0.64
1:B:2129:PRO:HA	1:B:2180:MET:CE	2.27	0.64
1:B:2328:TRP:HB3	1:B:2376:ILE:HG22	1.79	0.64
1:C:3433:LEU:HD12	1:C:3434:PRO:HD2	1.79	0.64
1:A:1321:PRO:HB3	1:A:1373:LEU:HD12	1.79	0.63
1:C:3249:PRO:O	1:C:3423:SER:HB3	1.98	0.63
1:C:3330:ILE:HG23	1:C:3373:LEU:HD23	1.77	0.62
1:C:3101:ASN:HB3	1:C:3220:VAL:HG21	1.81	0.62
1:C:3327:ASP:HB2	1:C:3379:PRO:HA	1.83	0.61
1:A:1330:ILE:HD12	1:A:1371:GLY:HA3	1.82	0.61
1:C:3272:LEU:HB2	1:C:3452:THR:HG21	1.82	0.60
1:B:2245:ARG:HB2	1:B:2433:LEU:HD21	1.84	0.60
1:C:3100:TYR:HA	1:C:3218:PHE:O	2.01	0.60
1:A:1245:ARG:HB2	1:A:1433:LEU:HD21	1.84	0.60
1:B:2269:ARG:HH22	1:B:2409:VAL:HG13	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1148:VAL:HG21	1:A:1160:VAL:HG22	1.84	0.59
1:B:2484:VAL:O	1:B:2515:PRO:HA	2.02	0.59
1:B:2387:VAL:HG21	1:B:2431:TYR:CE2	2.38	0.59
1:B:2105:GLY:HA3	1:B:2214:PHE:HA	1.85	0.59
1:B:2071:SER:H	1:B:2074:ASN:HD22	1.51	0.59
1:A:1241:LEU:HD22	1:A:1440:GLU:HG3	1.84	0.59
1:C:3071:SER:H	1:C:3074:ASN:ND2	2.00	0.59
1:A:1119:THR:HG21	1:A:1200:PHE:HE2	1.67	0.59
1:C:3343:TYR:CD1	1:C:3364:ILE:HD11	2.38	0.58
1:B:2192:THR:OG1	1:B:2196:THR:HG23	2.03	0.58
1:A:1057:ILE:HD12	1:A:1212:PRO:HA	1.85	0.58
1:A:1379:PRO:HB2	1:A:1382:PRO:HG3	1.86	0.58
1:A:1321:PRO:CB	1:A:1373:LEU:CD1	2.64	0.57
1:B:2162:LEU:HD13	1:B:2182:LEU:HG	1.86	0.57
1:B:2200:PHE:CE1	1:C:3191:ARG:HB3	2.39	0.57
1:B:2353:VAL:HG13	1:B:2396:GLY:HA3	1.84	0.57
1:C:3105:GLY:O	1:C:3164:ASP:HB2	2.04	0.57
1:A:1411:PRO:HD3	1:A:1418:LEU:HG	1.86	0.57
1:A:1071:SER:HB3	1:A:1201:VAL:HG12	1.87	0.57
1:A:1037:ALA:O	1:A:1166:ARG:NH2	2.38	0.57
1:C:3354:PRO:HD2	1:C:3396:GLY:H	1.69	0.57
1:C:3272:LEU:HB2	1:C:3452:THR:CG2	2.35	0.56
1:A:1122:LYS:NZ	1:A:1151:ASP:HB2	2.20	0.56
1:B:2265:PHE:O	1:B:2269:ARG:HD2	2.04	0.56
1:C:3425:MET:HG3	1:C:3426:PRO:HD2	1.87	0.56
1:A:1255:ILE:HD11	1:A:1492:PRO:HD2	1.87	0.56
1:A:1060:PHE:HZ	1:A:1108:ARG:HD3	1.70	0.56
1:A:1164:ASP:OD1	1:A:1171:HIS:HE1	1.89	0.56
1:A:1242:SER:HB3	1:B:2285:SER:HB3	1.87	0.56
1:A:1098:GLN:O	1:A:1222:PRO:HA	2.05	0.56
1:B:2061:VAL:HG11	1:C:3144:LEU:CD2	2.36	0.56
1:A:1328:TRP:HB3	1:A:1376:ILE:HG22	1.87	0.56
1:B:2098:GLN:O	1:B:2222:PRO:HA	2.06	0.56
1:A:1122:LYS:HD2	1:A:1151:ASP:HA	1.89	0.55
1:C:3459:LEU:HD11	1:C:3475:ALA:HB2	1.88	0.55
1:B:2113:LEU:O	1:B:2114:ALA:HB2	2.07	0.55
1:B:2334:GLN:HG3	1:B:2368:ASN:O	2.06	0.55
1:C:3321:PRO:HG3	1:C:3373:LEU:HD11	1.85	0.55
1:A:1221:PRO:HB3	1:C:3144:LEU:HD12	1.89	0.54
1:A:1422:MET:HE2	1:A:1434:PRO:HB3	1.89	0.54
1:B:2411:PRO:HD3	1:B:2418:LEU:HG	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3353:VAL:HG11	1:C:3356:LEU:HD12	1.88	0.54
1:A:1071:SER:H	1:A:1074:ASN:HD22	1.55	0.54
1:A:1125:VAL:HG12	1:A:1182:LEU:HD22	1.87	0.54
1:A:1105:GLY:HA3	1:A:1214:PHE:HA	1.90	0.54
1:C:3271:THR:OG1	1:C:3275:ARG:HB2	2.06	0.54
1:A:1106:ASN:HD22	1:A:1163:GLU:HA	1.72	0.54
1:B:2269:ARG:NH2	1:B:2409:VAL:HG13	2.22	0.54
1:A:1095:HIS:O	1:B:2095:HIS:HE1	1.91	0.54
1:A:1321:PRO:CG	1:A:1373:LEU:HD11	2.37	0.54
1:A:1321:PRO:HB2	1:A:1373:LEU:HD11	1.85	0.53
1:A:1244:SER:HB2	1:A:1268:GLY:CA	2.39	0.53
1:C:3292:GLY:O	1:C:3368:ASN:HA	2.08	0.53
1:A:1065:GLN:HG2	1:A:1506:TRP:HB2	1.90	0.53
1:B:2288:ALA:HB1	1:B:2373:LEU:HD12	1.91	0.53
1:C:3484:VAL:O	1:C:3515:PRO:HA	2.08	0.53
1:A:1460:HIS:ND1	1:A:1504:VAL:HG21	2.23	0.53
1:C:3211:SER:HB2	1:C:3212:PRO:HD2	1.91	0.53
1:C:3330:ILE:HG22	1:C:3373:LEU:HA	1.91	0.53
1:C:3352:PHE:C	1:C:3354:PRO:HD3	2.33	0.52
1:C:3334:GLN:HG3	1:C:3368:ASN:O	2.08	0.52
1:B:2164:ASP:OD2	1:B:2171:HIS:HE1	1.93	0.52
1:B:2238:LEU:HB3	1:B:2497:ILE:HG22	1.90	0.52
1:A:1050:ASN:HD22	1:A:1100:TYR:HE2	1.58	0.52
1:B:2012:VAL:HG12	1:C:3154:THR:HG21	1.91	0.52
1:C:3037:ALA:O	1:C:3166:ARG:NH2	2.42	0.52
1:C:3064:PRO:HB2	1:C:3505:SER:CB	2.38	0.52
1:C:3085:LEU:HB3	1:C:3179:THR:HB	1.90	0.52
1:B:2514:LYS:NZ	1:C:3076:PRO:HD2	2.25	0.52
1:A:1104:VAL:HG22	1:A:1170:PHE:HB2	1.90	0.52
1:B:2490:SER:HA	1:B:2494:GLN:NE2	2.25	0.52
1:B:2232:THR:OG1	1:B:2448:GLU:OE1	2.27	0.51
1:C:3326:CYS:HB3	1:C:3328:TRP:HD1	1.74	0.51
1:A:1036:VAL:HB	1:A:1106:ASN:HD21	1.75	0.51
1:A:1116:ASN:HB3	1:A:1118:PHE:H	1.75	0.51
1:A:1399:ILE:HB	1:C:3399:ILE:CG2	2.40	0.51
1:B:2330:ILE:HG22	1:B:2373:LEU:HA	1.92	0.51
1:B:2068:PHE:CZ	1:B:2204:GLY:HA3	2.46	0.51
1:C:3293:THR:HA	1:C:3367:GLY:O	2.11	0.51
1:B:2433:LEU:HD12	1:B:2434:PRO:HD2	1.91	0.51
1:A:1269:ARG:HH22	1:A:1409:VAL:HG13	1.76	0.51
1:B:2129:PRO:CA	1:B:2180:MET:HE2	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1305:ASP:HB2	1:A:1307:THR:OG1	2.11	0.50
1:B:2153:ARG:HH11	1:B:2153:ARG:HB2	1.76	0.50
1:B:2413:GLY:HA3	1:B:2416:GLU:CD	2.36	0.50
1:A:1460:HIS:CD2	1:A:1472:GLU:HG3	2.46	0.50
1:A:1119:THR:HG21	1:A:1200:PHE:CE2	2.45	0.50
1:A:1330:ILE:HG22	1:A:1373:LEU:HA	1.93	0.50
1:A:1387:VAL:HG21	1:A:1431:TYR:CE2	2.45	0.50
1:C:3378:PRO:HA	1:C:3387:VAL:HG22	1.93	0.50
1:B:2191:ARG:HH11	1:B:2191:ARG:HG2	1.77	0.50
1:B:2241:LEU:HD22	1:B:2440:GLU:HG3	1.92	0.50
1:A:1243:ASN:HB3	1:A:1246:ALA:O	2.11	0.50
1:A:1064:PRO:O	1:A:1505:SER:HB2	2.11	0.50
1:A:1124:ILE:HG12	1:A:1149:ILE:HG12	1.94	0.50
1:A:1242:SER:CB	1:B:2285:SER:HB3	2.42	0.50
1:B:2200:PHE:CD1	1:C:3191:ARG:HB3	2.47	0.49
1:B:2353:VAL:HG11	1:B:2356:LEU:HD12	1.93	0.49
1:C:3123:ILE:CD1	1:C:3184:CYS:SG	2.90	0.49
1:A:1118:PHE:N	1:A:1118:PHE:CD1	2.81	0.49
1:B:2422:MET:HG2	1:B:2434:PRO:HA	1.94	0.49
1:A:1029:ASP:HB3	1:A:1030:PRO:HD2	1.94	0.49
1:A:1484:VAL:O	1:A:1515:PRO:HA	2.12	0.49
1:C:3119:THR:HG22	1:C:3192:THR:HG22	1.95	0.49
1:A:1346:ASP:OD1	1:A:1349:PRO:HD3	2.11	0.49
1:C:3310:HIS:HB3	1:C:3312:PHE:CE2	2.48	0.49
1:B:2125:VAL:CG1	1:B:2182:LEU:HD22	2.43	0.49
1:C:3238:LEU:HB3	1:C:3497:ILE:HG22	1.95	0.49
1:B:2062:GLN:H	1:C:3143:THR:HG21	1.78	0.49
1:C:3399:ILE:HD12	1:C:3399:ILE:H	1.78	0.49
1:A:1110:ARG:HA	1:A:1158:ILE:O	2.13	0.48
1:A:1334:GLN:HG3	1:A:1368:ASN:O	2.13	0.48
1:B:2125:VAL:HG12	1:B:2182:LEU:HD22	1.95	0.48
1:C:3377:SER:OG	1:C:3378:PRO:HD2	2.13	0.48
1:C:3105:GLY:HA3	1:C:3214:PHE:HA	1.94	0.48
1:C:3164:ASP:OD1	1:C:3171:HIS:HE1	1.97	0.48
1:C:3123:ILE:HG12	1:C:3124:ILE:N	2.28	0.48
1:C:3459:LEU:HD23	1:C:3503:PHE:HA	1.94	0.48
1:A:1105:GLY:O	1:A:1164:ASP:HB2	2.14	0.48
1:A:1118:PHE:N	1:A:1118:PHE:HD1	2.12	0.48
1:C:3315:PRO:HD3	1:C:3404:HIS:O	2.13	0.48
1:B:2085:LEU:HB3	1:B:2179:THR:HB	1.96	0.48
1:B:2321:PRO:HG3	1:B:2373:LEU:HD11	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1352:PHE:C	1:A:1354:PRO:HD3	2.38	0.48
1:C:3050:ASN:ND2	1:C:3217:LEU:O	2.46	0.48
1:C:3299:ILE:HB	1:C:3359:ILE:HG13	1.95	0.48
1:A:1329:HIS:HE1	1:A:1377:SER:HB2	1.78	0.48
1:A:1324:GLY:HA3	1:A:1391:LYS:C	2.38	0.47
1:C:3375:TRP:CD1	1:C:3375:TRP:C	2.92	0.47
1:A:1221:PRO:HA	1:A:1222:PRO:HD3	1.78	0.47
1:A:1429:GLY:N	1:A:1431:TYR:HE1	2.13	0.47
1:B:2514:LYS:HZ2	1:C:3076:PRO:HD2	1.79	0.47
1:C:3457:ALA:HB3	1:C:3475:ALA:HB3	1.95	0.47
1:A:1122:LYS:HZ3	1:A:1151:ASP:HB2	1.78	0.47
1:A:1420:PHE:HB2	1:A:1422:MET:HE3	1.96	0.47
1:B:2221:PRO:HA	1:B:2222:PRO:HD2	1.57	0.47
1:B:2293:THR:HG22	1:B:2368:ASN:HB3	1.96	0.47
1:C:3324:GLY:HA3	1:C:3391:LYS:C	2.40	0.47
1:B:2229:ARG:HA	1:B:2230:PRO:HD3	1.74	0.47
1:C:3071:SER:HA	1:C:3201:VAL:HG12	1.97	0.47
1:C:3082:ASP:O	1:C:3083:LEU:HD23	2.15	0.46
1:A:1238:LEU:HB3	1:A:1497:ILE:HG22	1.97	0.46
1:A:1340:GLN:HB2	1:B:2375:TRP:CE3	2.49	0.46
1:A:1162:LEU:HD13	1:A:1182:LEU:HG	1.98	0.46
1:B:2077:GLY:O	1:B:2185:MET:HE2	2.15	0.46
1:A:1127:CYS:HB3	1:A:1180:MET:HE3	1.97	0.46
1:A:1071:SER:HB2	1:A:1072:PRO:HD2	1.98	0.46
1:A:1110:ARG:NH1	1:A:1159:GLU:OE1	2.49	0.46
1:A:1439:GLN:HA	1:A:1442:ILE:HD12	1.97	0.46
1:B:2090:ASN:HB3	1:B:2093:LEU:HD12	1.98	0.46
1:A:1107:MET:HE2	1:A:1107:MET:HB3	1.89	0.45
1:B:2107:MET:CE	1:B:2208:THR:HB	2.47	0.45
1:C:3422:MET:HA	1:C:3433:LEU:O	2.16	0.45
1:A:1376:ILE:CD1	1:A:1425:MET:HG2	2.47	0.45
1:A:1433:LEU:HA	1:A:1434:PRO:HD3	1.73	0.45
1:B:2431:TYR:N	1:B:2431:TYR:CD1	2.85	0.45
1:A:1242:SER:HB3	1:B:2285:SER:CB	2.46	0.45
1:A:1095:HIS:HE1	1:B:2095:HIS:O	1.99	0.45
1:A:1327:ASP:HB2	1:A:1379:PRO:CA	2.41	0.45
1:C:3269:ARG:HH11	1:C:3406:ALA:HB1	1.80	0.45
1:C:3472:GLU:HG2	1:C:3513:LEU:HD22	1.99	0.45
1:A:1031:LEU:HD12	1:A:1161:PRO:HG2	1.98	0.45
1:B:2191:ARG:HG2	1:B:2191:ARG:NH1	2.30	0.45
1:B:2293:THR:HA	1:B:2367:GLY:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2487:GLY:H	1:B:2490:SER:HB3	1.82	0.45
1:B:2050:ASN:HA	1:B:2051:PRO:HD2	1.87	0.45
1:B:2238:LEU:HB3	1:B:2497:ILE:CG2	2.46	0.45
1:C:3077:GLY:O	1:C:3185:MET:HE2	2.17	0.45
1:A:1292:GLY:O	1:A:1368:ASN:HA	2.17	0.45
1:A:1474:LYS:HG3	1:A:1513:LEU:HD21	1.98	0.45
1:C:3455:GLU:OE1	1:C:3509:ARG:HB2	2.18	0.44
1:A:1244:SER:HB2	1:A:1268:GLY:HA2	1.98	0.44
1:A:1453:VAL:O	1:A:1506:TRP:CD2	2.71	0.44
1:B:2473:PHE:CD1	1:B:2501:PHE:HE2	2.35	0.44
1:A:1061:VAL:HG12	1:A:1062:GLN:N	2.33	0.44
1:B:2222:PRO:O	1:B:2223:THR:CB	2.65	0.44
1:C:3255:ILE:HD11	1:C:3492:PRO:HD2	1.99	0.44
1:C:3238:LEU:HB3	1:C:3497:ILE:CG2	2.47	0.44
1:A:1071:SER:H	1:A:1074:ASN:ND2	2.14	0.44
1:B:2010:SER:OG	1:C:3029:ASP:HB2	2.17	0.44
1:B:2082:ASP:OD1	1:B:2181:ARG:HD2	2.18	0.44
1:C:3271:THR:HG22	1:C:3477:PRO:O	2.17	0.44
1:B:2104:VAL:HG13	1:B:2170:PHE:HA	1.99	0.43
1:C:3194:GLY:C	1:C:3196:THR:H	2.25	0.43
1:C:3280:THR:HA	1:C:3281:PRO:HD3	1.75	0.43
1:A:1422:MET:HA	1:A:1433:LEU:O	2.18	0.43
1:A:1490:SER:HA	1:A:1494:GLN:NE2	2.33	0.43
1:B:2252:SER:HA	1:B:2497:ILE:HD11	1.98	0.43
1:B:2352:PHE:C	1:B:2354:PRO:HD3	2.43	0.43
1:C:3282:VAL:HG23	1:C:3283:SER:N	2.33	0.43
1:B:2431:TYR:N	1:B:2431:TYR:HD1	2.17	0.43
1:B:2013:ASP:HB2	1:B:2155:LEU:HD22	2.00	0.43
1:A:1381:HIS:N	1:A:1382:PRO:HD2	2.34	0.43
1:A:1392:ILE:HD11	1:A:1395:TYR:HE1	1.84	0.43
1:B:2021:LEU:HB3	1:B:2117:ALA:HB1	2.00	0.43
1:B:2069:THR:HG23	1:B:2511:TYR:CE1	2.53	0.43
1:A:1252:SER:HA	1:A:1497:ILE:CD1	2.49	0.43
1:B:2037:ALA:O	1:B:2166:ARG:NH2	2.51	0.43
1:B:2323:LEU:C	1:B:2393:PRO:HD3	2.44	0.43
1:C:3192:THR:HG21	1:C:3200:PHE:CD2	2.52	0.43
1:A:1100:TYR:HA	1:A:1218:PHE:O	2.19	0.43
1:A:1487:GLY:H	1:A:1490:SER:HB3	1.83	0.43
1:B:2022:VAL:HG12	1:B:2153:ARG:HD2	2.00	0.43
1:B:2244:SER:HB3	1:B:2268:GLY:CA	2.49	0.43
1:C:3107:MET:HE3	1:C:3107:MET:HB3	1.95	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3320:PHE:HA	1:C:3321:PRO:HD2	1.85	0.43
1:C:3395:TYR:O	1:C:3396:GLY:C	2.61	0.43
1:A:1413:GLY:HA3	1:A:1416:GLU:CD	2.43	0.43
1:B:2064:PRO:O	1:B:2505:SER:HB2	2.19	0.43
1:C:3300:ASN:HA	1:C:3357:GLY:O	2.18	0.43
1:B:2039:SER:OG	1:B:2215:ASN:HB2	2.19	0.43
1:B:2453:VAL:O	1:B:2506:TRP:CD2	2.72	0.43
1:C:3053:ASP:HA	1:C:3054:PRO:HD2	1.87	0.43
1:A:1255:ILE:HG22	1:A:1256:SER:O	2.18	0.42
1:B:2459:LEU:HD23	1:B:2503:PHE:HA	2.01	0.42
1:C:3109:VAL:HG22	1:C:3208:THR:HG22	2.02	0.42
1:C:3282:VAL:HG23	1:C:3283:SER:H	1.84	0.42
1:A:1031:LEU:HD22	1:A:1146:PRO:HB2	2.01	0.42
1:A:1330:ILE:HG23	1:A:1373:LEU:HD23	1.75	0.42
1:A:1395:TYR:O	1:A:1396:GLY:C	2.62	0.42
1:B:2064:PRO:HG3	1:B:2088:HIS:O	2.19	0.42
1:B:2377:SER:OG	1:B:2378:PRO:HD2	2.19	0.42
1:C:3453:VAL:O	1:C:3506:TRP:CD2	2.72	0.42
1:B:2346:ASP:OD1	1:B:2349:PRO:HD3	2.19	0.42
1:A:1033:MET:HE1	1:A:1180:MET:HE1	2.00	0.42
1:B:2440:GLU:CD	1:B:2440:GLU:H	2.27	0.42
1:A:1243:ASN:OD1	1:A:1244:SER:N	2.53	0.42
1:A:1376:ILE:O	1:A:1428:PRO:HD2	2.20	0.42
1:B:2071:SER:HB2	1:B:2201:VAL:HG12	2.02	0.42
1:C:3329:HIS:HB2	1:C:3343:TYR:O	2.19	0.42
1:C:3244:SER:OG	1:C:3268:GLY:HA3	2.19	0.42
1:A:1263:VAL:HG12	1:A:1264:GLN:N	2.34	0.42
1:C:3031:LEU:HD23	1:C:3031:LEU:HA	1.80	0.42
1:C:3487:GLY:N	1:C:3490:SER:HB3	2.23	0.42
1:B:2284:LEU:HD23	1:B:2287:VAL:HG21	2.02	0.42
1:A:1264:GLN:HG3	1:A:1408:SER:HB3	2.02	0.41
1:A:1271:THR:HG22	1:A:1477:PRO:O	2.20	0.41
1:B:2071:SER:N	1:B:2074:ASN:HD22	2.16	0.41
1:C:3219:LEU:HD23	1:C:3219:LEU:HA	1.90	0.41
1:B:2292:GLY:O	1:B:2368:ASN:HA	2.20	0.41
1:A:1331:ASN:OD1	1:A:1340:GLN:HG3	2.21	0.41
1:B:2327:ASP:HB2	1:B:2379:PRO:CA	2.44	0.41
1:C:3286:HIS:O	1:C:3319:GLY:HA3	2.20	0.41
1:A:1119:THR:CG2	1:A:1200:PHE:CE2	3.03	0.41
1:A:1440:GLU:CD	1:A:1440:GLU:H	2.28	0.41
1:B:2252:SER:HA	1:B:2497:ILE:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2269:ARG:NH2	1:B:2407:PRO:O	2.53	0.41
1:B:2324:GLY:HA3	1:B:2391:LYS:C	2.45	0.41
1:B:2310:HIS:O	1:B:2313:GLU:HB3	2.21	0.41
1:C:3083:LEU:O	1:C:3181:ARG:HA	2.20	0.41
1:A:1271:THR:OG1	1:A:1275:ARG:HB2	2.20	0.41
1:A:1272:LEU:HB2	1:A:1452:THR:CG2	2.43	0.41
1:A:1399:ILE:CG1	1:C:3399:ILE:HG12	2.50	0.41
1:B:2182:LEU:HD23	1:B:2182:LEU:HA	1.78	0.41
1:C:3422:MET:HG2	1:C:3434:PRO:HA	2.02	0.41
1:A:1224:VAL:HG12	1:C:3135:HIS:CD2	2.55	0.41
1:B:2036:VAL:HG11	1:B:2164:ASP:O	2.21	0.41
1:B:2270:CYS:O	1:B:2479:GLY:HA3	2.21	0.41
1:C:3269:ARG:NH1	1:C:3406:ALA:HB1	2.36	0.41
1:A:1346:ASP:OD1	1:A:1381:HIS:HB2	2.20	0.41
1:A:1268:GLY:H	1:A:1280:THR:HG23	1.86	0.41
1:B:2395:TYR:O	1:B:2396:GLY:C	2.64	0.41
1:A:1192:THR:OG1	1:A:1196:THR:HG23	2.21	0.41
1:A:1294:SER:O	1:A:1366:SER:HA	2.21	0.41
1:A:1399:ILE:HG13	1:C:3399:ILE:HG12	2.03	0.41
1:B:2320:PHE:HA	1:B:2321:PRO:HD2	1.80	0.41
1:C:3268:GLY:H	1:C:3280:THR:HG23	1.86	0.41
1:C:3302:THR:HB	1:C:3307:THR:O	2.21	0.41
1:A:1042:ALA:HB2	1:A:1170:PHE:HB3	2.02	0.40
1:A:1277:VAL:HG12	1:A:1278:GLY:N	2.37	0.40
1:A:1421:PHE:CE1	1:A:1438:PRO:HD3	2.56	0.40
1:B:2175:ARG:O	1:B:2175:ARG:HG2	2.20	0.40
1:B:2014:GLY:O	1:B:2015:ALA:C	2.63	0.40
1:A:1458:LEU:HD11	1:A:1472:GLU:HG2	2.04	0.40
1:B:2110:ARG:HA	1:B:2158:ILE:O	2.22	0.40
1:C:3375:TRP:HD1	1:C:3376:ILE:N	2.19	0.40
1:C:3463:ASP:HB2	1:C:3470:LEU:HD11	2.04	0.40
1:A:1433:LEU:HD12	1:A:1434:PRO:CD	2.51	0.40
1:B:2128:ILE:HA	1:B:2129:PRO:HD2	1.92	0.40
1:C:3321:PRO:HB2	1:C:3373:LEU:CD1	2.41	0.40
1:C:3411:PRO:HA	1:C:3412:PRO:HD3	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	490/530 (92%)	444 (91%)	41 (8%)	5 (1%)	12	40
1	B	509/530 (96%)	460 (90%)	39 (8%)	10 (2%)	6	25
1	C	490/530 (92%)	445 (91%)	35 (7%)	10 (2%)	6	25
All	All	1489/1590 (94%)	1349 (91%)	115 (8%)	25 (2%)	7	28

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1383	SER
1	B	2114	ALA
1	C	3383	SER
1	A	1196	THR
1	A	1362	ASN
1	B	2223	THR
1	B	2486	ASN
1	C	3173	ASN
1	C	3196	THR
1	C	3362	ASN
1	C	3396	GLY
1	B	2040	SER
1	B	2196	THR
1	B	2222	PRO
1	B	2362	ASN
1	B	2363	GLY
1	B	2383	SER
1	C	3043	VAL
1	C	3363	GLY
1	C	3324	GLY
1	C	3397	SER
1	B	2396	GLY
1	C	3057	ILE
1	A	1363	GLY

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Mol	Chain	Res	Type
1	A	1396	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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	418/445 (94%)	392 (94%)	26 (6%)	16	43
1	B	431/445 (97%)	408 (95%)	23 (5%)	20	47
1	C	418/445 (94%)	394 (94%)	24 (6%)	18	45
All	All	1267/1335 (95%)	1194 (94%)	73 (6%)	18	45

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1041	THR
1	A	1045	THR
1	A	1084	SER
1	A	1101	ASN
1	A	1118	PHE
1	A	1124	ILE
1	A	1141	GLN
1	A	1154	THR
1	A	1179	THR
1	A	1180	MET
1	A	1181	ARG
1	A	1191	ARG
1	A	1220	VAL
1	A	1259	ASN
1	A	1264	GLN
1	A	1271	THR
1	A	1284	LEU
1	A	1307	THR
1	A	1329	HIS
1	A	1385	SER
1	A	1392	ILE

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Mol	Chain	Res	Type
1	A	1408	SER
1	A	1423	SER
1	A	1425	MET
1	A	1432	ASN
1	A	1440	GLU
1	B	2011	SER
1	B	2084	SER
1	B	2104	VAL
1	B	2126	SER
1	B	2139	ILE
1	B	2141	GLN
1	B	2153	ARG
1	B	2182	LEU
1	B	2220	VAL
1	B	2223	THR
1	B	2240	SER
1	B	2259	ASN
1	B	2273	ASP
1	B	2305	ASP
1	B	2307	THR
1	B	2329	HIS
1	B	2360	GLN
1	B	2387	VAL
1	B	2423	SER
1	B	2425	MET
1	B	2431	TYR
1	B	2440	GLU
1	B	2460	HIS
1	C	3031	LEU
1	C	3057	ILE
1	C	3101	ASN
1	C	3104	VAL
1	C	3107	MET
1	C	3108	ARG
1	C	3123	ILE
1	C	3135	HIS
1	C	3136	ASN
1	C	3152	VAL
1	C	3258	ASP
1	C	3271	THR
1	C	3305	ASP
1	C	3307	THR

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Mol	Chain	Res	Type
1	C	3338	SER
1	C	3345	VAL
1	C	3358	SER
1	C	3375	TRP
1	C	3387	VAL
1	C	3400	THR
1	C	3423	SER
1	C	3440	GLU
1	C	3509	ARG
1	C	3516	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1074	ASN
1	A	1095	HIS
1	A	1106	ASN
1	A	1116	ASN
1	A	1171	HIS
1	A	1286	HIS
1	A	1340	GLN
1	A	1342	GLN
1	A	1360	GLN
1	A	1394	ASN
1	B	2026	ASN
1	B	2062	GLN
1	B	2073	ASN
1	B	2074	ASN
1	B	2095	HIS
1	B	2116	ASN
1	B	2135	HIS
1	B	2171	HIS
1	B	2177	GLN
1	B	2342	GLN
1	B	2360	GLN
1	B	2444	HIS
1	B	2512	GLN
1	C	3050	ASN
1	C	3059	ASN
1	C	3074	ASN
1	C	3098	GLN
1	C	3135	HIS

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Mol	Chain	Res	Type
1	C	3171	HIS
1	C	3334	GLN
1	C	3337	HIS
1	C	3444	HIS
1	C	3469	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](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.