



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:28 AM UTC

PDB ID : 3IFP / pdb_00003ifp
Title : X-ray structure of amyloid beta peptide:antibody (Abeta1-7:12B4) complex
Authors : Weis, W.I.; Feinberg, H.; Basi, G.S.; Schenk, D.
Deposited on : 2009-07-24
Resolution : 2.95 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:

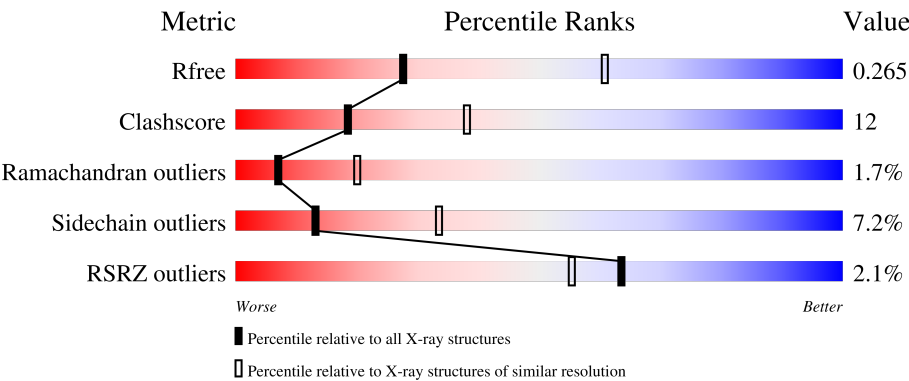
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	226	
1	C	226	
1	E	226	
1	H	226	
2	B	219	

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Mol	Chain	Length	Quality of chain
2	D	219	% 73% 25%
2	F	219	2% 73% 26%
2	L	219	3% 70% 28%
3	P	7	71% 14% 14%
3	Q	7	14% 71% 14% 14%
3	R	7	71% 14% 14%
3	S	7	71% 29%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13667 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 12B4 FAB antibody heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	222	Total	C	N	O	S	0	0	0
			1702	1079	281	336	6			
1	A	216	Total	C	N	O	S	0	0	0
			1662	1057	273	326	6			
1	C	216	Total	C	N	O	S	0	0	0
			1662	1057	273	326	6			
1	E	216	Total	C	N	O	S	0	0	0
			1662	1057	273	326	6			

- Molecule 2 is a protein called 12B4 FAB antibody light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	218	Total	C	N	O	S	0	0	0
			1692	1058	287	341	6			
2	B	218	Total	C	N	O	S	0	0	0
			1692	1058	287	341	6			
2	D	218	Total	C	N	O	S	0	0	0
			1692	1058	287	341	6			
2	F	218	Total	C	N	O	S	0	0	0
			1692	1058	287	341	6			

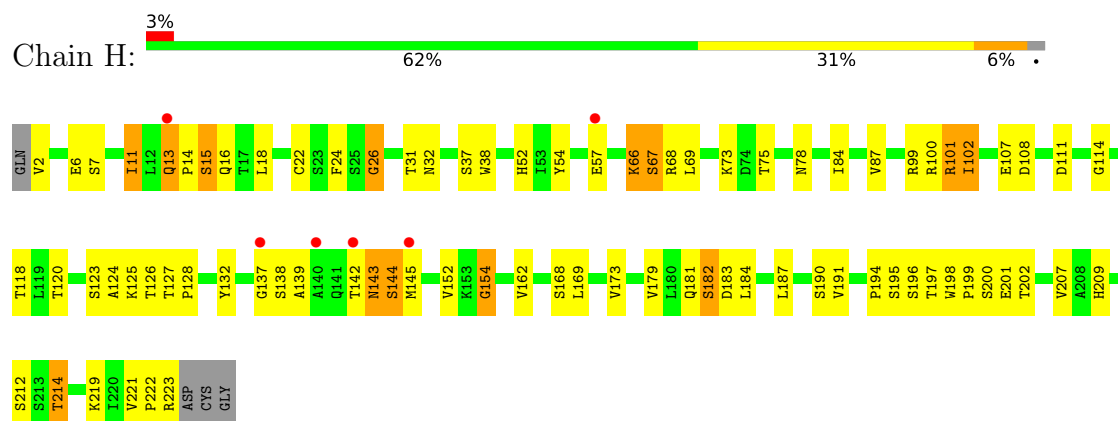
- Molecule 3 is a protein called Amyloid beta A4 protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	6	Total	C	N	O	0	0	0
			55	33	11	11			
3	Q	6	Total	C	N	O	0	0	0
			55	33	11	11			
3	R	6	Total	C	N	O	0	0	0
			55	33	11	11			
3	S	5	Total	C	N	O	0	0	0
			46	29	10	7			

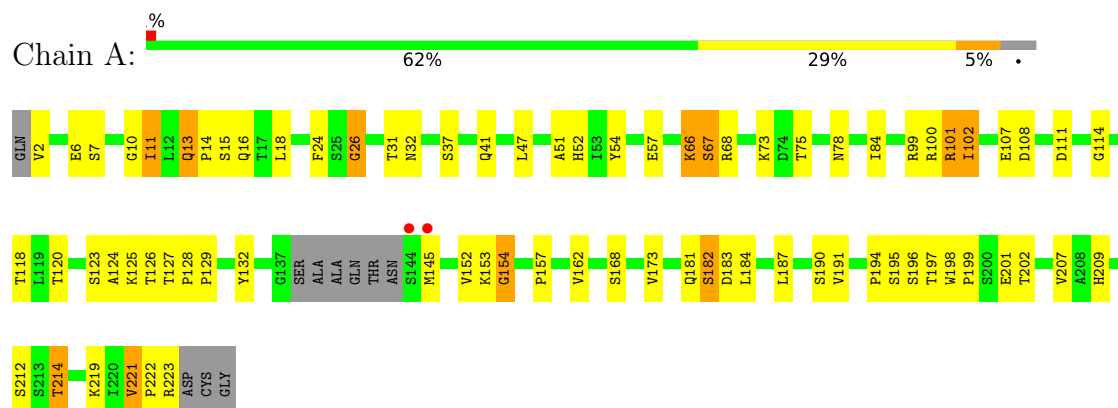
3 Residue-property plots [i](#)

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.

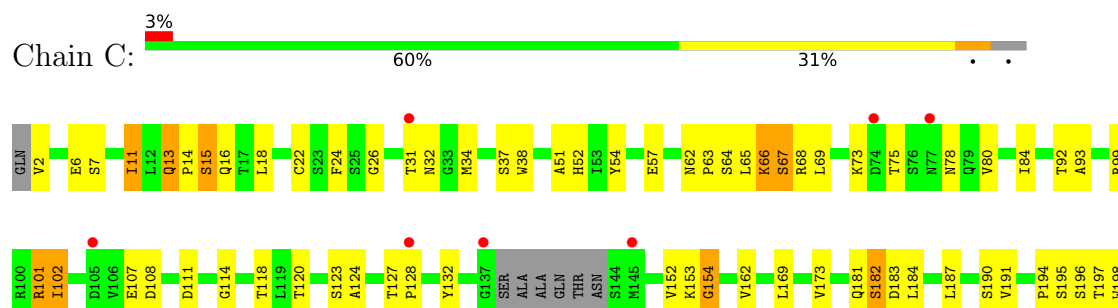
• Molecule 1: 12B4 FAB antibody heavy chain



• Molecule 1: 12B4 FAB antibody heavy chain

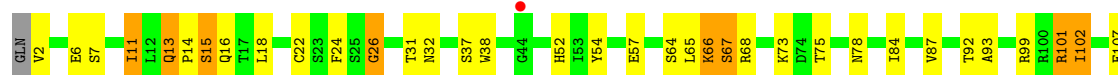


• Molecule 1: 12B4 FAB antibody heavy chain

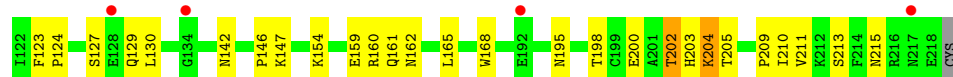




• Molecule 1: 12B4 FAB antibody heavy chain



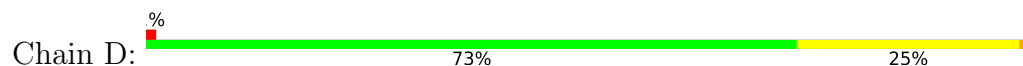
• Molecule 2: 12B4 FAB antibody light chain



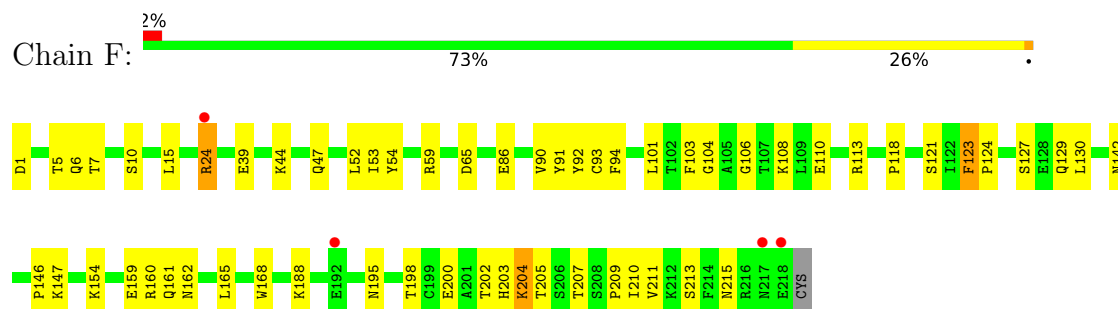
• Molecule 2: 12B4 FAB antibody light chain



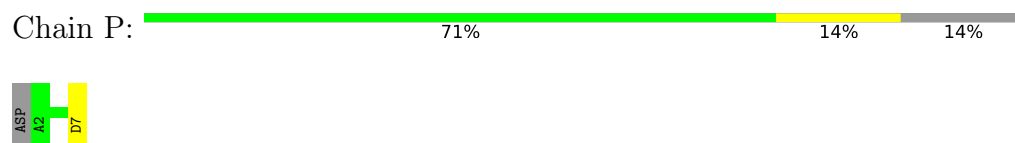
• Molecule 2: 12B4 FAB antibody light chain



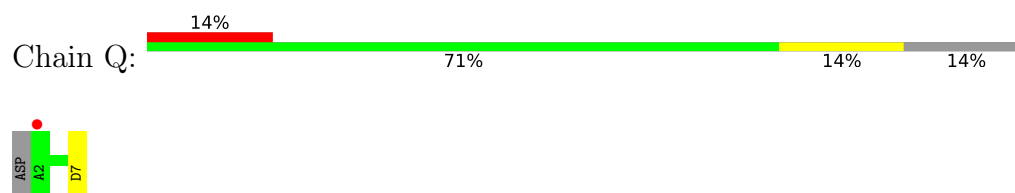
- Molecule 2: 12B4 FAB antibody light chain



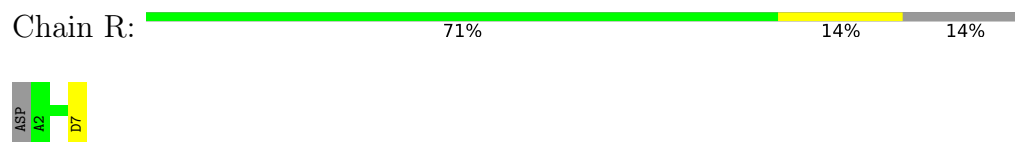
- Molecule 3: Amyloid beta A4 protein



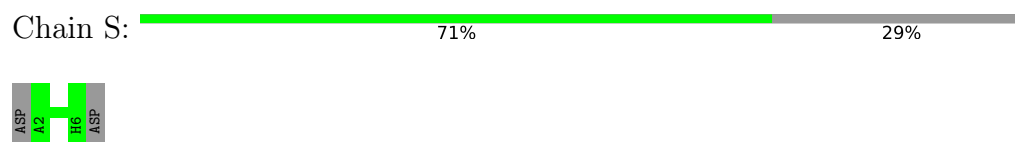
- Molecule 3: Amyloid beta A4 protein



- Molecule 3: Amyloid beta A4 protein



- Molecule 3: Amyloid beta A4 protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	78.86Å 79.24Å 94.10Å 68.66° 65.28° 78.47°	Depositor
Resolution (Å)	81.13 – 2.95 81.13 – 2.95	Depositor EDS
% Data completeness (in resolution range)	96.0 (81.13-2.95) 96.0 (81.13-2.95)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.96Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.235 , 0.269 0.231 , 0.265	Depositor DCC
R_{free} test set	1945 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.768	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 31.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.007 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	13667	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

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	0.58	0/1706	0.90	3/2334 (0.1%)
1	C	0.53	0/1706	0.88	2/2334 (0.1%)
1	E	0.54	0/1706	0.90	3/2334 (0.1%)
1	H	0.57	0/1747	0.90	3/2392 (0.1%)
2	B	0.57	0/1730	0.90	1/2348 (0.0%)
2	D	0.55	0/1730	0.89	4/2348 (0.2%)
2	F	0.57	0/1730	0.91	3/2348 (0.1%)
2	L	0.56	0/1730	0.90	2/2348 (0.1%)
3	P	0.42	0/56	0.65	0/72
3	Q	0.44	0/56	0.64	0/72
3	R	0.45	0/56	0.58	0/72
3	S	0.46	0/47	0.58	0/61
All	All	0.56	0/14000	0.89	21/19063 (0.1%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	123	PHE	CA-C-N	5.97	124.02	119.66
2	F	123	PHE	C-N-CA	5.97	124.02	119.66
2	D	142	ASN	N-CA-C	5.92	118.61	109.07
1	H	209	HIS	CA-C-N	5.85	125.32	119.24
1	H	209	HIS	C-N-CA	5.85	125.32	119.24
2	L	142	ASN	N-CA-C	5.77	118.30	108.90
1	A	209	HIS	CA-C-N	5.73	125.20	119.24
1	A	209	HIS	C-N-CA	5.73	125.20	119.24
2	D	123	PHE	CA-C-N	5.71	123.83	119.66
2	D	123	PHE	C-N-CA	5.71	123.83	119.66
1	E	209	HIS	CA-C-N	5.53	124.99	119.24
1	E	209	HIS	C-N-CA	5.53	124.99	119.24
1	C	209	HIS	CA-C-N	5.52	124.98	119.24
1	C	209	HIS	C-N-CA	5.52	124.98	119.24
2	F	142	ASN	N-CA-C	5.46	117.87	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	26	GLY	N-CA-C	-5.46	108.20	114.69
1	E	26	GLY	N-CA-C	-5.42	108.24	114.69
2	B	142	ASN	N-CA-C	5.20	117.97	109.59
1	A	26	GLY	N-CA-C	-5.04	108.69	114.69
2	D	162	ASN	N-CA-C	5.04	117.82	109.46
2	L	102	THR	N-CA-C	5.02	117.01	109.23

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	1662	0	1631	48	2
1	C	1662	0	1631	50	0
1	E	1662	0	1631	44	0
1	H	1702	0	1668	50	2
2	B	1692	0	1636	37	0
2	D	1692	0	1636	36	0
2	F	1692	0	1636	35	0
2	L	1692	0	1636	37	0
3	P	55	0	43	0	0
3	Q	55	0	43	0	0
3	R	55	0	43	0	0
3	S	46	0	39	0	0
All	All	13667	0	13273	322	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 12.

All (322) 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)
2:F:6:GLN:NE2	2:F:93:CYS:H	1.80	0.78
1:C:212:SER:OG	1:C:214:THR:HG23	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:6:GLN:NE2	2:B:93:CYS:H	1.82	0.77
2:D:6:GLN:NE2	2:D:93:CYS:H	1.82	0.76
2:L:6:GLN:NE2	2:L:93:CYS:H	1.83	0.76
1:H:212:SER:OG	1:H:214:THR:HG23	1.88	0.74
1:E:212:SER:OG	1:E:214:THR:HG23	1.87	0.73
1:A:37:SER:HB3	1:A:52:HIS:ND1	2.05	0.72
1:H:37:SER:HB3	1:H:52:HIS:ND1	2.05	0.72
1:H:142:THR:OG1	1:H:143:ASN:HB2	1.89	0.72
2:B:146:PRO:HG2	2:B:204:LYS:HE3	1.72	0.71
1:A:212:SER:OG	1:A:214:THR:HG23	1.91	0.71
2:F:203:HIS:CE1	2:F:205:THR:HG23	2.26	0.70
1:H:173:VAL:HG22	1:H:191:VAL:HG23	1.74	0.70
2:D:203:HIS:CE1	2:D:205:THR:HG23	2.26	0.70
1:H:202:THR:HB	1:H:219:LYS:HE3	1.74	0.69
1:C:37:SER:HB3	1:C:52:HIS:ND1	2.08	0.69
1:E:37:SER:HB3	1:E:52:HIS:ND1	2.08	0.69
1:E:202:THR:HB	1:E:219:LYS:HE3	1.74	0.69
2:B:203:HIS:CE1	2:B:205:THR:HG23	2.29	0.68
2:B:6:GLN:HE22	2:B:92:TYR:HA	1.59	0.68
2:F:146:PRO:HG2	2:F:204:LYS:HE3	1.74	0.67
1:H:137:GLY:O	1:H:139:ALA:N	2.27	0.67
2:B:44:LYS:HB2	2:B:47:GLN:HG3	1.76	0.67
1:C:124:ALA:HB2	1:C:183:ASP:HB3	1.75	0.67
1:A:173:VAL:HG22	1:A:191:VAL:HG23	1.77	0.67
2:L:203:HIS:CE1	2:L:205:THR:HG23	2.30	0.67
1:A:124:ALA:HB2	1:A:183:ASP:HB3	1.77	0.67
2:F:44:LYS:HB2	2:F:47:GLN:HG3	1.76	0.67
2:L:146:PRO:HG2	2:L:204:LYS:HE3	1.74	0.67
2:D:6:GLN:HE22	2:D:92:TYR:HA	1.60	0.66
2:L:44:LYS:HB2	2:L:47:GLN:HG3	1.77	0.66
1:C:202:THR:HB	1:C:219:LYS:HE3	1.77	0.66
2:D:44:LYS:HB2	2:D:47:GLN:HG3	1.77	0.66
1:E:187:LEU:HD12	1:E:187:LEU:C	2.20	0.66
1:A:162:VAL:HG22	1:A:207:VAL:HG22	1.77	0.66
2:D:146:PRO:HG2	2:D:204:LYS:HE3	1.76	0.65
1:A:187:LEU:C	1:A:187:LEU:HD12	2.22	0.65
1:H:187:LEU:C	1:H:187:LEU:HD12	2.22	0.65
1:C:187:LEU:C	1:C:187:LEU:HD12	2.21	0.65
2:F:6:GLN:HE22	2:F:92:TYR:HA	1.62	0.65
1:H:162:VAL:HG22	1:H:207:VAL:HG22	1.78	0.64
1:A:202:THR:HB	1:A:219:LYS:HE3	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:124:ALA:HB2	1:E:183:ASP:HB3	1.79	0.64
1:E:173:VAL:HG22	1:E:191:VAL:HG23	1.79	0.64
2:D:39:GLU:HG2	2:D:54:TYR:HA	1.79	0.64
1:E:162:VAL:HG22	1:E:207:VAL:HG22	1.80	0.64
2:L:6:GLN:HE22	2:L:92:TYR:HA	1.63	0.64
1:A:66:LYS:O	1:A:67:SER:HB2	1.99	0.63
2:F:39:GLU:HG2	2:F:54:TYR:HA	1.80	0.63
1:C:173:VAL:HG22	1:C:191:VAL:HG23	1.80	0.63
1:H:66:LYS:O	1:H:67:SER:HB2	1.99	0.63
1:C:162:VAL:HG22	1:C:207:VAL:HG22	1.81	0.63
1:E:66:LYS:O	1:E:67:SER:HB2	1.99	0.63
2:B:39:GLU:HG2	2:B:54:TYR:HA	1.82	0.61
1:C:197:THR:O	1:C:201:GLU:HB2	2.00	0.61
1:E:132:TYR:CZ	2:F:129:GLN:HG3	2.36	0.61
1:C:132:TYR:CZ	2:D:129:GLN:HG3	2.36	0.60
1:E:2:VAL:HA	1:E:26:GLY:HA3	1.82	0.60
2:B:90:VAL:HG22	2:B:108:LYS:HD2	1.83	0.60
1:E:197:THR:O	1:E:201:GLU:HB2	2.01	0.60
1:H:124:ALA:HB2	1:H:183:ASP:HB3	1.84	0.60
1:H:197:THR:O	1:H:201:GLU:HB2	2.01	0.60
1:A:197:THR:O	1:A:201:GLU:HB2	2.01	0.59
2:L:39:GLU:HG2	2:L:54:TYR:HA	1.84	0.59
1:A:2:VAL:HA	1:A:26:GLY:HA3	1.84	0.59
1:C:2:VAL:HA	1:C:26:GLY:HA3	1.83	0.59
1:H:132:TYR:CZ	2:L:129:GLN:HG3	2.38	0.58
1:C:57:GLU:HG3	1:C:73:LYS:HE3	1.86	0.58
1:H:2:VAL:HA	1:H:26:GLY:HA3	1.86	0.58
1:C:101:ARG:O	1:C:108:ASP:HA	2.04	0.57
1:C:66:LYS:O	1:C:67:SER:HB2	2.05	0.57
1:H:142:THR:HA	1:H:143:ASN:CB	2.35	0.56
1:A:132:TYR:CZ	2:B:129:GLN:HG3	2.40	0.56
1:E:101:ARG:O	1:E:108:ASP:HA	2.05	0.56
1:C:222:PRO:O	1:C:223:ARG:HB2	2.05	0.56
1:E:68:ARG:O	1:E:84:ILE:HA	2.07	0.55
1:E:57:GLU:HG3	1:E:73:LYS:HE3	1.89	0.55
1:H:57:GLU:HG3	1:H:73:LYS:HE3	1.88	0.55
1:A:68:ARG:O	1:A:84:ILE:HA	2.07	0.55
1:A:67:SER:HB2	2:D:189:ASP:OD1	2.07	0.54
1:A:11:ILE:HD11	1:A:157:PRO:HB3	1.89	0.54
2:D:90:VAL:HG22	2:D:108:LYS:HD2	1.89	0.54
2:F:90:VAL:HG22	2:F:108:LYS:HD2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:ARG:O	1:C:84:ILE:HA	2.08	0.54
1:E:222:PRO:O	1:E:223:ARG:HB2	2.08	0.53
1:H:101:ARG:O	1:H:108:ASP:HA	2.08	0.53
1:A:101:ARG:O	1:A:108:ASP:HA	2.09	0.53
1:E:54:TYR:HE2	1:E:102:ILE:HD13	1.73	0.52
2:D:154:LYS:HE3	2:D:200:GLU:OE2	2.09	0.52
2:L:90:VAL:HG22	2:L:108:LYS:HD2	1.92	0.52
1:A:57:GLU:HG3	1:A:73:LYS:HE3	1.91	0.52
1:A:18:LEU:HD23	1:A:84:ILE:HD12	1.92	0.51
2:B:195:ASN:HD21	2:B:215:ASN:HB3	1.75	0.51
2:F:203:HIS:ND1	2:F:205:THR:HG23	2.24	0.51
2:D:159:GLU:HG3	2:D:160:ARG:N	2.25	0.51
2:D:203:HIS:ND1	2:D:205:THR:HG23	2.24	0.51
2:B:59:ARG:CD	2:B:65:ASP:HA	2.41	0.51
1:C:54:TYR:HE2	1:C:102:ILE:HD13	1.76	0.51
1:H:68:ARG:O	1:H:84:ILE:HA	2.10	0.50
1:C:99:ARG:NH2	1:C:111:ASP:OD2	2.43	0.50
1:H:222:PRO:O	1:H:223:ARG:HB2	2.11	0.50
2:L:94:PHE:HB2	2:L:103:PHE:CD2	2.46	0.50
1:A:154:GLY:HA2	1:A:184:LEU:HB3	1.94	0.50
1:H:101:ARG:CZ	1:H:107:GLU:HB3	2.42	0.50
2:B:91:TYR:O	2:B:106:GLY:HA2	2.11	0.50
2:F:154:LYS:HE3	2:F:200:GLU:OE2	2.11	0.50
2:D:59:ARG:CD	2:D:65:ASP:HA	2.42	0.50
1:H:18:LEU:HD23	1:H:84:ILE:HD12	1.94	0.50
1:E:99:ARG:HH21	1:E:111:ASP:CG	2.20	0.50
2:L:6:GLN:NE2	2:L:104:GLY:HA3	2.26	0.49
2:B:159:GLU:HG3	2:B:160:ARG:N	2.26	0.49
2:L:159:GLU:HG3	2:L:160:ARG:N	2.26	0.49
2:F:154:LYS:HB2	2:F:198:THR:HB	1.94	0.49
2:L:108:LYS:HG2	2:L:110:GLU:CD	2.37	0.49
2:B:154:LYS:HB2	2:B:198:THR:HB	1.93	0.49
2:B:154:LYS:HE3	2:B:200:GLU:OE2	2.13	0.49
1:C:99:ARG:HH21	1:C:111:ASP:CG	2.20	0.49
2:F:159:GLU:HG3	2:F:160:ARG:N	2.27	0.49
2:L:154:LYS:HB2	2:L:198:THR:HB	1.93	0.49
2:B:52:LEU:C	2:B:53:ILE:HD12	2.38	0.49
1:H:142:THR:CA	1:H:143:ASN:HB2	2.43	0.49
2:B:198:THR:HG23	2:B:211:VAL:HG13	1.94	0.49
2:L:52:LEU:C	2:L:53:ILE:HD12	2.38	0.48
2:F:198:THR:HG23	2:F:211:VAL:HG13	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:59:ARG:CD	2:L:65:ASP:HA	2.43	0.48
2:L:154:LYS:HE3	2:L:200:GLU:OE2	2.13	0.48
1:H:181:GLN:O	1:H:182:SER:CB	2.61	0.48
2:L:198:THR:HG23	2:L:211:VAL:HG13	1.95	0.48
1:E:181:GLN:O	1:E:182:SER:CB	2.60	0.48
2:D:52:LEU:C	2:D:53:ILE:HD12	2.39	0.48
1:E:132:TYR:CE2	2:F:129:GLN:HG3	2.49	0.48
2:B:161:GLN:H	2:B:161:GLN:CD	2.22	0.48
1:C:132:TYR:CE2	2:D:129:GLN:HG3	2.49	0.48
2:D:94:PHE:HB2	2:D:103:PHE:CD2	2.49	0.48
2:F:59:ARG:CD	2:F:65:ASP:HA	2.44	0.48
2:L:165:LEU:HA	2:L:165:LEU:HD23	1.68	0.47
2:D:41:TYR:CZ	2:D:51:LEU:HD13	2.48	0.47
2:D:91:TYR:O	2:D:106:GLY:HA2	2.14	0.47
1:E:99:ARG:NH2	1:E:111:ASP:OD2	2.45	0.47
1:A:152:VAL:HG22	1:A:207:VAL:HG21	1.96	0.47
1:C:154:GLY:HA2	1:C:184:LEU:HB3	1.96	0.47
1:E:101:ARG:CZ	1:E:107:GLU:HB3	2.43	0.47
2:L:195:ASN:HD21	2:L:215:ASN:HB3	1.80	0.47
1:A:11:ILE:HD11	1:A:157:PRO:CB	2.44	0.47
2:B:203:HIS:ND1	2:B:205:THR:HG23	2.29	0.47
1:H:31:THR:O	1:H:32:ASN:C	2.56	0.47
1:A:67:SER:HB3	2:D:189:ASP:OD2	2.15	0.47
1:A:181:GLN:O	1:A:182:SER:CB	2.62	0.47
2:F:52:LEU:C	2:F:53:ILE:HD12	2.40	0.47
2:F:195:ASN:HD21	2:F:215:ASN:HB3	1.80	0.47
2:L:6:GLN:HE21	2:L:104:GLY:HA3	1.80	0.47
2:L:91:TYR:O	2:L:106:GLY:HA2	2.14	0.47
1:C:18:LEU:HD23	1:C:84:ILE:HD12	1.96	0.47
2:D:195:ASN:HD21	2:D:215:ASN:HB3	1.79	0.47
1:E:24:PHE:CZ	1:E:78:ASN:HA	2.50	0.47
1:H:173:VAL:CG2	1:H:191:VAL:HG23	2.42	0.47
1:A:99:ARG:NH2	1:A:111:ASP:OD2	2.46	0.47
2:F:91:TYR:O	2:F:106:GLY:HA2	2.15	0.47
2:D:161:GLN:H	2:D:161:GLN:CD	2.22	0.46
1:A:41:GLN:HB2	1:A:47:LEU:HD23	1.98	0.46
2:D:198:THR:HG23	2:D:211:VAL:HG13	1.97	0.46
1:C:11:ILE:HA	1:C:120:THR:O	2.16	0.46
1:E:14:PRO:O	1:E:15:SER:CB	2.63	0.46
1:A:66:LYS:O	1:A:67:SER:CB	2.64	0.46
2:B:94:PHE:HB2	2:B:103:PHE:CD2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:PRO:O	1:C:15:SER:CB	2.63	0.46
1:H:99:ARG:HH21	1:H:111:ASP:CG	2.24	0.46
1:A:222:PRO:O	1:A:223:ARG:HB2	2.15	0.46
1:C:127:THR:HA	1:C:128:PRO:HD3	1.78	0.46
2:F:94:PHE:HB2	2:F:103:PHE:CD2	2.51	0.46
1:C:24:PHE:CZ	1:C:78:ASN:HA	2.51	0.46
2:D:154:LYS:HB2	2:D:198:THR:HB	1.98	0.46
1:A:173:VAL:CG2	1:A:191:VAL:HG23	2.44	0.46
1:E:11:ILE:HA	1:E:120:THR:O	2.16	0.46
1:H:154:GLY:HA2	1:H:184:LEU:HB3	1.98	0.46
1:A:99:ARG:HH21	1:A:111:ASP:CG	2.24	0.46
1:C:173:VAL:CG2	1:C:191:VAL:HG23	2.46	0.46
1:C:181:GLN:O	1:C:182:SER:CB	2.63	0.46
1:H:181:GLN:HG3	2:L:165:LEU:CD1	2.46	0.45
1:H:11:ILE:HA	1:H:120:THR:O	2.16	0.45
1:H:66:LYS:O	1:H:67:SER:CB	2.65	0.45
1:A:24:PHE:CZ	1:A:78:ASN:HA	2.51	0.45
1:C:101:ARG:CZ	1:C:107:GLU:HB3	2.45	0.45
1:E:22:CYS:HB2	1:E:38:TRP:CZ2	2.51	0.45
1:H:13:GLN:HG3	1:H:16:GLN:NE2	2.31	0.45
1:E:18:LEU:HD23	1:E:84:ILE:HD12	1.97	0.45
1:A:54:TYR:HE2	1:A:102:ILE:HD13	1.80	0.45
2:L:202:THR:CG2	2:L:209:PRO:HG3	2.46	0.45
1:C:101:ARG:H	1:C:101:ARG:HG2	1.58	0.45
2:F:161:GLN:HE21	2:F:162:ASN:H	1.64	0.45
2:F:94:PHE:CD2	2:F:94:PHE:C	2.95	0.45
1:H:143:ASN:O	1:H:144:SER:HB3	2.17	0.45
1:A:11:ILE:HA	1:A:120:THR:O	2.17	0.45
2:L:130:LEU:HD23	2:L:130:LEU:HA	1.81	0.45
2:D:161:GLN:HG2	2:D:162:ASN:N	2.32	0.45
2:F:123:PHE:HA	2:F:124:PRO:HD2	1.75	0.45
2:L:203:HIS:ND1	2:L:205:THR:HG23	2.32	0.44
1:A:132:TYR:CE2	2:B:129:GLN:HG3	2.52	0.44
2:B:108:LYS:HG2	2:B:110:GLU:CD	2.42	0.44
1:A:31:THR:O	1:A:32:ASN:C	2.60	0.44
1:A:37:SER:HB2	1:A:51:ALA:O	2.18	0.44
2:B:59:ARG:HD3	2:B:65:ASP:HA	1.99	0.44
2:B:111:LEU:HB2	2:B:171:GLN:HE22	1.81	0.44
1:C:153:LYS:NZ	1:C:181:GLN:HE22	2.16	0.44
1:E:173:VAL:CG2	1:E:191:VAL:HG23	2.47	0.44
1:H:54:TYR:HE2	1:H:102:ILE:HD13	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:123:PHE:HA	2:B:124:PRO:HD2	1.77	0.44
2:B:165:LEU:HA	2:B:165:LEU:HD23	1.66	0.44
1:E:31:THR:O	1:E:32:ASN:C	2.59	0.44
1:H:99:ARG:HG2	1:H:100:ARG:N	2.33	0.44
1:E:154:GLY:HA2	1:E:184:LEU:HB3	2.00	0.44
2:L:161:GLN:HG2	2:L:162:ASN:N	2.33	0.44
1:A:127:THR:HA	1:A:128:PRO:HD3	1.77	0.44
2:D:15:LEU:HD23	2:D:15:LEU:HA	1.83	0.44
2:D:130:LEU:HD23	2:D:130:LEU:HA	1.82	0.44
1:H:24:PHE:CZ	1:H:78:ASN:HA	2.53	0.44
2:L:121:SER:HB2	2:L:123:PHE:CE2	2.53	0.44
1:A:181:GLN:HG3	2:B:165:LEU:CD1	2.48	0.43
1:A:198:TRP:CG	1:A:199:PRO:HA	2.53	0.43
1:C:22:CYS:HB2	1:C:38:TRP:CZ2	2.53	0.43
1:E:153:LYS:NZ	1:E:181:GLN:HE22	2.16	0.43
1:H:14:PRO:O	1:H:15:SER:CB	2.65	0.43
2:L:59:ARG:HD3	2:L:65:ASP:HA	2.00	0.43
2:L:161:GLN:CD	2:L:161:GLN:H	2.27	0.43
2:F:108:LYS:HG2	2:F:110:GLU:CD	2.44	0.43
1:E:54:TYR:CE2	1:E:102:ILE:HD13	2.53	0.43
2:F:161:GLN:HG2	2:F:162:ASN:N	2.33	0.43
1:A:99:ARG:HG2	1:A:100:ARG:N	2.33	0.43
1:A:194:PRO:C	1:A:196:SER:H	2.27	0.43
2:D:123:PHE:HA	2:D:124:PRO:HD2	1.77	0.43
1:A:6:GLU:OE2	1:A:114:GLY:HA3	2.18	0.43
2:B:189:ASP:OD1	1:C:67:SER:HB2	2.18	0.43
1:C:31:THR:O	1:C:32:ASN:C	2.59	0.43
2:F:24:ARG:CZ	2:F:24:ARG:HB3	2.49	0.43
2:D:54:TYR:CE1	2:D:58:ASN:HB3	2.54	0.43
1:H:132:TYR:CE2	2:L:129:GLN:HG3	2.54	0.42
1:H:198:TRP:CG	1:H:199:PRO:HA	2.54	0.42
2:L:123:PHE:HA	2:L:124:PRO:HD2	1.76	0.42
2:L:202:THR:HG22	2:L:209:PRO:HG3	2.01	0.42
1:A:101:ARG:CZ	1:A:107:GLU:HB3	2.49	0.42
1:C:199:PRO:O	1:C:200:SER:C	2.62	0.42
1:E:187:LEU:C	1:E:187:LEU:CD1	2.89	0.42
1:H:142:THR:CA	1:H:143:ASN:CB	2.97	0.42
1:H:194:PRO:C	1:H:196:SER:H	2.28	0.42
1:H:199:PRO:O	1:H:200:SER:C	2.61	0.42
1:E:6:GLU:OE2	1:E:114:GLY:HA3	2.18	0.42
1:E:198:TRP:CG	1:E:199:PRO:HA	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:69:LEU:CD2	1:H:84:ILE:HG12	2.49	0.42
2:L:118:PRO:HD2	2:L:210:ILE:CD1	2.49	0.42
1:H:99:ARG:NH2	1:H:111:ASP:OD2	2.47	0.42
2:L:38:LEU:HD13	2:L:76:PHE:CD2	2.55	0.42
2:F:118:PRO:HD2	2:F:210:ILE:CD1	2.49	0.42
2:B:40:TRP:CZ3	2:B:93:CYS:HB3	2.54	0.42
2:D:94:PHE:CD2	2:D:94:PHE:C	2.98	0.42
1:C:54:TYR:CE2	1:C:102:ILE:HD13	2.55	0.42
1:C:152:VAL:HG22	1:C:207:VAL:HG21	2.01	0.42
2:D:24:ARG:CZ	2:D:24:ARG:HB3	2.50	0.42
2:D:161:GLN:HE21	2:D:162:ASN:H	1.67	0.42
1:E:181:GLN:O	1:E:182:SER:HB2	2.20	0.42
1:A:14:PRO:O	1:A:15:SER:CB	2.67	0.42
1:A:153:LYS:NZ	1:A:181:GLN:HE22	2.18	0.42
2:B:6:GLN:NE2	2:B:104:GLY:HA3	2.33	0.42
1:A:221:VAL:HA	1:A:222:PRO:HD3	1.91	0.42
1:C:198:TRP:CG	1:C:199:PRO:HA	2.55	0.42
1:E:194:PRO:C	1:E:196:SER:H	2.26	0.42
2:F:6:GLN:NE2	2:F:104:GLY:HA3	2.35	0.42
2:F:161:GLN:CD	2:F:161:GLN:H	2.27	0.42
1:H:101:ARG:H	1:H:101:ARG:HG2	1.62	0.42
1:H:187:LEU:C	1:H:187:LEU:CD1	2.91	0.42
1:E:13:GLN:HG3	1:E:16:GLN:NE2	2.35	0.42
1:E:127:THR:HA	1:E:128:PRO:HD3	1.77	0.42
2:F:15:LEU:HD23	2:F:15:LEU:HA	1.80	0.42
1:A:13:GLN:HG3	1:A:16:GLN:NE2	2.35	0.41
2:D:6:GLN:NE2	2:D:104:GLY:HA3	2.34	0.41
1:C:13:GLN:HG3	1:C:16:GLN:NE2	2.35	0.41
1:C:187:LEU:C	1:C:187:LEU:CD1	2.90	0.41
2:D:108:LYS:HG2	2:D:110:GLU:CD	2.46	0.41
1:H:125:LYS:O	1:H:126:THR:C	2.63	0.41
1:A:187:LEU:C	1:A:187:LEU:CD1	2.91	0.41
1:C:181:GLN:HG3	2:D:165:LEU:CD1	2.50	0.41
2:B:24:ARG:CZ	2:B:24:ARG:HB3	2.49	0.41
1:E:64:SER:O	1:E:65:LEU:HD23	2.20	0.41
1:H:6:GLU:OE2	1:H:114:GLY:HA3	2.20	0.41
2:B:130:LEU:HD23	2:B:130:LEU:HA	1.84	0.41
1:C:62:ASN:HA	1:C:63:PRO:HD3	1.85	0.41
1:C:64:SER:O	1:C:65:LEU:HD23	2.21	0.41
1:E:92:THR:O	1:E:93:ALA:HB2	2.20	0.41
1:A:128:PRO:HA	1:A:129:PRO:HD3	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:TRP:HE1	1:C:80:VAL:HG12	1.86	0.41
1:E:11:ILE:HD11	1:E:157:PRO:HB3	2.03	0.41
2:L:40:TRP:CZ3	2:L:93:CYS:HB3	2.55	0.41
2:L:41:TYR:CZ	2:L:51:LEU:HD13	2.56	0.41
2:B:54:TYR:CE1	2:B:58:ASN:HB3	2.55	0.41
2:B:59:ARG:HD3	2:B:67:PHE:O	2.20	0.41
2:B:195:ASN:O	2:B:216:ARG:N	2.51	0.41
2:D:59:ARG:HD3	2:D:65:ASP:HA	2.01	0.41
1:C:2:VAL:HG22	1:C:26:GLY:C	2.46	0.41
1:C:194:PRO:C	1:C:196:SER:H	2.27	0.41
2:D:59:ARG:HD3	2:D:67:PHE:O	2.21	0.41
1:E:181:GLN:HG3	2:F:165:LEU:CD1	2.51	0.41
2:L:15:LEU:HD23	2:L:15:LEU:HA	1.81	0.40
1:C:34:MET:SD	1:C:99:ARG:HD2	2.61	0.40
1:C:37:SER:HB2	1:C:51:ALA:O	2.21	0.40
1:C:92:THR:O	1:C:93:ALA:HB2	2.20	0.40
2:F:6:GLN:HE21	2:F:104:GLY:HA3	1.85	0.40
2:F:130:LEU:O	2:F:188:LYS:HD3	2.21	0.40
2:B:118:PRO:HD2	2:B:210:ILE:CD1	2.51	0.40
1:H:127:THR:HA	1:H:128:PRO:HD3	1.78	0.40
2:B:130:LEU:HD21	2:B:135:ALA:HB2	2.04	0.40
1:E:125:LYS:O	1:E:126:THR:C	2.64	0.40
2:F:6:GLN:HE22	2:F:93:CYS:H	1.64	0.40
2:F:59:ARG:HD3	2:F:65:ASP:HA	2.01	0.40
1:H:22:CYS:HB2	1:H:38:TRP:CZ2	2.56	0.40
1:H:152:VAL:HG22	1:H:207:VAL:HG21	2.03	0.40
1:H:221:VAL:HA	1:H:222:PRO:HD3	1.90	0.40
1:A:125:LYS:O	1:A:126:THR:C	2.63	0.40
1:C:6:GLU:OE2	1:C:114:GLY:HA3	2.21	0.40
1:C:69:LEU:CD2	1:C:84:ILE:HG12	2.52	0.40
1:E:199:PRO:O	1:E:200:SER:C	2.65	0.40
2:B:41:TYR:CZ	2:B:51:LEU:HD13	2.57	0.40
2:F:207:THR:O	2:F:209:PRO:HD3	2.22	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:H:126:THR:OG1	1:A:126:THR:OG1[1_554]	2.17	0.03
1:H:212:SER:O	1:A:10:GLY:O[1_554]	2.17	0.03

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	212/226 (94%)	191 (90%)	15 (7%)	6 (3%)	4	11
1	C	212/226 (94%)	192 (91%)	13 (6%)	7 (3%)	3	7
1	E	212/226 (94%)	191 (90%)	14 (7%)	7 (3%)	3	7
1	H	220/226 (97%)	195 (89%)	15 (7%)	10 (4%)	2	4
2	B	216/219 (99%)	203 (94%)	13 (6%)	0	100	100
2	D	216/219 (99%)	203 (94%)	13 (6%)	0	100	100
2	F	216/219 (99%)	205 (95%)	11 (5%)	0	100	100
2	L	216/219 (99%)	204 (94%)	12 (6%)	0	100	100
3	P	4/7 (57%)	4 (100%)	0	0	100	100
3	Q	4/7 (57%)	4 (100%)	0	0	100	100
3	R	4/7 (57%)	4 (100%)	0	0	100	100
3	S	3/7 (43%)	3 (100%)	0	0	100	100
All	All	1735/1808 (96%)	1599 (92%)	106 (6%)	30 (2%)	7	20

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	67	SER
1	H	138	SER
1	H	182	SER
1	A	67	SER
1	A	182	SER
1	C	67	SER
1	C	182	SER
1	E	67	SER
1	E	182	SER
1	H	66	LYS
1	H	143	ASN
1	H	144	SER

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Mol	Chain	Res	Type
1	H	154	GLY
1	A	66	LYS
1	A	154	GLY
1	C	66	LYS
1	C	154	GLY
1	E	66	LYS
1	E	154	GLY
1	H	11	ILE
1	A	11	ILE
1	C	11	ILE
1	C	15	SER
1	E	11	ILE
1	E	15	SER
1	H	15	SER
1	H	102	ILE
1	A	102	ILE
1	C	102	ILE
1	E	102	ILE

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	193/200 (96%)	181 (94%)	12 (6%)	16	39
1	C	193/200 (96%)	182 (94%)	11 (6%)	18	42
1	E	193/200 (96%)	180 (93%)	13 (7%)	15	36
1	H	197/200 (98%)	183 (93%)	14 (7%)	13	33
2	B	195/196 (100%)	180 (92%)	15 (8%)	12	30
2	D	195/196 (100%)	181 (93%)	14 (7%)	13	33
2	F	195/196 (100%)	180 (92%)	15 (8%)	12	30
2	L	195/196 (100%)	179 (92%)	16 (8%)	10	27
3	P	5/6 (83%)	4 (80%)	1 (20%)	1	3
3	Q	5/6 (83%)	4 (80%)	1 (20%)	1	3

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	R	5/6 (83%)	4 (80%)	1 (20%)	1	3
3	S	4/6 (67%)	4 (100%)	0	100	100
All	All	1575/1608 (98%)	1462 (93%)	113 (7%)	13	33

All (113) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	7	SER
1	H	13	GLN
1	H	75	THR
1	H	87	VAL
1	H	101	ARG
1	H	118	THR
1	H	123	SER
1	H	145	MET
1	H	168	SER
1	H	169	LEU
1	H	179	VAL
1	H	190	SER
1	H	195	SER
1	H	214	THR
2	L	1	ASP
2	L	5	THR
2	L	7	THR
2	L	10	SER
2	L	24	ARG
2	L	86	GLU
2	L	101	LEU
2	L	113	ARG
2	L	119	THR
2	L	121	SER
2	L	127	SER
2	L	147	LYS
2	L	168	TRP
2	L	202	THR
2	L	204	LYS
2	L	213	SER
3	P	7	ASP
1	A	7	SER
1	A	13	GLN
1	A	75	THR

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Mol	Chain	Res	Type
1	A	101	ARG
1	A	118	THR
1	A	123	SER
1	A	145	MET
1	A	168	SER
1	A	190	SER
1	A	195	SER
1	A	214	THR
1	A	221	VAL
2	B	1	ASP
2	B	5	THR
2	B	7	THR
2	B	10	SER
2	B	14	SER
2	B	24	ARG
2	B	86	GLU
2	B	101	LEU
2	B	113	ARG
2	B	121	SER
2	B	127	SER
2	B	147	LYS
2	B	202	THR
2	B	204	LYS
2	B	213	SER
3	Q	7	ASP
1	C	7	SER
1	C	13	GLN
1	C	75	THR
1	C	101	ARG
1	C	118	THR
1	C	123	SER
1	C	169	LEU
1	C	190	SER
1	C	195	SER
1	C	214	THR
1	C	221	VAL
2	D	1	ASP
2	D	5	THR
2	D	7	THR
2	D	10	SER
2	D	24	ARG
2	D	86	GLU

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Mol	Chain	Res	Type
2	D	101	LEU
2	D	113	ARG
2	D	121	SER
2	D	127	SER
2	D	147	LYS
2	D	202	THR
2	D	204	LYS
2	D	213	SER
3	R	7	ASP
1	E	7	SER
1	E	13	GLN
1	E	75	THR
1	E	87	VAL
1	E	101	ARG
1	E	118	THR
1	E	123	SER
1	E	168	SER
1	E	169	LEU
1	E	190	SER
1	E	195	SER
1	E	214	THR
1	E	221	VAL
2	F	1	ASP
2	F	5	THR
2	F	7	THR
2	F	10	SER
2	F	24	ARG
2	F	86	GLU
2	F	101	LEU
2	F	113	ARG
2	F	121	SER
2	F	127	SER
2	F	147	LYS
2	F	168	TRP
2	F	202	THR
2	F	204	LYS
2	F	213	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	16	GLN

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Mol	Chain	Res	Type
1	H	32	ASN
1	H	79	GLN
1	H	181	GLN
2	L	6	GLN
2	L	47	GLN
2	L	58	ASN
2	L	98	HIS
2	L	161	GLN
2	L	162	ASN
2	L	195	ASN
1	A	13	GLN
1	A	16	GLN
1	A	32	ASN
1	A	79	GLN
1	A	181	GLN
2	B	6	GLN
2	B	47	GLN
2	B	58	ASN
2	B	98	HIS
2	B	161	GLN
2	B	162	ASN
2	B	195	ASN
1	C	16	GLN
1	C	32	ASN
1	C	79	GLN
1	C	181	GLN
2	D	6	GLN
2	D	47	GLN
2	D	58	ASN
2	D	161	GLN
2	D	162	ASN
2	D	195	ASN
1	E	13	GLN
1	E	16	GLN
1	E	32	ASN
1	E	79	GLN
1	E	181	GLN
2	F	6	GLN
2	F	47	GLN
2	F	58	ASN
2	F	161	GLN
2	F	162	ASN

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Mol	Chain	Res	Type
2	F	195	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 ⓘ

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	216/226 (95%)	0.15	2 (0%) 81 76	15, 20, 33, 51	0
1	C	216/226 (95%)	0.31	7 (3%) 50 43	20, 29, 55, 65	0
1	E	216/226 (95%)	0.27	5 (2%) 61 53	14, 25, 55, 78	0
1	H	222/226 (98%)	0.15	6 (2%) 56 48	14, 23, 39, 63	0
2	B	218/219 (99%)	0.26	2 (0%) 81 76	15, 23, 40, 57	0
2	D	218/219 (99%)	0.29	3 (1%) 73 67	20, 30, 47, 92	0
2	F	218/219 (99%)	0.22	4 (1%) 67 59	15, 23, 52, 105	0
2	L	218/219 (99%)	0.24	7 (3%) 50 43	15, 24, 50, 91	0
3	P	6/7 (85%)	0.37	0 100 100	22, 30, 39, 43	0
3	Q	6/7 (85%)	0.48	1 (16%) 4 4	27, 31, 40, 54	0
3	R	6/7 (85%)	0.14	0 100 100	36, 36, 40, 45	0
3	S	5/7 (71%)	0.16	0 100 100	32, 32, 37, 38	0
All	All	1765/1808 (97%)	0.24	37 (2%) 63 55	14, 25, 49, 105	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	137	GLY	3.2
2	F	217	ASN	3.1
2	D	217	ASN	2.9
2	F	218	GLU	2.9
1	E	221	VAL	2.9
1	E	222	PRO	2.8
2	L	128	GLU	2.8
3	Q	2	ALA	2.7
1	C	128	PRO	2.5
1	H	140	ALA	2.5
2	B	195	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
2	D	218	GLU	2.5
1	C	74	ASP	2.5
2	F	24	ARG	2.4
1	A	144	SER	2.3
1	C	77	ASN	2.3
2	B	217	ASN	2.3
2	F	192	GLU	2.3
2	L	134	GLY	2.3
1	H	145	MET	2.2
1	E	44	GLY	2.2
1	A	145	MET	2.2
1	E	128	PRO	2.2
1	C	137	GLY	2.2
1	H	57	GLU	2.2
2	L	217	ASN	2.2
2	L	86	GLU	2.2
2	L	192	GLU	2.2
1	H	13	GLN	2.1
1	C	145	MET	2.1
1	H	137	GLY	2.1
2	D	24	ARG	2.1
2	L	75	ASP	2.1
2	L	34	GLY	2.0
1	H	142	THR	2.0
1	C	31	THR	2.0
1	C	105	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.