



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 10:33 AM UTC

PDB ID : 4LEE / pdb_00004lee
Title : Structure of the Als3 adhesin from *Candida albicans*, residues 1-313 (mature sequence), triple mutant in the binding cavity: K59M, A116V, Y301F
Authors : Lin, J.; Garnett, J.A.; Cota, E.
Deposited on : 2013-06-25
Resolution : 3.00 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

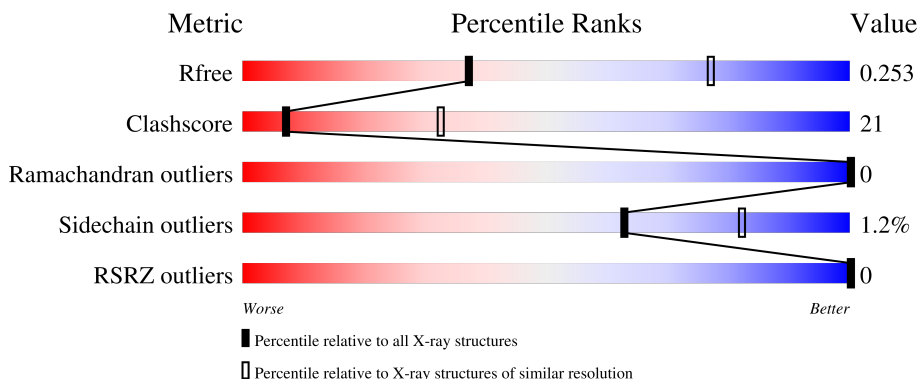
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	314	72% 25% ..
1	B	314	70% 28% .
1	C	314	72% 27% .
1	D	314	78% 20% .

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 9277 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 Agglutinin-like protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2309	1460	372	465	12			
1	B	313	Total	C	N	O	S	0	0	0
			2314	1464	373	465	12			
1	C	313	Total	C	N	O	S	0	0	0
			2318	1465	372	470	11			
1	D	313	Total	C	N	O	S	0	0	0
			2336	1480	371	473	12			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP O74623
A	59	MET	LYS	engineered mutation	UNP O74623
A	116	VAL	ALA	engineered mutation	UNP O74623
A	298	PHE	TYR	engineered mutation	UNP O74623
B	0	GLY	-	expression tag	UNP O74623
B	59	MET	LYS	engineered mutation	UNP O74623
B	116	VAL	ALA	engineered mutation	UNP O74623
B	298	PHE	TYR	engineered mutation	UNP O74623
C	0	GLY	-	expression tag	UNP O74623
C	59	MET	LYS	engineered mutation	UNP O74623
C	116	VAL	ALA	engineered mutation	UNP O74623
C	298	PHE	TYR	engineered mutation	UNP O74623
D	0	GLY	-	expression tag	UNP O74623
D	59	MET	LYS	engineered mutation	UNP O74623
D	116	VAL	ALA	engineered mutation	UNP O74623
D	298	PHE	TYR	engineered mutation	UNP O74623

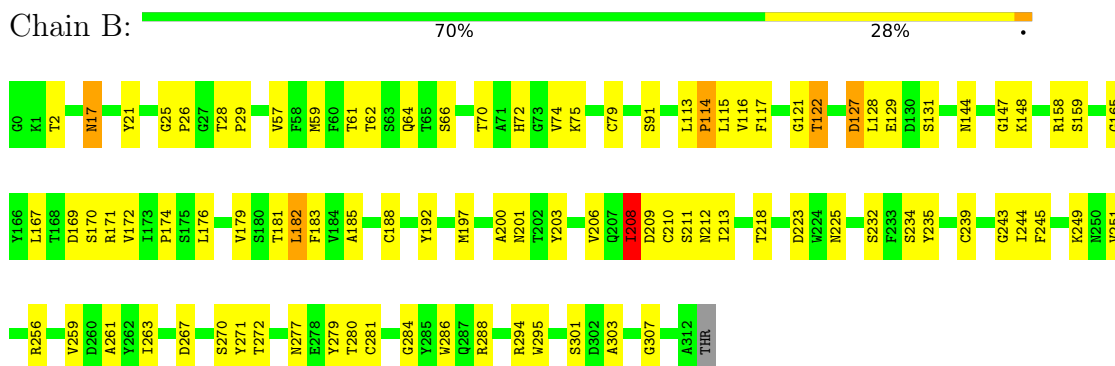
3 Residue-property plots

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.

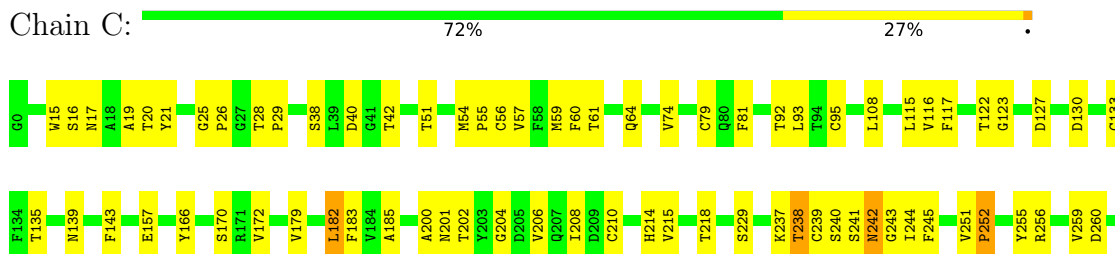
• Molecule 1: Agglutinin-like protein 3

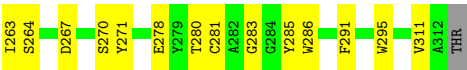


• Molecule 1: Agglutinin-like protein 3

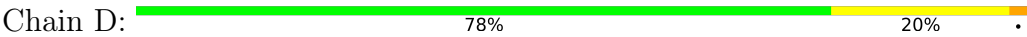


• Molecule 1: Agglutinin-like protein 3





● Molecule 1: Agglutinin-like protein 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	112.56Å 67.09Å 112.69Å 90.00° 103.55° 90.00°	Depositor
Resolution (Å)	48.33 – 3.00 48.33 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.0 (48.33-3.00) 97.0 (48.33-3.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.229 , 0.242 0.241 , 0.253	Depositor DCC
R_{free} test set	1634 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	46.2	Xtriage
Anisotropy	0.554	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 16.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.447 for l,-k,h	Xtriage
Reported twinning fraction	0.492 for H, K, L 0.508 for -L, -K, -H	Depositor
Outliers	1 of 32253 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9277	wwPDB-VP
Average B, all atoms (Å ²)	45.0	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 25.46 % 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 3.1595e-03. 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 [i](#)

5.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.69	3/2367 (0.1%)	1.01	16/3246 (0.5%)
1	B	0.71	5/2373 (0.2%)	0.95	10/3255 (0.3%)
1	C	0.72	2/2377 (0.1%)	0.98	8/3261 (0.2%)
1	D	0.64	1/2396 (0.0%)	0.98	7/3283 (0.2%)
All	All	0.69	11/9513 (0.1%)	0.98	41/13045 (0.3%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	54	MET	SD-CE	-7.43	1.60	1.79
1	A	70	THR	C-O	-6.34	1.16	1.23
1	D	121	GLY	C-O	-6.14	1.15	1.23
1	B	114	PRO	N-CD	5.95	1.56	1.47
1	B	113	LEU	CA-C	-5.86	1.46	1.53
1	B	148	LYS	C-O	-5.57	1.17	1.24
1	A	120	GLY	C-O	-5.55	1.18	1.23
1	C	238	THR	CA-C	5.46	1.59	1.53
1	B	176	LEU	C-O	-5.30	1.17	1.24
1	B	113	LEU	C-O	-5.10	1.17	1.25
1	C	252	PRO	N-CD	5.07	1.54	1.47

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	20	THR	N-CA-C	9.91	123.47	112.97
1	C	238	THR	N-CA-C	8.24	120.91	107.32
1	D	234	SER	N-CA-C	7.87	120.87	108.67
1	D	122	THR	N-CA-C	-7.82	97.25	107.73
1	A	20	THR	N-CA-CB	-7.52	99.97	110.65
1	B	129	GLU	N-CA-C	-7.50	103.05	111.07
1	D	172	VAL	CB-CA-C	-7.41	101.84	111.25
1	C	311	VAL	N-CA-C	7.37	118.91	108.36
1	B	212	ASN	N-CA-C	7.15	120.59	107.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	20	THR	N-CA-C	6.94	120.33	110.68
1	C	242	ASN	N-CA-C	6.90	120.18	111.69
1	B	127	ASP	N-CA-C	6.89	118.79	111.28
1	A	118	ASN	CA-C-N	-6.76	113.87	122.26
1	A	118	ASN	C-N-CA	-6.76	113.87	122.26
1	A	234	SER	N-CA-C	6.68	120.18	108.56
1	B	277	ASN	N-CA-C	6.58	120.03	109.83
1	A	235	TYR	N-CA-C	6.52	119.78	109.81
1	C	59	MET	N-CA-C	6.50	118.95	109.07
1	B	122	THR	N-CA-CB	6.47	121.43	110.49
1	A	249	LYS	N-CA-C	6.38	118.77	109.07
1	A	249	LYS	N-CA-CB	-6.15	101.39	111.66
1	D	123	GLY	N-CA-C	5.89	127.14	113.18
1	B	208	ILE	N-CA-C	5.89	117.22	109.21
1	A	234	SER	CB-CA-C	-5.88	102.85	112.03
1	A	250	ASN	N-CA-C	5.79	119.66	112.24
1	A	251	VAL	N-CA-C	5.75	121.30	108.88
1	A	70	THR	N-CA-C	5.73	118.91	107.62
1	A	19	ALA	N-CA-C	-5.72	98.61	110.80
1	D	311	VAL	N-CA-C	5.70	116.51	108.36
1	A	208	ILE	N-CA-C	5.55	116.58	108.58
1	C	238	THR	N-CA-CB	-5.51	101.33	111.36
1	B	17	ASN	CB-CA-C	5.46	118.03	109.84
1	B	17	ASN	N-CA-C	-5.43	102.26	110.24
1	A	235	TYR	N-CA-CB	-5.41	102.93	111.05
1	C	202	THR	N-CA-C	5.36	120.06	112.93
1	A	105	ILE	N-CA-C	5.36	116.00	109.30
1	C	157	GLU	N-CA-C	-5.22	102.33	110.42
1	D	282	ALA	N-CA-C	5.19	119.24	113.01
1	B	169	ASP	CA-C-O	-5.14	114.36	120.43
1	D	238	THR	N-CA-C	5.06	117.33	107.71
1	B	251	VAL	N-CA-C	5.02	119.72	108.88

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	2309	0	2144	93	0
1	B	2314	0	2138	95	2
1	C	2318	0	2141	102	2
1	D	2336	0	2172	82	0
All	All	9277	0	8595	369	2

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 21.

All (369) 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:D:60:PHE:CD2	1:D:115:LEU:HD21	1.74	1.22
1:B:280:THR:HG1	1:B:286:TRP:CD1	1.60	1.19
1:C:79:CYS:CB	1:C:95:CYS:SG	2.31	1.17
1:C:79:CYS:SG	1:C:95:CYS:CB	2.33	1.15
1:C:210:CYS:SG	1:C:239:CYS:SG	1.38	1.13
1:D:280:THR:OG1	1:D:286:TRP:HD1	1.27	1.13
1:A:202:THR:HG21	1:A:272:THR:HB	1.15	1.13
1:C:210:CYS:CB	1:C:239:CYS:SG	2.39	1.10
1:D:280:THR:HG1	1:D:286:TRP:CD1	1.68	1.10
1:C:122:THR:HG22	1:C:123:GLY:H	0.94	1.09
1:D:119:VAL:HG11	1:D:224:TRP:CE3	1.89	1.06
1:C:122:THR:HG22	1:C:123:GLY:N	1.71	1.04
1:D:281:CYS:HB2	1:D:285:TYR:HB3	1.39	1.04
1:A:76:TYR:CD1	1:A:100:THR:HG23	1.90	1.04
1:D:267:ASP:HB3	1:D:270:SER:O	1.56	1.04
1:C:208:ILE:HG12	1:C:263:ILE:HG12	1.39	1.03
1:A:119:VAL:HG11	1:A:224:TRP:CE3	1.93	1.02
1:B:280:THR:OG1	1:B:286:TRP:HD1	1.41	1.01
1:B:167:LEU:HD21	1:B:225:ASN:ND2	1.76	1.00
1:C:122:THR:CG2	1:C:123:GLY:H	1.75	0.99
1:C:206:VAL:HG13	1:C:263:ILE:HG21	1.45	0.98
1:B:29:PRO:HB3	1:B:116:VAL:CG2	1.95	0.96
1:A:60:PHE:CE2	1:A:115:LEU:HD21	2.03	0.93
1:D:267:ASP:OD2	1:D:270:SER:HB3	1.68	0.93
1:A:119:VAL:HG12	1:A:119:VAL:O	1.70	0.91
1:B:167:LEU:CD2	1:B:225:ASN:HB3	2.02	0.90
1:C:206:VAL:HG13	1:C:263:ILE:CG2	2.03	0.89
1:D:60:PHE:CD2	1:D:115:LEU:CD2	2.56	0.89
1:A:202:THR:CG2	1:A:272:THR:HB	2.04	0.88
1:A:218:THR:HG22	1:A:220:GLY:O	1.74	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:60:PHE:CE2	1:D:115:LEU:HD21	2.08	0.88
1:B:62:THR:HG22	1:B:301:SER:O	1.74	0.87
1:D:280:THR:OG1	1:D:286:TRP:CD1	2.20	0.87
1:A:76:TYR:CD1	1:A:100:THR:CG2	2.58	0.86
1:C:54:MET:HB2	1:C:143:PHE:CE1	2.11	0.85
1:A:52:LEU:HD23	1:A:54:MET:HE3	1.59	0.85
1:A:60:PHE:CD2	1:A:115:LEU:CD2	2.60	0.84
1:B:280:THR:HG23	1:B:286:TRP:HB2	1.58	0.83
1:D:280:THR:HG1	1:D:286:TRP:HD1	0.87	0.83
1:A:60:PHE:CD2	1:A:115:LEU:HD21	2.14	0.82
1:B:197:MET:HE2	1:B:259:VAL:HG22	1.59	0.82
1:C:208:ILE:HG21	1:C:244:ILE:HD13	1.61	0.82
1:D:119:VAL:HG12	1:D:119:VAL:O	1.78	0.82
1:B:183:PHE:CD2	1:B:197:MET:HE3	2.15	0.81
1:C:252:PRO:HG2	1:C:255:TYR:CD1	2.16	0.81
1:A:251:VAL:HG13	1:A:252:PRO:HD2	1.61	0.81
1:C:60:PHE:CE2	1:C:115:LEU:HD21	2.16	0.81
1:C:206:VAL:CG1	1:C:263:ILE:HG21	2.11	0.80
1:D:59:MET:HE1	1:D:174:PRO:HG3	1.63	0.80
1:A:252:PRO:HG2	1:A:255:TYR:CD1	2.17	0.80
1:B:167:LEU:CD2	1:B:225:ASN:ND2	2.45	0.79
1:A:70:THR:HG22	1:A:70:THR:O	1.83	0.79
1:D:249:LYS:HG2	1:D:250:ASN:H	1.45	0.79
1:A:76:TYR:HD1	1:A:100:THR:HG23	1.48	0.78
1:A:202:THR:HG22	1:A:272:THR:O	1.84	0.78
1:B:29:PRO:HB3	1:B:116:VAL:HG21	1.66	0.78
1:A:202:THR:HG21	1:A:272:THR:CB	2.08	0.77
1:C:122:THR:HB	1:C:127:ASP:OD2	1.85	0.76
1:A:251:VAL:HG13	1:A:252:PRO:CD	2.16	0.76
1:C:38:SER:CB	1:C:108:LEU:CD2	2.63	0.76
1:B:167:LEU:HD23	1:B:225:ASN:HB3	1.65	0.76
1:B:172:VAL:HG12	1:B:179:VAL:HG22	1.66	0.75
1:C:38:SER:HB2	1:C:108:LEU:CD2	2.17	0.75
1:C:215:VAL:HG22	1:C:259:VAL:HG23	1.69	0.75
1:D:208:ILE:HG21	1:D:244:ILE:CD1	2.17	0.74
1:C:54:MET:HB2	1:C:143:PHE:CD1	2.22	0.74
1:D:58:PHE:CE1	1:D:171:ARG:NH1	2.55	0.74
1:A:60:PHE:CE2	1:A:115:LEU:CD2	2.71	0.74
1:A:119:VAL:HG11	1:A:224:TRP:CZ3	2.22	0.74
1:D:118:ASN:O	1:D:130:ASP:HB3	1.88	0.73
1:D:58:PHE:CE2	1:D:171:ARG:HD2	2.24	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:ILE:CG2	1:D:244:ILE:HD11	2.18	0.73
1:D:208:ILE:HG21	1:D:244:ILE:HG13	1.70	0.73
1:C:38:SER:HA	1:C:108:LEU:HD23	1.71	0.72
1:A:218:THR:CG2	1:A:220:GLY:O	2.38	0.72
1:C:79:CYS:SG	1:C:95:CYS:SG	0.82	0.72
1:A:60:PHE:CD2	1:A:115:LEU:HD23	2.25	0.71
1:D:239:CYS:SG	1:D:244:ILE:HG12	2.29	0.71
1:C:38:SER:OG	1:C:108:LEU:HD21	1.90	0.71
1:C:122:THR:CG2	1:C:214:HIS:ND1	2.54	0.70
1:C:280:THR:OG1	1:C:286:TRP:HD1	1.74	0.70
1:D:208:ILE:CG2	1:D:244:ILE:CD1	2.70	0.70
1:A:66:SER:HB2	1:A:79:CYS:O	1.92	0.69
1:C:122:THR:HG23	1:C:260:ASP:CG	2.16	0.69
1:B:21:TYR:OH	1:B:29:PRO:HD2	1.93	0.69
1:B:280:THR:HG1	1:B:286:TRP:HD1	0.71	0.69
1:C:251:VAL:HG12	1:C:252:PRO:HD2	1.74	0.69
1:C:280:THR:HG1	1:C:286:TRP:CD1	2.11	0.69
1:A:119:VAL:O	1:A:119:VAL:CG1	2.41	0.69
1:C:251:VAL:HG12	1:C:252:PRO:CD	2.23	0.68
1:C:57:VAL:HG11	1:C:115:LEU:HD22	1.74	0.68
1:C:60:PHE:CD2	1:C:115:LEU:HD21	2.29	0.68
1:B:167:LEU:CD2	1:B:225:ASN:CB	2.72	0.68
1:A:72:HIS:C	1:A:74:VAL:H	2.02	0.68
1:A:271:TYR:HE2	1:A:273:LEU:HD11	1.59	0.67
1:B:288:ARG:O	1:B:288:ARG:HG3	1.94	0.67
1:B:57:VAL:HG13	1:B:115:LEU:HD13	1.75	0.67
1:C:79:CYS:CB	1:C:95:CYS:HG	1.92	0.67
1:C:251:VAL:CG1	1:C:252:PRO:CD	2.73	0.67
1:B:74:VAL:HG12	1:B:75:LYS:H	1.61	0.66
1:A:59:MET:HE1	1:A:174:PRO:HG2	1.77	0.66
1:D:119:VAL:CG1	1:D:224:TRP:CE3	2.74	0.66
1:D:249:LYS:CG	1:D:250:ASN:H	2.09	0.66
1:B:61:THR:CG2	1:B:116:VAL:HG23	2.26	0.66
1:B:29:PRO:CB	1:B:116:VAL:CG2	2.72	0.65
1:B:172:VAL:HA	1:B:179:VAL:HG22	1.76	0.65
1:C:278:GLU:OE2	1:C:286:TRP:NE1	2.29	0.65
1:D:58:PHE:CZ	1:D:171:ARG:HD2	2.31	0.65
1:A:162:ASP:OD1	1:A:164:LYS:HG3	1.96	0.65
1:A:52:LEU:CD2	1:A:54:MET:HE3	2.26	0.65
1:C:281:CYS:HB2	1:C:285:TYR:CB	2.27	0.65
1:B:17:ASN:OD1	1:B:21:TYR:HD2	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:267:ASP:OD2	1:D:270:SER:CB	2.43	0.64
1:D:208:ILE:HG21	1:D:244:ILE:CG1	2.27	0.64
1:D:26:PRO:HG2	1:D:119:VAL:HG21	1.79	0.64
1:B:172:VAL:HG12	1:B:179:VAL:CG2	2.28	0.64
1:B:201:ASN:CB	1:B:206:VAL:HB	2.27	0.64
1:B:213:ILE:HG21	1:B:244:ILE:HD12	1.79	0.64
1:B:59:MET:HE1	1:B:174:PRO:HG2	1.79	0.64
1:C:56:CYS:SG	1:C:133:CYS:CB	2.86	0.63
1:A:239:CYS:HB2	1:A:244:ILE:HD13	1.79	0.63
1:D:59:MET:CE	1:D:174:PRO:HG3	2.29	0.63
1:A:26:PRO:HG2	1:A:119:VAL:HG21	1.81	0.62
1:B:64:GLN:OE1	1:B:303:ALA:HB3	1.99	0.62
1:C:210:CYS:CA	1:C:239:CYS:SG	2.87	0.62
1:C:206:VAL:CG1	1:C:263:ILE:CG2	2.76	0.62
1:A:102:THR:HB	1:A:103:PRO:HD2	1.81	0.62
1:C:280:THR:HG23	1:C:286:TRP:HB2	1.82	0.62
1:A:119:VAL:HG11	1:A:224:TRP:HE3	1.60	0.61
1:A:294:ARG:NH1	1:A:296:THR:CG2	2.63	0.61
1:B:279:TYR:CD2	1:B:281:CYS:SG	2.94	0.61
1:C:251:VAL:CG1	1:C:252:PRO:HD2	2.30	0.61
1:A:192:TYR:HB2	1:A:251:VAL:HB	1.82	0.61
1:B:167:LEU:HD23	1:B:225:ASN:CB	2.30	0.61
1:C:60:PHE:CD2	1:C:115:LEU:CD2	2.83	0.61
1:C:271:TYR:CZ	1:C:295:TRP:HB2	2.35	0.60
1:B:121:GLY:O	1:B:122:THR:HG23	2.01	0.60
1:B:208:ILE:HD13	1:B:243:GLY:HA2	1.84	0.59
1:C:280:THR:HG1	1:C:286:TRP:HD1	1.48	0.59
1:D:268:VAL:HG12	1:D:268:VAL:O	2.01	0.59
1:A:144:ASN:OD1	1:A:145:ASP:N	2.36	0.59
1:D:239:CYS:SG	1:D:244:ILE:CD1	2.90	0.59
1:D:176:LEU:O	1:D:178:LYS:HG3	2.03	0.58
1:D:119:VAL:HG11	1:D:224:TRP:CZ3	2.36	0.58
1:A:105:ILE:HD13	1:A:311:VAL:HG22	1.86	0.58
1:B:29:PRO:CA	1:B:116:VAL:HG22	2.33	0.58
1:D:279:TYR:CE2	1:D:285:TYR:HE2	2.21	0.58
1:B:21:TYR:CZ	1:B:29:PRO:HD2	2.38	0.58
1:C:170:SER:OG	1:C:179:VAL:CG1	2.50	0.58
1:D:213:ILE:HG21	1:D:244:ILE:HD13	1.86	0.58
1:B:167:LEU:CD2	1:B:225:ASN:HD22	2.16	0.57
1:B:279:TYR:HD2	1:B:281:CYS:SG	2.27	0.57
1:B:203:TYR:OH	1:B:294:ARG:NH2	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:239:CYS:SG	1:D:244:ILE:HD11	2.43	0.57
1:B:213:ILE:HG22	1:B:261:ALA:HB2	1.85	0.57
1:A:239:CYS:HB2	1:A:244:ILE:CD1	2.34	0.57
1:D:60:PHE:HA	1:D:115:LEU:HD23	1.85	0.57
1:B:29:PRO:CB	1:B:116:VAL:HG21	2.34	0.57
1:B:209:ASP:OD1	1:B:210:CYS:N	2.38	0.57
1:C:237:LYS:C	1:C:238:THR:HG23	2.30	0.57
1:B:167:LEU:HD21	1:B:225:ASN:CG	2.30	0.57
1:C:54:MET:CB	1:C:143:PHE:CD1	2.88	0.57
1:A:271:TYR:CE2	1:A:273:LEU:HD11	2.39	0.56
1:B:167:LEU:HD22	1:B:225:ASN:HB3	1.83	0.56
1:B:29:PRO:HA	1:B:116:VAL:HG22	1.87	0.56
1:D:121:GLY:O	1:D:122:THR:CG2	2.53	0.56
1:A:72:HIS:C	1:A:74:VAL:N	2.62	0.56
1:D:122:THR:HG22	1:D:260:ASP:OD1	2.06	0.56
1:D:267:ASP:C	1:D:269:ASN:H	2.14	0.56
1:B:167:LEU:CD2	1:B:225:ASN:CG	2.79	0.56
1:B:17:ASN:OD1	1:B:21:TYR:CD2	2.60	0.56
1:D:119:VAL:O	1:D:119:VAL:CG1	2.51	0.55
1:B:61:THR:HG21	1:B:116:VAL:HG23	1.86	0.55
1:A:59:MET:HE1	1:A:174:PRO:CG	2.37	0.55
1:A:251:VAL:CG1	1:A:252:PRO:CD	2.83	0.55
1:D:58:PHE:CD1	1:D:171:ARG:NH1	2.74	0.55
1:A:72:HIS:O	1:A:74:VAL:N	2.39	0.55
1:B:213:ILE:HG21	1:B:244:ILE:CD1	2.35	0.55
1:C:259:VAL:HG13	1:C:259:VAL:O	2.07	0.54
1:A:208:ILE:HD13	1:A:243:GLY:HA2	1.89	0.54
1:D:119:VAL:HG11	1:D:224:TRP:HE3	1.65	0.54
1:C:38:SER:CB	1:C:108:LEU:HD21	2.37	0.54
1:B:59:MET:HE1	1:B:174:PRO:CG	2.38	0.54
1:C:38:SER:HB2	1:C:108:LEU:HD22	1.88	0.54
1:A:271:TYR:HE2	1:A:273:LEU:CD1	2.20	0.54
1:C:17:ASN:OD1	1:C:21:TYR:HD2	1.90	0.54
1:D:280:THR:HA	1:D:285:TYR:O	2.08	0.53
1:C:185:ALA:O	1:C:256:ARG:NH2	2.30	0.53
1:C:55:PRO:O	1:C:56:CYS:HB2	2.07	0.53
1:C:280:THR:HA	1:C:285:TYR:O	2.08	0.53
1:D:121:GLY:C	1:D:122:THR:HG23	2.34	0.53
1:C:263:ILE:CG2	1:C:264:SER:N	2.72	0.53
1:D:281:CYS:SG	1:D:285:TYR:HD2	2.32	0.53
1:A:294:ARG:HH12	1:A:296:THR:CG2	2.21	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:SER:CA	1:C:108:LEU:HD23	2.39	0.53
1:B:62:THR:HG23	1:B:114:PRO:HG2	1.91	0.52
1:C:263:ILE:HG22	1:C:264:SER:N	2.24	0.52
1:A:210:CYS:HG	1:A:239:CYS:HG	1.54	0.52
1:C:239:CYS:HA	1:C:244:ILE:HG13	1.91	0.52
1:B:201:ASN:HB2	1:B:206:VAL:HB	1.91	0.52
1:C:122:THR:HG23	1:C:260:ASP:CB	2.40	0.52
1:D:239:CYS:SG	1:D:244:ILE:CG1	2.96	0.52
1:B:121:GLY:O	1:B:122:THR:CG2	2.58	0.52
1:A:119:VAL:HG12	1:A:222:ASN:ND2	2.25	0.51
1:B:201:ASN:HB3	1:B:206:VAL:HB	1.91	0.51
1:D:269:ASN:OD1	1:D:297:GLY:HA3	2.11	0.51
1:A:25:GLY:O	1:A:28:THR:OG1	2.27	0.51
1:A:201:ASN:HB2	1:A:206:VAL:HB	1.93	0.51
1:D:231:GLU:HG2	1:D:232:SER:H	1.76	0.51
1:A:270:SER:HB2	1:A:294:ARG:NH2	2.26	0.51
1:A:293:LEU:HD23	1:A:295:TRP:CZ2	2.45	0.51
1:C:130:ASP:O	1:C:133:CYS:SG	2.69	0.50
1:D:201:ASN:HD21	1:D:204:GLY:HA2	1.77	0.50
1:B:66:SER:HB2	1:B:79:CYS:O	2.11	0.50
1:C:201:ASN:HD21	1:C:204:GLY:HA2	1.77	0.50
1:A:26:PRO:HB2	1:A:117:PHE:HB2	1.94	0.49
1:D:25:GLY:O	1:D:28:THR:OG1	2.27	0.49
1:B:158:ARG:HG2	1:B:159:SER:N	2.26	0.49
1:B:271:TYR:CZ	1:B:295:TRP:HB2	2.46	0.49
1:A:76:TYR:CE1	1:A:100:THR:HG23	2.43	0.49
1:C:271:TYR:CZ	1:C:295:TRP:CB	2.95	0.49
1:D:105:ILE:HD13	1:D:311:VAL:HG22	1.94	0.49
1:A:81:PHE:CD1	1:A:93:LEU:HD21	2.48	0.49
1:A:185:ALA:O	1:A:256:ARG:NH2	2.33	0.49
1:D:208:ILE:HG21	1:D:244:ILE:HD11	1.87	0.49
1:A:251:VAL:CG1	1:A:252:PRO:HD2	2.38	0.49
1:B:59:MET:O	1:B:115:LEU:HB2	2.12	0.49
1:C:38:SER:OG	1:C:108:LEU:CD2	2.60	0.49
1:A:294:ARG:NH1	1:A:296:THR:HG22	2.27	0.48
1:C:60:PHE:CE2	1:C:115:LEU:CD2	2.93	0.48
1:B:158:ARG:HG3	1:B:223:ASP:O	2.13	0.48
1:C:25:GLY:O	1:C:28:THR:OG1	2.27	0.48
1:C:251:VAL:HG12	1:C:252:PRO:N	2.28	0.48
1:C:210:CYS:HA	1:C:239:CYS:SG	2.53	0.48
1:B:25:GLY:O	1:B:28:THR:OG1	2.26	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:VAL:CG1	1:A:252:PRO:N	2.76	0.48
1:D:234:SER:HB2	1:D:249:LYS:HB3	1.95	0.48
1:A:218:THR:HG21	1:A:227:PRO:HA	1.95	0.48
1:B:209:ASP:OD1	1:B:211:SER:N	2.47	0.48
1:D:185:ALA:O	1:D:256:ARG:NH2	2.32	0.48
1:D:279:TYR:HE2	1:D:285:TYR:HE2	1.61	0.48
1:A:84:GLY:HA2	1:A:88:MET:HG3	1.95	0.48
1:A:128:LEU:HD13	1:B:232:SER:OG	2.13	0.48
1:B:172:VAL:CG1	1:B:179:VAL:HG22	2.42	0.48
1:C:38:SER:CB	1:C:108:LEU:HD23	2.43	0.48
1:D:163:PRO:HA	1:D:167:LEU:HD21	1.96	0.48
1:B:267:ASP:O	1:B:270:SER:O	2.32	0.47
1:B:57:VAL:O	1:B:91:SER:OG	2.31	0.47
1:D:281:CYS:SG	1:D:285:TYR:CD2	3.07	0.47
1:C:15:TRP:CZ2	1:C:17:ASN:CG	2.92	0.47
1:D:269:ASN:O	1:D:296:THR:HA	2.15	0.47
1:A:200:ALA:HB2	1:A:245:PHE:HZ	1.78	0.47
1:B:170:SER:HB2	1:B:179:VAL:CG1	2.45	0.47
1:B:170:SER:HB3	1:B:181:THR:HB	1.96	0.47
1:C:16:SER:O	1:C:17:ASN:C	2.57	0.47
1:C:208:ILE:HG12	1:C:263:ILE:CG1	2.27	0.47
1:B:213:ILE:HG22	1:B:261:ALA:CB	2.45	0.47
1:D:249:LYS:HG2	1:D:250:ASN:N	2.23	0.47
1:A:119:VAL:HG12	1:A:222:ASN:HD21	1.80	0.47
1:A:125:SER:HB2	1:B:235:TYR:CE1	2.50	0.47
1:C:51:THR:HG23	1:C:92:THR:OG1	2.15	0.47
1:D:271:TYR:CZ	1:D:295:TRP:HB2	2.50	0.47
1:A:267:ASP:O	1:A:270:SER:O	2.32	0.46
1:B:29:PRO:CB	1:B:116:VAL:HG22	2.44	0.46
1:B:179:VAL:O	1:B:263:ILE:HG12	2.15	0.46
1:C:240:SER:C	1:C:242:ASN:N	2.70	0.46
1:C:240:SER:O	1:C:241:SER:C	2.57	0.46
1:A:17:ASN:CG	1:A:19:ALA:O	2.59	0.46
1:A:121:GLY:C	1:A:122:THR:CG2	2.88	0.46
1:D:121:GLY:O	1:D:122:THR:HG23	2.15	0.46
1:B:29:PRO:HB3	1:B:116:VAL:HG22	1.90	0.46
1:D:249:LYS:CG	1:D:250:ASN:N	2.78	0.46
1:B:2:THR:O	1:B:2:THR:HG22	2.16	0.46
1:C:122:THR:HG21	1:C:214:HIS:ND1	2.28	0.46
1:D:54:MET:HE2	1:D:91:SER:HB2	1.96	0.46
1:B:185:ALA:O	1:B:256:ARG:NH2	2.33	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:208:ILE:HG13	1:A:263:ILE:HG22	1.98	0.46
1:B:171:ARG:C	1:B:179:VAL:HG13	2.40	0.46
1:A:163:PRO:HA	1:A:167:LEU:HD21	1.98	0.46
1:C:251:VAL:HG13	1:C:252:PRO:CD	2.46	0.46
1:A:121:GLY:C	1:A:122:THR:HG23	2.41	0.46
1:C:56:CYS:SG	1:C:133:CYS:HB2	2.56	0.45
1:C:243:GLY:C	1:C:244:ILE:HD12	2.42	0.45
1:B:72:HIS:CD2	1:B:307:GLY:O	2.70	0.45
1:A:183:PHE:CD2	1:A:197:MET:HE3	2.52	0.45
1:B:234:SER:HB2	1:B:249:LYS:HG2	1.98	0.45
1:A:192:TYR:HD1	1:A:251:VAL:HG11	1.81	0.45
1:B:122:THR:OG1	1:B:127:ASP:OD2	2.29	0.45
1:D:267:ASP:OD2	1:D:270:SER:N	2.50	0.45
1:D:267:ASP:C	1:D:269:ASN:N	2.73	0.45
1:C:170:SER:OG	1:C:179:VAL:HG12	2.15	0.45
1:C:200:ALA:HB2	1:C:245:PHE:CZ	2.52	0.45
1:A:280:THR:HG1	1:A:286:TRP:CD1	2.34	0.45
1:D:54:MET:HE3	1:D:57:VAL:HG11	1.99	0.45
1:A:57:VAL:HG11	1:A:115:LEU:HD22	1.98	0.45
1:A:192:TYR:CD1	1:A:251:VAL:HG11	2.52	0.45
1:B:70:THR:HG22	1:B:75:LYS:HB2	1.99	0.45
1:B:188:CYS:HB2	1:B:192:TYR:CZ	2.52	0.45
1:C:64:GLN:O	1:C:81:PHE:CZ	2.70	0.45
1:C:122:THR:HG22	1:C:214:HIS:ND1	2.30	0.45
1:D:57:VAL:O	1:D:91:SER:OG	2.31	0.44
1:B:158:ARG:CG	1:B:159:SER:N	2.81	0.44
1:C:60:PHE:CD2	1:C:115:LEU:HD23	2.52	0.44
1:C:200:ALA:HB2	1:C:245:PHE:HZ	1.82	0.44
1:D:54:MET:CE	1:D:57:VAL:HG11	2.47	0.44
1:D:267:ASP:O	1:D:270:SER:O	2.35	0.44
1:B:144:ASN:OD1	1:B:147:GLY:N	2.50	0.44
1:C:26:PRO:HB2	1:C:117:PHE:HB2	1.99	0.44
1:D:121:GLY:C	1:D:122:THR:CG2	2.90	0.44
1:D:183:PHE:HE2	1:D:197:MET:HB2	1.82	0.44
1:A:202:THR:HG23	1:A:202:THR:O	2.16	0.44
1:B:128:LEU:O	1:B:131:SER:OG	2.28	0.44
1:D:200:ALA:HB2	1:D:245:PHE:CZ	2.53	0.44
1:C:116:VAL:O	1:C:116:VAL:HG13	2.17	0.44
1:A:121:GLY:O	1:A:122:THR:HG22	2.18	0.44
1:B:272:THR:OG1	1:B:294:ARG:HG3	2.18	0.44
1:A:200:ALA:HB2	1:A:245:PHE:CZ	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:VAL:HG12	1:A:252:PRO:N	2.34	0.43
1:A:72:HIS:O	1:A:74:VAL:HG23	2.18	0.43
1:B:57:VAL:HG22	1:B:117:PHE:CE1	2.54	0.43
1:B:218:THR:HG22	1:B:256:ARG:HB2	2.00	0.43
1:C:64:GLN:O	1:C:81:PHE:HZ	2.01	0.43
1:C:218:THR:HA	1:C:229:SER:O	2.19	0.43
1:B:200:ALA:HB2	1:B:245:PHE:HZ	1.83	0.43
1:A:31:TRP:HB2	1:A:115:LEU:O	2.19	0.43
1:C:122:THR:CB	1:C:127:ASP:OD2	2.63	0.43
1:B:165:GLY:O	1:B:167:LEU:HD12	2.19	0.42
1:D:188:CYS:HB2	1:D:192:TYR:CZ	2.54	0.42
1:A:59:MET:CE	1:A:174:PRO:HG2	2.47	0.42
1:B:279:TYR:CE2	1:B:281:CYS:SG	3.12	0.42
1:C:182:LEU:HG	1:C:183:PHE:N	2.35	0.42
1:C:237:LYS:O	1:C:238:THR:HG23	2.19	0.42
1:C:79:CYS:HB3	1:C:93:LEU:HD22	2.00	0.42
1:A:52:LEU:HD12	1:A:144:ASN:O	2.18	0.42
1:D:200:ALA:HB2	1:D:245:PHE:HZ	1.83	0.42
1:B:218:THR:O	1:B:218:THR:HG23	2.19	0.42
1:C:267:ASP:O	1:C:270:SER:O	2.38	0.42
1:C:166:TYR:CE1	1:C:291:PHE:CZ	3.07	0.42
1:D:29:PRO:HA	1:D:30:THR:HA	1.86	0.42
1:C:15:TRP:CH2	1:C:17:ASN:CG	2.98	0.42
1:D:10:PHE:CD2	1:D:152:ILE:HG22	2.55	0.42
1:B:172:VAL:O	1:B:172:VAL:HG23	2.20	0.42
1:C:40:ASP:OD1	1:C:42:THR:HG22	2.20	0.42
1:D:185:ALA:HB2	1:D:197:MET:HE2	2.01	0.42
1:B:200:ALA:HB2	1:B:245:PHE:CZ	2.54	0.41
1:B:239:CYS:SG	1:B:244:ILE:HD12	2.60	0.41
1:A:46:PRO:HB3	1:A:99:ASN:HA	2.02	0.41
1:B:26:PRO:HB2	1:B:117:PHE:HB2	2.02	0.41
1:D:185:ALA:HB2	1:D:197:MET:CE	2.50	0.41
1:A:57:VAL:O	1:A:91:SER:OG	2.31	0.41
1:A:135:THR:O	1:A:139:ASN:ND2	2.46	0.41
1:C:17:ASN:CG	1:C:19:ALA:O	2.64	0.41
1:A:118:ASN:O	1:A:130:ASP:HB3	2.20	0.41
1:A:128:LEU:CD1	1:B:232:SER:OG	2.69	0.41
1:C:252:PRO:HD2	1:C:255:TYR:HB2	2.02	0.41
1:D:116:VAL:O	1:D:116:VAL:HG13	2.19	0.41
1:B:182:LEU:HG	1:B:183:PHE:N	2.36	0.41
1:C:271:TYR:CD1	1:C:271:TYR:C	2.99	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:GLY:O	1:A:122:THR:CG2	2.69	0.40
1:B:183:PHE:HD2	1:B:197:MET:HE3	1.76	0.40
1:C:135:THR:O	1:C:139:ASN:ND2	2.46	0.40
1:C:79:CYS:HA	1:C:95:CYS:HA	2.03	0.40
1:C:81:PHE:HB3	1:C:93:LEU:HD23	2.04	0.40
1:A:280:THR:OG1	1:A:286:TRP:HD1	2.04	0.40
1:D:182:LEU:HG	1:D:183:PHE:N	2.35	0.40
1:A:293:LEU:CD2	1:A:295:TRP:CZ2	3.04	0.40
1:D:29:PRO:HB3	1:D:61:THR:HG21	2.03	0.40
1:A:201:ASN:CB	1:A:206:VAL:HB	2.51	0.40
1:C:29:PRO:HB3	1:C:61:THR:HG21	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:GLY:O	1:C:283:GLY:CA[2_545]	1.89	0.31
1:B:284:GLY:O	1:C:283:GLY:C[2_545]	2.06	0.14

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	311/314 (99%)	295 (95%)	16 (5%)	0	100	100
1	B	311/314 (99%)	287 (92%)	24 (8%)	0	100	100
1	C	311/314 (99%)	289 (93%)	22 (7%)	0	100	100
1	D	311/314 (99%)	295 (95%)	16 (5%)	0	100	100
All	All	1244/1256 (99%)	1166 (94%)	78 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	249/264 (94%)	246 (99%)	3 (1%)	63	82
1	B	249/264 (94%)	247 (99%)	2 (1%)	73	86
1	C	250/264 (95%)	247 (99%)	3 (1%)	63	82
1	D	255/264 (97%)	251 (98%)	4 (2%)	55	79
All	All	1003/1056 (95%)	991 (99%)	12 (1%)	63	82

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	THR
1	A	208	ILE
1	A	244	ILE
1	B	182	LEU
1	B	208	ILE
1	C	74	VAL
1	C	172	VAL
1	C	182	LEU
1	D	54	MET
1	D	172	VAL
1	D	182	LEU
1	D	308	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	207	GLN
1	A	269	ASN
1	A	287	GLN
1	B	72	HIS
1	B	214	HIS
1	B	225	ASN
1	B	242	ASN

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Mol	Chain	Res	Type
1	D	214	HIS

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	313/314 (99%)	-0.92	0 100 100	24, 42, 58, 69	0
1	B	313/314 (99%)	-0.91	0 100 100	27, 39, 70, 93	0
1	C	313/314 (99%)	-0.88	0 100 100	33, 50, 75, 106	0
1	D	313/314 (99%)	-0.87	0 100 100	30, 41, 69, 81	0
All	All	1252/1256 (99%)	-0.89	0 100 100	24, 43, 70, 106	0

There are no RSRZ outliers to report.

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.