



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:05 PM UTC

PDB ID : 1EOH / pdb_00001eoh
Title : GLUTATHIONE TRANSFERASE P1-1
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Deposited on : 2000-03-22
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

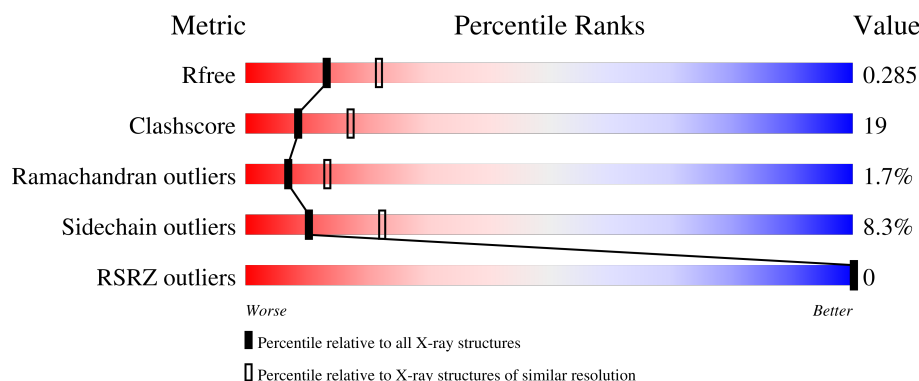
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	209	 67% 27% 5%
1	B	209	 64% 32% 4% 1%
1	C	209	 58% 35% 6%
1	D	209	 60% 33% 6%
1	E	209	 65% 31% 4% 1%

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Mol	Chain	Length	Quality of chain
1	F	209	 66% 27% 7%
1	G	209	 63% 28% 9%
1	H	209	 58% 35% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13580 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 GLUTATHIONE S-TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	0	0
			1635	1051	272	306	6			
1	B	209	Total	C	N	O	S	0	0	0
			1635	1051	272	306	6			
1	C	209	Total	C	N	O	S	0	0	0
			1635	1051	272	306	6			
1	D	209	Total	C	N	O	S	0	0	0
			1635	1051	272	306	6			
1	E	209	Total	C	N	O	S	0	0	0
			1635	1051	272	306	6			
1	F	209	Total	C	N	O	S	0	0	0
			1635	1051	272	306	6			
1	G	209	Total	C	N	O	S	0	0	0
			1635	1051	272	306	6			
1	H	209	Total	C	N	O	S	0	0	0
			1635	1051	272	306	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	152	ALA	ASP	engineered mutation	UNP P09211
B	152	ALA	ASP	engineered mutation	UNP P09211
C	152	ALA	ASP	engineered mutation	UNP P09211
D	152	ALA	ASP	engineered mutation	UNP P09211
E	152	ALA	ASP	engineered mutation	UNP P09211
F	152	ALA	ASP	engineered mutation	UNP P09211
G	152	ALA	ASP	engineered mutation	UNP P09211
H	152	ALA	ASP	engineered mutation	UNP P09211

- Molecule 2 is water.

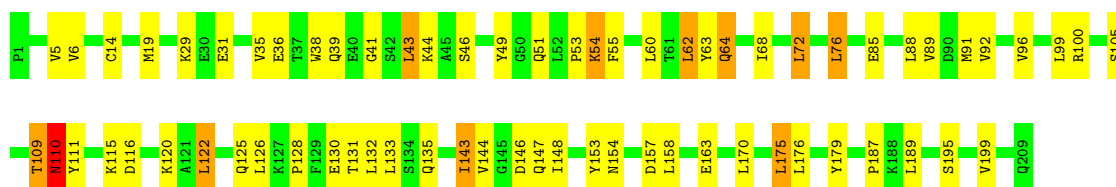
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	72	Total 72	O 72	0	0
2	B	96	Total 96	O 96	0	0
2	C	51	Total 51	O 51	0	0
2	D	54	Total 54	O 54	0	0
2	E	70	Total 70	O 70	0	0
2	F	69	Total 69	O 69	0	0
2	G	32	Total 32	O 32	0	0
2	H	56	Total 56	O 56	0	0

3 Residue-property plots

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

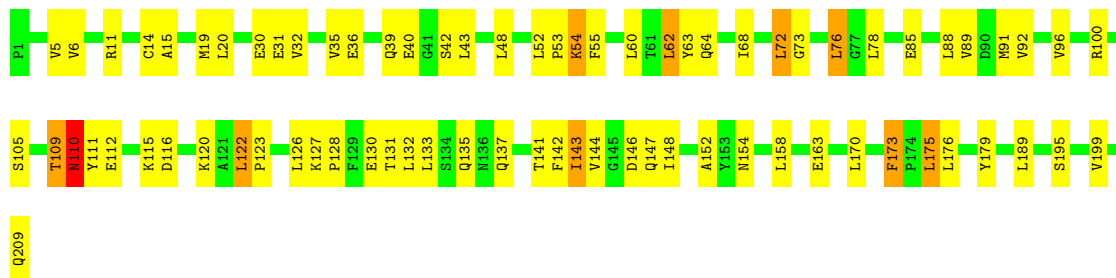
• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain A: 



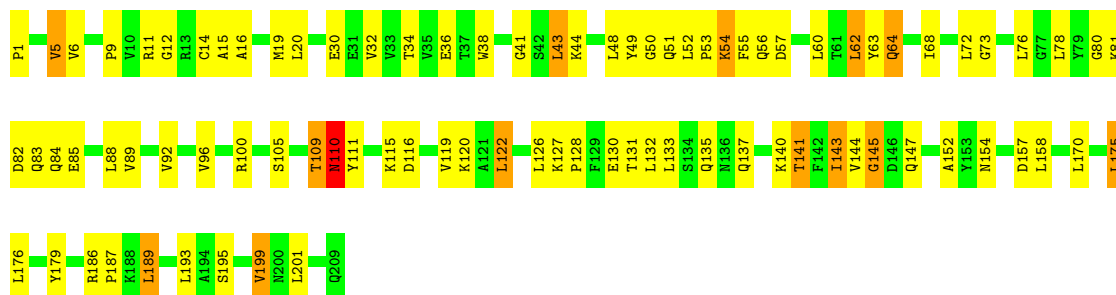
• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain B: 



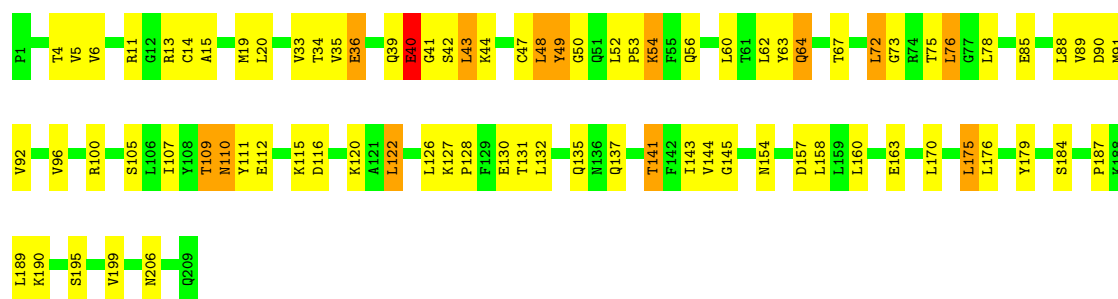
• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain C: 



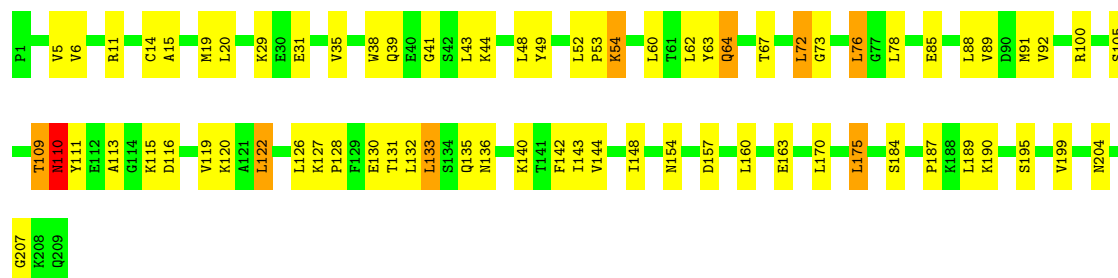
• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain D: 



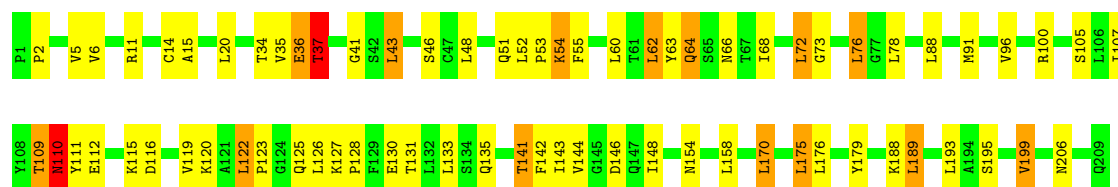
• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain E: 65% 31% .



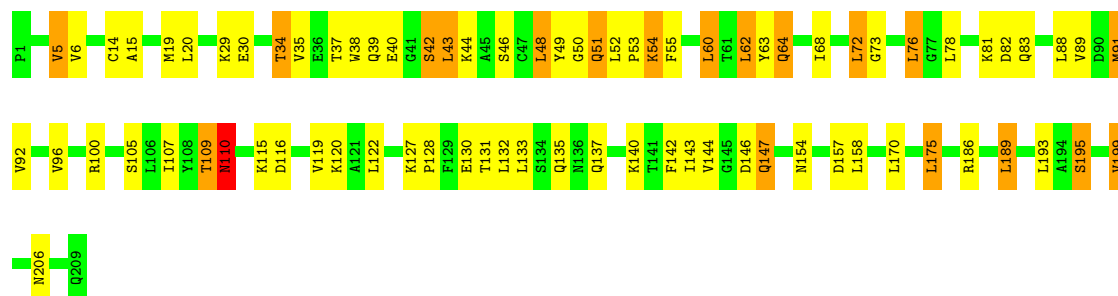
• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain F: 66% 27% 7% .



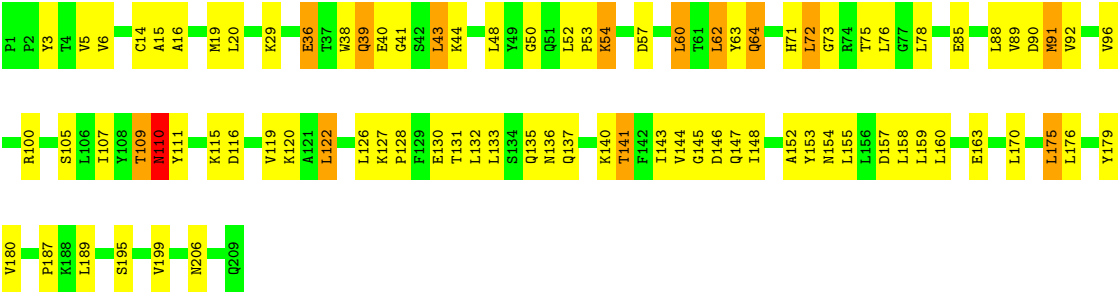
• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain G: 63% 28% 9%



• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain H: 58% 35% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.97Å 84.02Å 236.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.50) 95.7 (20.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.52 (at 2.50Å)	Xtriage
Refinement program	CNS 0.4	Depositor
R, R_{free}	0.231 , 0.286 0.233 , 0.285	Depositor DCC
R_{free} test set	2923 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	0.477	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 17.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.146 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13580	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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 27.39 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.2164e-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

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	0.53	0/1670	0.96	3/2265 (0.1%)
1	B	0.58	0/1670	0.94	6/2265 (0.3%)
1	C	0.54	0/1670	0.96	3/2265 (0.1%)
1	D	0.52	0/1670	0.92	3/2265 (0.1%)
1	E	0.53	0/1670	0.95	3/2265 (0.1%)
1	F	0.54	0/1670	0.94	5/2265 (0.2%)
1	G	0.54	0/1670	0.95	7/2265 (0.3%)
1	H	0.59	0/1670	0.98	2/2265 (0.1%)
All	All	0.55	0/13360	0.95	32/18120 (0.2%)

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	41	GLY	N-CA-C	8.90	123.10	113.58
1	E	41	GLY	N-CA-C	7.88	124.45	115.08
1	E	54	LYS	N-CA-C	-7.22	97.96	109.59
1	F	41	GLY	N-CA-C	6.88	124.48	115.47
1	F	37	THR	N-CA-C	-6.50	104.11	111.07
1	B	42	SER	N-CA-C	6.29	118.66	111.11
1	H	43	LEU	N-CA-C	-6.26	103.75	112.45
1	H	54	LYS	N-CA-C	-6.22	99.57	109.59
1	B	54	LYS	N-CA-C	-6.04	99.55	109.40
1	A	54	LYS	N-CA-C	-6.00	99.93	109.59
1	D	112	GLU	N-CA-C	5.93	117.42	111.07
1	C	199	VAL	N-CA-C	5.83	116.56	110.62
1	G	54	LYS	N-CA-C	-5.57	100.63	109.59
1	A	163	GLU	N-CA-C	-5.54	104.87	111.69
1	B	163	GLU	N-CA-C	-5.51	104.91	111.69
1	A	41	GLY	N-CA-C	5.45	122.61	115.47
1	D	90	ASP	N-CA-C	-5.40	105.29	111.07
1	G	195	SER	CA-C-N	5.38	125.02	119.05
1	G	195	SER	C-N-CA	5.38	125.02	119.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	54	LYS	N-CA-C	-5.34	100.69	109.40
1	E	113	ALA	N-CA-C	5.32	117.50	111.11
1	D	54	LYS	N-CA-C	-5.29	100.56	109.07
1	B	173	PHE	CA-C-N	5.26	125.07	119.28
1	B	173	PHE	C-N-CA	5.26	125.07	119.28
1	G	89	VAL	N-CA-C	-5.14	105.31	110.30
1	B	32	VAL	N-CA-C	5.12	115.66	108.89
1	F	54	LYS	N-CA-C	-5.12	100.56	108.90
1	G	51	GLN	N-CA-C	5.10	116.74	109.14
1	G	199	VAL	N-CA-C	5.08	115.80	110.62
1	F	148	ILE	N-CA-C	5.08	116.11	109.21
1	F	199	VAL	N-CA-C	5.05	115.77	110.62
1	G	48	LEU	N-CA-C	5.05	117.17	111.11

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	1635	0	1644	52	0
1	B	1635	0	1644	56	0
1	C	1635	0	1644	77	0
1	D	1635	0	1644	75	0
1	E	1635	0	1644	57	0
1	F	1635	0	1644	60	0
1	G	1635	0	1644	73	0
1	H	1635	0	1644	73	1
2	A	72	0	0	3	0
2	B	96	0	0	11	0
2	C	51	0	0	4	0
2	D	54	0	0	6	0
2	E	70	0	0	9	0
2	F	69	0	0	8	1
2	G	32	0	0	1	0
2	H	56	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	13580	0	13152	494	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 19.

All (494) 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:133:LEU:HD13	1:B:143:ILE:HD12	1.35	1.04
1:G:48:LEU:CD2	1:H:91:MET:HG2	1.87	1.03
1:C:92:VAL:HG21	1:C:143:ILE:HD11	1.46	0.98
1:F:34:THR:HG23	1:F:36:GLU:OE1	1.62	0.97
1:C:132:LEU:HG	1:D:48:LEU:HD11	1.49	0.94
1:F:130:GLU:OE2	1:F:175:LEU:HB2	1.69	0.93
1:G:105:SER:O	1:G:109:THR:HB	1.73	0.89
1:G:132:LEU:HG	1:H:48:LEU:HD11	1.56	0.87
1:H:130:GLU:OE2	1:H:175:LEU:HB2	1.73	0.87
1:C:105:SER:O	1:C:109:THR:HB	1.76	0.86
1:D:130:GLU:OE2	1:D:175:LEU:HB2	1.78	0.82
1:E:105:SER:O	1:E:109:THR:HB	1.79	0.81
1:E:6:VAL:HB	1:E:54:LYS:HB3	1.63	0.81
1:G:48:LEU:HD23	1:H:91:MET:HG2	1.63	0.80
1:C:48:LEU:CD2	1:D:91:MET:HG2	2.11	0.80
1:B:130:GLU:OE2	1:B:175:LEU:HB2	1.81	0.80
1:C:130:GLU:OE2	1:C:175:LEU:HB2	1.81	0.80
1:F:105:SER:O	1:F:109:THR:HB	1.81	0.79
1:F:170:LEU:HB2	2:F:238:HOH:O	1.83	0.78
1:A:105:SER:O	1:A:109:THR:HB	1.84	0.77
1:A:6:VAL:HB	1:A:54:LYS:HB3	1.67	0.76
1:A:130:GLU:OE2	1:A:175:LEU:HB2	1.85	0.75
1:E:130:GLU:OE2	1:E:175:LEU:HB2	1.87	0.75
1:B:85:GLU:O	1:B:89:VAL:HG23	1.87	0.75
1:B:105:SER:O	1:B:109:THR:HB	1.87	0.74
1:G:100:ARG:HH21	1:G:154:ASN:ND2	1.86	0.74
1:D:120:LYS:HG2	2:D:243:HOH:O	1.87	0.73
1:G:92:VAL:HG21	1:G:143:ILE:HG12	1.70	0.73
1:C:85:GLU:HB3	1:C:144:VAL:HG11	1.70	0.73
1:C:143:ILE:HG13	1:C:144:VAL:HG23	1.71	0.73
1:H:105:SER:O	1:H:109:THR:HB	1.88	0.73
1:E:35:VAL:HG23	2:E:222:HOH:O	1.89	0.72
1:F:143:ILE:HG23	1:F:144:VAL:HG23	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:6:VAL:HB	1:H:54:LYS:HB3	1.70	0.71
1:B:100:ARG:HH21	1:B:154:ASN:ND2	1.87	0.71
1:G:130:GLU:OE2	1:G:175:LEU:HB2	1.90	0.71
1:H:144:VAL:HA	2:H:229:HOH:O	1.90	0.71
1:G:88:LEU:HD23	1:G:144:VAL:HG13	1.73	0.71
1:B:143:ILE:HG23	1:B:144:VAL:HG23	1.73	0.70
1:B:152:ALA:HB2	2:B:269:HOH:O	1.91	0.70
1:C:137:GLN:O	1:C:140:LYS:HE2	1.90	0.70
2:E:261:HOH:O	1:F:60:LEU:HD21	1.90	0.70
1:D:39:GLN:C	1:D:41:GLY:H	1.98	0.70
1:C:195:SER:O	1:C:199:VAL:HG23	1.92	0.69
1:H:96:VAL:HG13	1:H:158:LEU:HD22	1.74	0.69
1:D:100:ARG:HH21	1:D:154:ASN:ND2	1.91	0.68
1:G:38:TRP:HH2	1:G:51:GLN:HA	1.59	0.68
1:D:105:SER:O	1:D:109:THR:HB	1.94	0.68
1:B:35:VAL:O	1:B:39:GLN:HG3	1.93	0.67
1:G:6:VAL:HB	1:G:54:LYS:HB3	1.76	0.67
1:E:100:ARG:HH21	1:E:154:ASN:ND2	1.92	0.67
1:C:54:LYS:HG3	1:C:63:TYR:CE1	2.28	0.67
1:F:36:GLU:OE1	1:F:37:THR:N	2.25	0.67
1:C:43:LEU:HD22	1:C:43:LEU:O	1.94	0.67
1:G:49:TYR:HD2	1:G:64:GLN:HE21	1.40	0.67
1:C:6:VAL:HB	1:C:54:LYS:HB3	1.76	0.67
1:C:85:GLU:O	1:C:89:VAL:HG23	1.95	0.67
1:G:49:TYR:CD1	1:H:128:PRO:HB3	2.29	0.67
1:B:6:VAL:HB	1:B:54:LYS:HB3	1.77	0.66
1:D:39:GLN:O	1:D:41:GLY:N	2.29	0.66
1:G:48:LEU:HD21	1:H:91:MET:HG2	1.77	0.66
1:G:52:LEU:HB3	1:G:53:PRO:HA	1.76	0.66
1:H:29:LYS:HD2	2:H:228:HOH:O	1.95	0.66
1:D:39:GLN:C	1:D:41:GLY:N	2.45	0.66
1:D:63:TYR:O	1:D:64:GLN:HB2	1.95	0.66
1:F:36:GLU:N	1:F:36:GLU:CD	2.52	0.66
1:D:6:VAL:HB	1:D:54:LYS:HB3	1.78	0.66
1:E:73:GLY:HA2	1:E:78:LEU:HB2	1.78	0.66
1:E:88:LEU:HD12	1:E:91:MET:HE3	1.78	0.66
1:F:116:ASP:O	1:F:120:LYS:HG3	1.96	0.66
1:C:73:GLY:HA2	1:C:78:LEU:HB2	1.78	0.66
1:A:100:ARG:HH21	1:A:154:ASN:ND2	1.94	0.65
1:E:63:TYR:O	1:E:64:GLN:HB2	1.95	0.64
1:B:143:ILE:HG22	2:B:269:HOH:O	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:96:VAL:HG13	1:G:158:LEU:HD22	1.79	0.64
1:C:81:LYS:NZ	1:C:82:ASP:OD2	2.30	0.64
1:E:48:LEU:HD23	1:E:49:TYR:CE2	2.33	0.64
1:G:131:THR:O	1:G:135:GLN:HG3	1.98	0.64
1:F:63:TYR:O	1:F:64:GLN:HB2	1.96	0.64
1:C:48:LEU:HD23	1:D:91:MET:HG2	1.80	0.64
1:A:35:VAL:O	1:A:39:GLN:HG3	1.98	0.63
1:H:137:GLN:HB3	1:H:140:LYS:HE2	1.79	0.63
1:G:81:LYS:NZ	1:G:82:ASP:OD2	2.30	0.63
1:F:51:GLN:OE1	2:F:222:HOH:O	2.15	0.63
1:A:63:TYR:O	1:A:64:GLN:HB2	1.97	0.63
1:B:73:GLY:HA2	1:B:78:LEU:HB2	1.81	0.63
1:B:110:ASN:HA	2:B:248:HOH:O	1.99	0.63
1:A:133:LEU:HD13	1:A:143:ILE:HG13	1.80	0.63
1:H:60:LEU:HD22	1:H:62:LEU:CD1	2.30	0.62
1:G:73:GLY:HA2	1:G:78:LEU:HB2	1.82	0.62
1:D:43:LEU:HD22	1:D:47:CYS:SG	2.40	0.62
1:F:43:LEU:HD23	1:F:46:SER:OG	1.99	0.62
1:H:137:GLN:O	1:H:140:LYS:HE2	1.99	0.62
1:G:49:TYR:HD2	1:G:64:GLN:NE2	1.97	0.62
1:C:85:GLU:OE2	1:C:145:GLY:HA3	2.00	0.61
1:G:29:LYS:HE2	1:G:30:GLU:O	2.00	0.61
1:C:1:PRO:N	1:C:57:ASP:OD1	2.32	0.60
1:E:88:LEU:HG	1:E:144:VAL:HG22	1.83	0.60
1:G:49:TYR:HD1	1:H:128:PRO:HB3	1.67	0.60
1:D:41:GLY:HA2	1:D:44:LYS:HB3	1.84	0.60
1:D:34:THR:OG1	1:D:36:GLU:HG2	2.01	0.60
1:C:127:LYS:HB3	1:C:128:PRO:HD3	1.84	0.60
1:A:88:LEU:HD23	1:A:144:VAL:HG22	1.83	0.60
1:C:19:MET:CE	1:C:157:ASP:HB2	2.32	0.60
1:F:131:THR:O	1:F:135:GLN:HG3	2.01	0.60
1:G:142:PHE:N	1:G:142:PHE:CD2	2.69	0.60
1:D:85:GLU:O	1:D:89:VAL:HG23	2.02	0.59
1:C:5:VAL:HG22	1:C:30:GLU:OE1	2.01	0.59
1:D:116:ASP:O	1:D:120:LYS:HG3	2.02	0.59
1:C:132:LEU:CG	1:D:48:LEU:HD11	2.28	0.59
1:G:63:TYR:O	1:G:64:GLN:HB2	2.03	0.59
1:B:63:TYR:O	1:B:64:GLN:HB2	2.00	0.59
1:H:20:LEU:C	1:H:20:LEU:HD23	2.28	0.59
1:F:195:SER:O	1:F:199:VAL:HG23	2.01	0.59
1:H:92:VAL:HG22	1:H:132:LEU:HD13	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:110:ASN:N	1:H:110:ASN:HD22	1.99	0.59
1:D:109:THR:HG23	2:D:252:HOH:O	2.02	0.58
1:B:147:GLN:NE2	2:B:218:HOH:O	2.35	0.58
1:F:36:GLU:N	1:F:36:GLU:OE2	2.36	0.58
1:B:133:LEU:CD1	1:B:143:ILE:HD12	2.23	0.58
1:F:96:VAL:HG13	1:F:158:LEU:HD22	1.86	0.58
1:C:83:GLN:OE1	1:D:75:THR:HG22	2.02	0.58
1:F:73:GLY:HA2	1:F:78:LEU:HB2	1.86	0.58
1:A:14:CYS:SG	1:A:53:PRO:HB3	2.44	0.58
1:E:88:LEU:CD2	1:E:144:VAL:HG22	2.34	0.58
1:G:100:ARG:HH21	1:G:154:ASN:HD21	1.48	0.58
1:C:48:LEU:HD11	1:D:132:LEU:HG	1.85	0.58
2:E:221:HOH:O	1:F:64:GLN:HB3	2.04	0.58
1:H:127:LYS:HB3	1:H:128:PRO:HD3	1.84	0.58
1:F:60:LEU:HD22	1:F:62:LEU:CD1	2.34	0.57
1:G:49:TYR:CD2	1:G:64:GLN:NE2	2.72	0.57
1:H:85:GLU:O	1:H:89:VAL:HG23	2.03	0.57
1:E:92:VAL:HG22	1:E:132:LEU:HD13	1.86	0.57
1:C:63:TYR:O	1:C:64:GLN:HB2	2.03	0.57
1:E:130:GLU:CD	1:E:175:LEU:HB2	2.30	0.57
1:G:195:SER:O	1:G:199:VAL:HG23	2.04	0.57
1:H:63:TYR:O	1:H:64:GLN:HB2	2.04	0.57
1:F:14:CYS:SG	1:F:53:PRO:HB3	2.44	0.57
1:B:88:LEU:HA	1:B:91:MET:HE2	1.86	0.56
1:G:142:PHE:HD2	1:G:142:PHE:H	1.52	0.56
1:D:96:VAL:HG13	1:D:158:LEU:HD22	1.85	0.56
1:B:131:THR:O	1:B:135:GLN:HG3	2.06	0.56
1:G:116:ASP:O	1:G:120:LYS:HG3	2.05	0.56
1:A:122:LEU:HD22	1:A:126:LEU:HG	1.86	0.56
1:B:14:CYS:SG	1:B:53:PRO:HB3	2.46	0.56
1:B:195:SER:O	1:B:199:VAL:HG23	2.06	0.56
1:D:143:ILE:HG23	1:D:144:VAL:HG23	1.87	0.56
1:E:116:ASP:O	1:E:120:LYS:HG3	2.05	0.56
1:G:34:THR:HG23	1:G:37:THR:OG1	2.06	0.56
1:C:176:LEU:O	1:C:179:TYR:HB3	2.06	0.56
1:E:195:SER:O	1:E:199:VAL:HG23	2.05	0.56
1:H:116:ASP:O	1:H:120:LYS:HG3	2.06	0.56
1:B:123:PRO:HG3	2:B:229:HOH:O	2.04	0.56
1:F:36:GLU:CD	1:F:37:THR:H	2.13	0.56
1:D:195:SER:O	1:D:199:VAL:HG23	2.06	0.55
1:G:137:GLN:O	1:G:140:LYS:HE3	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:LEU:HD12	1:B:91:MET:HE3	1.88	0.55
1:D:11:ARG:NH1	2:D:236:HOH:O	2.39	0.55
1:F:34:THR:HG23	1:F:36:GLU:CD	2.29	0.55
1:H:38:TRP:CZ2	1:H:44:LYS:HG3	2.41	0.55
1:A:64:GLN:HB3	2:A:246:HOH:O	2.06	0.55
1:F:60:LEU:HD22	1:F:62:LEU:HD13	1.88	0.55
2:E:221:HOH:O	1:F:66:ASN:HB2	2.06	0.55
1:G:54:LYS:HG3	1:G:63:TYR:CE1	2.42	0.55
1:H:19:MET:CE	1:H:157:ASP:HB2	2.37	0.55
1:D:67:THR:HG23	2:D:213:HOH:O	2.05	0.55
1:B:52:LEU:HB3	1:B:53:PRO:HA	1.87	0.55
1:D:92:VAL:HG21	1:D:143:ILE:HG13	1.89	0.55
1:G:115:LYS:O	1:G:119:VAL:HG23	2.07	0.55
1:F:6:VAL:HB	1:F:54:LYS:HB3	1.88	0.55
1:F:52:LEU:HB3	1:F:53:PRO:HA	1.89	0.55
1:G:88:LEU:HG	1:G:144:VAL:HG22	1.89	0.55
1:E:122:LEU:HD22	1:E:126:LEU:HG	1.87	0.55
1:H:36:GLU:O	1:H:39:GLN:HB3	2.05	0.55
1:H:122:LEU:HD22	1:H:126:LEU:HG	1.89	0.55
1:B:15:ALA:O	1:B:19:MET:HG3	2.06	0.55
1:D:88:LEU:HG	1:D:144:VAL:HG22	1.89	0.55
1:H:60:LEU:HD22	1:H:62:LEU:HD13	1.88	0.55
1:C:52:LEU:HB3	1:C:53:PRO:HA	1.89	0.54
1:E:14:CYS:SG	1:E:53:PRO:HB3	2.47	0.54
1:G:19:MET:CE	1:G:157:ASP:HB2	2.37	0.54
1:A:92:VAL:HG22	1:A:132:LEU:HD13	1.90	0.54
1:D:130:GLU:CD	1:D:175:LEU:HB2	2.32	0.54
1:E:11:ARG:HB3	1:E:15:ALA:HB2	1.90	0.54
1:E:54:LYS:HG3	1:E:63:TYR:CE1	2.43	0.54
1:A:116:ASP:O	1:A:120:LYS:HG3	2.07	0.54
1:D:100:ARG:HH21	1:D:154:ASN:HD21	1.56	0.54
1:D:127:LYS:HB3	1:D:128:PRO:HD3	1.89	0.54
1:F:112:GLU:HG2	2:F:248:HOH:O	2.07	0.54
1:G:60:LEU:HD22	1:G:62:LEU:CD1	2.37	0.54
1:F:100:ARG:HH21	1:F:154:ASN:ND2	2.06	0.54
1:B:78:LEU:HD22	1:B:148:ILE:HG23	1.89	0.54
1:G:92:VAL:HG22	1:G:132:LEU:HD13	1.89	0.54
1:E:88:LEU:HD23	1:E:144:VAL:HG22	1.90	0.54
1:G:127:LYS:HB3	1:G:128:PRO:HD3	1.90	0.54
1:B:112:GLU:HB2	2:B:270:HOH:O	2.09	0.53
1:C:36:GLU:HB3	2:C:236:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:44:LYS:O	1:H:50:GLY:HA2	2.08	0.53
1:C:84:GLN:HG3	1:D:60:LEU:HD21	1.91	0.53
1:C:92:VAL:HG22	1:C:132:LEU:HD13	1.91	0.53
1:G:48:LEU:CD2	1:H:91:MET:CG	2.75	0.53
1:A:111:TYR:O	1:A:115:LYS:HB2	2.09	0.53
1:H:15:ALA:O	1:H:19:MET:HG3	2.07	0.53
1:B:137:GLN:OE1	1:B:141:THR:HG21	2.08	0.53
1:A:143:ILE:CG2	1:A:144:VAL:HG23	2.39	0.53
1:E:88:LEU:CG	1:E:144:VAL:HG22	2.39	0.53
1:H:148:ILE:HG21	1:H:153:TYR:OH	2.08	0.53
1:E:88:LEU:HD23	1:E:144:VAL:HG13	1.90	0.53
1:H:52:LEU:HB3	1:H:53:PRO:HA	1.91	0.53
1:A:29:LYS:HZ3	1:A:31:GLU:HG3	1.74	0.53
1:C:96:VAL:HG13	1:C:158:LEU:HD22	1.91	0.53
1:B:116:ASP:O	1:B:120:LYS:HG3	2.08	0.52
1:D:52:LEU:HB3	1:D:53:PRO:HA	1.90	0.52
1:D:88:LEU:HA	1:D:91:MET:HE2	1.89	0.52
1:E:143:ILE:HG23	1:E:144:VAL:HG23	1.91	0.52
1:A:72:LEU:O	1:A:76:LEU:HB2	2.09	0.52
1:C:14:CYS:SG	1:C:53:PRO:HB3	2.50	0.52
1:A:19:MET:CE	1:A:157:ASP:HB2	2.39	0.52
1:A:88:LEU:CD2	1:A:144:VAL:HG22	2.40	0.52
1:C:85:GLU:CB	1:C:144:VAL:HG11	2.39	0.52
1:E:38:TRP:CH2	1:E:44:LYS:HA	2.45	0.52
1:C:43:LEU:HD22	1:C:43:LEU:C	2.35	0.51
1:C:43:LEU:HD13	1:C:43:LEU:C	2.35	0.51
1:G:44:LYS:O	1:G:50:GLY:HA2	2.09	0.51
1:G:92:VAL:HG21	1:G:143:ILE:CG1	2.37	0.51
1:C:44:LYS:O	1:C:50:GLY:HA2	2.10	0.51
1:G:15:ALA:O	1:G:19:MET:HG3	2.10	0.51
1:B:100:ARG:HH21	1:B:154:ASN:HD21	1.54	0.51
1:E:35:VAL:O	1:E:39:GLN:HB2	2.10	0.51
1:E:78:LEU:HD22	1:E:148:ILE:HG23	1.93	0.51
1:C:122:LEU:HD22	1:C:126:LEU:HG	1.92	0.51
1:C:100:ARG:HH21	1:C:154:ASN:ND2	2.08	0.51
1:G:130:GLU:CD	1:G:175:LEU:HB2	2.36	0.51
1:E:20:LEU:C	1:E:20:LEU:HD23	2.36	0.51
1:A:60:LEU:HD22	1:A:62:LEU:HD13	1.93	0.51
1:H:131:THR:O	1:H:135:GLN:HG3	2.11	0.50
1:H:143:ILE:HG22	1:H:152:ALA:HB2	1.93	0.50
1:E:29:LYS:HD2	2:E:220:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:88:LEU:HA	1:H:91:MET:HE2	1.92	0.50
1:C:43:LEU:HD13	1:C:44:LYS:N	2.26	0.50
1:D:92:VAL:HG11	1:D:143:ILE:HG12	1.93	0.50
1:F:123:PRO:HG3	2:F:264:HOH:O	2.10	0.50
1:G:132:LEU:HD11	1:H:48:LEU:HD21	1.94	0.50
1:B:20:LEU:HD23	1:B:20:LEU:C	2.36	0.50
1:C:189:LEU:HD22	1:C:193:LEU:HG	1.92	0.50
1:G:38:TRP:CH2	1:G:51:GLN:HA	2.43	0.50
1:D:131:THR:O	1:D:135:GLN:HG3	2.12	0.50
1:D:88:LEU:CD2	1:D:144:VAL:HG22	2.41	0.50
1:G:19:MET:HE2	1:G:157:ASP:HB2	1.93	0.50
1:F:127:LYS:HB3	1:F:128:PRO:HD3	1.94	0.50
1:A:60:LEU:HD22	1:A:62:LEU:CD1	2.42	0.49
1:F:20:LEU:C	1:F:20:LEU:HD23	2.37	0.49
1:H:130:GLU:CD	1:H:175:LEU:HB2	2.36	0.49
1:D:15:ALA:O	1:D:19:MET:HG3	2.12	0.49
1:F:122:LEU:HD22	1:F:126:LEU:HG	1.93	0.49
1:H:71:HIS:HB2	2:H:241:HOH:O	2.12	0.49
1:E:67:THR:HG23	2:F:273:HOH:O	2.12	0.49
1:B:141:THR:HG23	1:B:142:PHE:N	2.28	0.49
1:E:131:THR:O	1:E:135:GLN:HG3	2.12	0.49
1:F:189:LEU:HD22	1:F:193:LEU:HG	1.94	0.49
1:A:125:GLN:NE2	2:A:219:HOH:O	2.45	0.49
1:C:88:LEU:O	1:C:92:VAL:HG23	2.13	0.49
1:D:92:VAL:HG21	1:D:143:ILE:CG1	2.43	0.48
1:C:80:GLY:C	1:C:147:GLN:OE1	2.57	0.48
1:D:85:GLU:OE2	1:D:145:GLY:HA3	2.12	0.48
1:E:19:MET:HE2	2:E:219:HOH:O	2.12	0.48
1:H:41:GLY:HA2	1:H:44:LYS:CB	2.43	0.48
1:H:133:LEU:HD21	1:H:179:TYR:HB2	1.96	0.48
1:G:63:TYR:O	1:G:64:GLN:CB	2.61	0.48
1:B:133:LEU:HD21	1:B:179:TYR:HB2	1.96	0.48
1:E:63:TYR:O	1:E:64:GLN:CB	2.61	0.48
1:B:40:GLU:OE1	1:B:40:GLU:C	2.57	0.48
1:C:16:ALA:HB3	2:C:218:HOH:O	2.12	0.48
1:F:54:LYS:HG3	1:F:63:TYR:CE1	2.49	0.48
1:B:130:GLU:OE1	1:B:173:PHE:HB3	2.15	0.47
1:B:142:PHE:CD2	1:B:146:ASP:C	2.91	0.47
1:D:20:LEU:C	1:D:20:LEU:HD23	2.39	0.47
1:H:115:LYS:O	1:H:119:VAL:HG23	2.14	0.47
1:C:143:ILE:HG21	1:C:152:ALA:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:176:LEU:O	1:D:179:TYR:HB3	2.14	0.47
1:D:39:GLN:O	1:D:40:GLU:C	2.57	0.47
1:D:40:GLU:C	1:D:42:SER:H	2.21	0.47
1:E:115:LYS:O	1:E:119:VAL:HG23	2.14	0.47
1:F:63:TYR:O	1:F:64:GLN:CB	2.62	0.47
1:F:55:PHE:HB2	1:F:68:ILE:HD13	1.95	0.47
1:B:60:LEU:HD22	1:B:62:LEU:CD1	2.45	0.47
1:G:14:CYS:SG	1:G:53:PRO:HB3	2.55	0.47
1:A:38:TRP:CH2	1:A:44:LYS:HA	2.49	0.47
1:C:19:MET:HE2	1:C:157:ASP:HB2	1.97	0.47
1:C:111:TYR:O	1:C:115:LYS:HB2	2.15	0.47
1:D:110:ASN:O	1:D:110:ASN:CG	2.57	0.47
1:F:2:PRO:HA	2:F:246:HOH:O	2.14	0.47
1:H:155:LEU:O	1:H:159:LEU:HG	2.14	0.47
1:C:11:ARG:HB3	1:C:15:ALA:HB2	1.96	0.47
1:D:11:ARG:HB3	1:D:15:ALA:HB2	1.97	0.47
1:H:36:GLU:CD	1:H:36:GLU:H	2.23	0.47
1:H:38:TRP:O	1:H:41:GLY:N	2.47	0.47
1:A:88:LEU:HA	1:A:91:MET:HE2	1.97	0.47
1:A:100:ARG:HH21	1:A:154:ASN:HD21	1.63	0.47
1:F:110:ASN:N	1:F:110:ASN:HD22	2.13	0.47
1:H:111:TYR:O	1:H:115:LYS:HB2	2.15	0.47
1:C:132:LEU:HD11	1:D:48:LEU:HD21	1.97	0.47
1:C:48:LEU:HD11	1:D:132:LEU:CG	2.45	0.46
1:D:19:MET:CE	1:D:157:ASP:HB2	2.45	0.46
1:D:48:LEU:C	1:D:50:GLY:H	2.23	0.46
1:H:176:LEU:O	1:H:180:VAL:HG23	2.16	0.46
1:A:111:TYR:CZ	1:A:115:LYS:HD2	2.51	0.46
1:A:130:GLU:CD	1:A:175:LEU:HB2	2.40	0.46
1:G:88:LEU:O	1:G:92:VAL:HG23	2.15	0.46
1:B:96:VAL:HG13	1:B:158:LEU:HD22	1.98	0.46
1:D:107:ILE:HD13	1:D:111:TYR:HD1	1.80	0.46
1:E:160:LEU:O	1:E:163:GLU:HB2	2.15	0.46
1:C:92:VAL:HA	1:D:49:TYR:OH	2.15	0.46
1:D:184:SER:O	1:D:190:LYS:HB2	2.15	0.46
1:F:111:TYR:CZ	1:F:115:LYS:HD2	2.51	0.46
1:F:125:GLN:O	1:F:128:PRO:HD2	2.16	0.46
1:F:133:LEU:HD13	1:F:143:ILE:HD12	1.96	0.46
1:H:88:LEU:O	1:H:92:VAL:HG23	2.15	0.46
1:D:92:VAL:HG22	1:D:132:LEU:HD13	1.97	0.46
1:D:154:ASN:ND2	2:D:230:HOH:O	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:49:TYR:HD1	1:H:128:PRO:CB	2.26	0.46
1:E:140:LYS:HE2	2:E:262:HOH:O	2.16	0.46
1:G:83:GLN:CD	1:H:75:THR:HG22	2.41	0.46
1:C:131:THR:O	1:C:135:GLN:HG3	2.16	0.46
1:E:19:MET:CE	1:E:157:ASP:HB2	2.46	0.46
1:F:130:GLU:CD	1:F:175:LEU:HB2	2.38	0.46
1:H:85:GLU:OE2	1:H:145:GLY:CA	2.64	0.46
1:A:43:LEU:O	1:A:44:LYS:C	2.59	0.45
1:B:11:ARG:HB3	1:B:15:ALA:HB2	1.97	0.45
1:H:3:TYR:CE1	1:H:57:ASP:OD1	2.70	0.45
1:H:41:GLY:HA2	1:H:44:LYS:HB3	1.98	0.45
1:H:160:LEU:O	1:H:163:GLU:HB2	2.17	0.45
1:G:147:GLN:HE21	1:G:147:GLN:HB2	1.57	0.45
1:E:136:ASN:OD1	1:E:142:PHE:O	2.33	0.45
1:F:176:LEU:O	1:F:179:TYR:HB3	2.16	0.45
1:G:88:LEU:HD11	1:G:132:LEU:HD22	1.99	0.45
1:B:88:LEU:HD23	1:B:144:VAL:HG22	1.99	0.45
1:H:14:CYS:SG	1:H:53:PRO:HB3	2.55	0.45
1:F:72:LEU:O	1:F:76:LEU:HB2	2.16	0.45
1:B:11:ARG:NH2	2:B:272:HOH:O	2.38	0.45
1:E:85:GLU:O	1:E:89:VAL:HG23	2.17	0.45
1:H:195:SER:O	1:H:199:VAL:HG23	2.17	0.45
1:B:130:GLU:CD	1:B:175:LEU:HB2	2.42	0.45
1:D:34:THR:CB	1:D:36:GLU:HG2	2.46	0.45
1:A:55:PHE:HB2	1:A:68:ILE:CD1	2.47	0.45
1:D:73:GLY:HA2	1:D:78:LEU:HB2	1.99	0.45
1:E:48:LEU:CD2	1:F:91:MET:HG2	2.47	0.45
1:H:39:GLN:O	1:H:39:GLN:HG3	2.07	0.45
1:H:88:LEU:HD12	1:H:91:MET:HE3	1.98	0.45
1:A:43:LEU:O	1:A:46:SER:N	2.50	0.45
1:F:55:PHE:HB2	1:F:68:ILE:CD1	2.47	0.45
1:G:132:LEU:HG	1:H:48:LEU:CD1	2.39	0.45
1:A:63:TYR:O	1:A:64:GLN:CB	2.64	0.45
1:G:35:VAL:O	1:G:39:GLN:HB2	2.17	0.45
1:A:128:PRO:HA	2:A:274:HOH:O	2.17	0.44
1:E:88:LEU:CD1	1:E:91:MET:HE3	2.45	0.44
1:A:143:ILE:HG22	1:A:144:VAL:HG23	1.99	0.44
1:C:55:PHE:HB2	1:C:68:ILE:HD13	2.00	0.44
1:E:184:SER:O	1:E:190:LYS:HB2	2.17	0.44
1:G:68:ILE:HG22	1:G:72:LEU:HD22	1.99	0.44
1:A:92:VAL:O	1:A:96:VAL:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:110:ASN:O	1:F:110:ASN:CG	2.60	0.44
1:G:48:LEU:HD23	1:H:91:MET:CG	2.40	0.44
1:C:38:TRP:CH2	1:C:51:GLN:HA	2.52	0.44
1:C:82:ASP:OD1	1:C:85:GLU:HG3	2.18	0.44
1:D:72:LEU:O	1:D:76:LEU:HB2	2.17	0.44
1:E:133:LEU:HD12	1:E:133:LEU:HA	1.82	0.44
1:G:142:PHE:HZ	1:G:186:ARG:NH1	2.16	0.44
1:C:15:ALA:O	1:C:19:MET:HG3	2.18	0.44
1:F:115:LYS:O	1:F:119:VAL:HG23	2.17	0.44
1:A:54:LYS:HG3	1:A:63:TYR:CE1	2.53	0.44
1:D:63:TYR:O	1:D:64:GLN:CB	2.61	0.44
1:D:137:GLN:OE1	1:D:141:THR:HG21	2.18	0.44
1:H:19:MET:HE1	1:H:157:ASP:HB2	2.00	0.44
1:A:110:ASN:O	1:A:110:ASN:CG	2.60	0.44
1:C:85:GLU:HB3	1:C:144:VAL:CG1	2.43	0.44
1:C:115:LYS:O	1:C:119:VAL:HG23	2.18	0.44
1:B:92:VAL:HG22	1:B:132:LEU:HD13	2.00	0.43
1:D:13:ARG:HD3	2:D:244:HOH:O	2.18	0.43
1:D:88:LEU:CG	1:D:144:VAL:HG22	2.47	0.43
1:D:88:LEU:HD23	1:D:144:VAL:HG22	1.99	0.43
1:D:111:TYR:CZ	1:D:115:LYS:HD2	2.53	0.43
1:E:110:ASN:O	1:E:110:ASN:CG	2.59	0.43
1:E:72:LEU:O	1:E:76:LEU:HB2	2.18	0.43
1:B:209:GLN:HB3	2:B:266:HOH:O	2.18	0.43
1:C:9:PRO:HD3	1:C:32:VAL:HG13	2.00	0.43
1:E:110:ASN:N	1:E:110:ASN:HD22	2.17	0.43
1:G:110:ASN:HD22	1:G:110:ASN:N	2.16	0.43
1:B:31:GLU:HG3	2:B:279:HOH:O	2.17	0.43
1:G:20:LEU:HD23	1:G:20:LEU:C	2.43	0.43
1:G:88:LEU:HD12	1:G:91:MET:HE3	1.99	0.43
1:E:15:ALA:O	1:E:19:MET:HG3	2.19	0.43
1:E:127:LYS:HB3	1:E:128:PRO:HD3	2.00	0.43
1:G:83:GLN:OE1	1:H:75:THR:HG22	2.19	0.43
1:A:148:ILE:HD13	1:A:153:TYR:CZ	2.54	0.43
1:B:122:LEU:HD22	1:B:126:LEU:HG	2.01	0.43
1:D:107:ILE:O	1:D:206:ASN:ND2	2.52	0.43
1:F:11:ARG:HB3	1:F:15:ALA:HB2	2.00	0.43
1:F:142:PHE:CE2	1:F:146:ASP:O	2.72	0.43
1:G:43:LEU:HA	1:G:46:SER:OG	2.18	0.43
1:H:63:TYR:O	1:H:64:GLN:CB	2.67	0.43
1:E:111:TYR:CZ	1:E:115:LYS:HD2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:107:ILE:O	1:F:206:ASN:ND2	2.49	0.43
1:A:109:THR:O	1:A:110:ASN:HB3	2.18	0.43
1:B:30:GLU:HG3	2:B:239:HOH:O	2.18	0.43
1:D:110:ASN:N	1:D:110:ASN:HD22	2.16	0.43
2:G:223:HOH:O	1:H:90:ASP:HB3	2.18	0.43
1:C:56:GLN:HB3	2:C:226:HOH:O	2.18	0.43
1:D:4:THR:HB	1:D:56:GLN:HB3	2.01	0.43
1:A:49:TYR:O	1:A:51:GLN:HG3	2.20	0.42
1:D:110:ASN:O	1:D:110:ASN:ND2	2.52	0.42
1:E:207:GLY:HA2	2:E:259:HOH:O	2.18	0.42
1:G:60:LEU:HD22	1:G:62:LEU:HD13	2.01	0.42
1:A:133:LEU:CD1	1:A:143:ILE:HG13	2.48	0.42
1:B:175:LEU:HD12	1:B:175:LEU:HA	1.77	0.42
1:G:55:PHE:HB2	1:G:68:ILE:CD1	2.49	0.42
1:H:107:ILE:O	1:H:206:ASN:ND2	2.52	0.42
1:C:34:THR:C	1:C:36:GLU:N	2.76	0.42
1:C:63:TYR:O	1:C:64:GLN:CB	2.67	0.42
1:G:49:TYR:CD1	1:H:128:PRO:CB	2.99	0.42
1:A:72:LEU:HD12	1:A:72:LEU:HA	1.88	0.42
1:C:116:ASP:O	1:C:120:LYS:HG3	2.20	0.42
1:A:195:SER:O	1:A:199:VAL:HG23	2.19	0.42
1:B:110:ASN:N	1:B:110:ASN:HD22	2.16	0.42
1:E:204:ASN:ND2	2:E:229:HOH:O	2.44	0.42
1:G:76:LEU:HD12	1:G:76:LEU:HA	1.91	0.42
1:A:55:PHE:HB2	1:A:68:ILE:HD13	2.02	0.42
1:A:131:THR:O	1:A:135:GLN:HG3	2.20	0.42
1:C:60:LEU:HD22	1:C:62:LEU:CD1	2.50	0.42
1:F:36:GLU:CD	1:F:37:THR:N	2.77	0.42
1:H:88:LEU:HD11	1:H:132:LEU:HD22	2.01	0.42
1:C:20:LEU:C	1:C:20:LEU:HD23	2.45	0.42
1:E:48:LEU:HD22	1:F:91:MET:HG2	2.02	0.42
1:E:110:ASN:O	1:E:110:ASN:ND2	2.53	0.42
1:F:88:LEU:HD23	1:F:144:VAL:HG22	2.02	0.42
1:B:111:TYR:CZ	1:B:115:LYS:HD2	2.55	0.41
1:C:11:ARG:O	1:C:12:GLY:C	2.62	0.41
1:C:81:LYS:HG2	1:C:85:GLU:CD	2.45	0.41
1:H:89:VAL:HG23	1:H:144:VAL:HG11	2.02	0.41
1:C:34:THR:OG1	1:C:36:GLU:HB2	2.20	0.41
1:C:175:LEU:HD12	1:C:175:LEU:HA	1.88	0.41
1:D:6:VAL:HG12	1:D:33:VAL:HG21	2.02	0.41
1:E:91:MET:HG2	1:F:48:LEU:CD2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:72:LEU:HD12	1:H:72:LEU:HA	1.88	0.41
1:F:188:LYS:NZ	2:F:274:HOH:O	2.46	0.41
1:A:92:VAL:HG21	1:A:143:ILE:HG21	2.03	0.41
1:B:39:GLN:O	1:B:40:GLU:C	2.63	0.41
1:C:133:LEU:HD22	1:C:176:LEU:HD23	2.02	0.41
1:G:107:ILE:O	1:G:206:ASN:ND2	2.50	0.41
1:A:175:LEU:HD12	1:A:175:LEU:HA	1.85	0.41
1:C:130:GLU:CD	1:C:175:LEU:HB2	2.43	0.41
1:D:111:TYR:O	1:D:115:LYS:HB2	2.21	0.41
1:D:122:LEU:HD22	1:D:126:LEU:HG	2.02	0.41
1:F:130:GLU:OE1	1:F:176:LEU:HG	2.20	0.41
1:G:110:ASN:O	1:G:110:ASN:ND2	2.54	0.41
1:A:85:GLU:O	1:A:89:VAL:HG23	2.21	0.41
1:C:133:LEU:HD12	1:C:133:LEU:HA	1.91	0.41
1:G:189:LEU:CD2	1:G:193:LEU:HG	2.50	0.41
1:C:110:ASN:HD22	1:C:110:ASN:N	2.18	0.41
1:E:52:LEU:HB3	1:E:53:PRO:HA	2.01	0.41
1:G:109:THR:O	1:G:110:ASN:HB3	2.21	0.41
1:B:72:LEU:O	1:B:76:LEU:HB2	2.21	0.41
1:C:201:LEU:HD13	2:C:255:HOH:O	2.21	0.41
1:E:38:TRP:CZ2	1:E:44:LYS:HG3	2.56	0.41
1:H:73:GLY:HA2	1:H:78:LEU:HB2	2.03	0.41
1:B:63:TYR:O	1:B:64:GLN:CB	2.67	0.41
1:D:41:GLY:O	1:D:42:SER:C	2.64	0.41
1:F:88:LEU:HD23	1:F:144:VAL:HG13	2.03	0.41
1:H:16:ALA:HA	1:H:19:MET:HE3	2.03	0.41
1:A:91:MET:HG2	1:B:48:LEU:HD23	2.03	0.41
1:A:122:LEU:HD22	1:A:126:LEU:CG	2.51	0.41
1:E:72:LEU:HD12	1:E:72:LEU:HA	1.90	0.41
1:H:85:GLU:OE2	1:H:145:GLY:HA2	2.20	0.41
1:H:136:ASN:O	1:H:137:GLN:C	2.64	0.41
1:C:137:GLN:OE1	1:C:141:THR:HG21	2.21	0.40
1:E:60:LEU:HG	2:F:224:HOH:O	2.20	0.40
1:F:88:LEU:HD12	1:F:91:MET:HE3	2.03	0.40
1:G:110:ASN:O	1:G:110:ASN:CG	2.64	0.40
1:C:110:ASN:O	1:C:110:ASN:CG	2.63	0.40
1:A:99:LEU:HD23	1:A:158:LEU:HD21	2.04	0.40
1:B:55:PHE:HB2	1:B:68:ILE:HD13	2.03	0.40
1:G:40:GLU:OE1	1:G:42:SER:HB3	2.22	0.40
1:H:100:ARG:HH21	1:H:154:ASN:ND2	2.18	0.40
1:A:88:LEU:HG	1:A:144:VAL:HG22	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:LYS:HD3	2:B:298:HOH:O	2.21	0.40
1:B:176:LEU:O	1:B:179:TYR:HB3	2.21	0.40
1:C:186:ARG:HA	1:C:187:PRO:HD3	1.95	0.40
1:D:14:CYS:SG	1:D:53:PRO:HB3	2.61	0.40
1:D:160:LEU:O	1:D:163:GLU:HB2	2.21	0.40
1:F:133:LEU:HD21	1:F:179:TYR:HB2	2.02	0.40
1:A:176:LEU:O	1:A:179:TYR:HB3	2.21	0.40
1:B:127:LYS:HB3	1:B:128:PRO:HD3	2.04	0.40
1:C:49:TYR:CD1	1:D:128:PRO:HB3	2.56	0.40
1:G:5:VAL:HG22	1:G:30:GLU:OE1	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:40:GLU:CG	2:F:215:HOH:O[1_565]	2.02	0.18

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	207/209 (99%)	198 (96%)	6 (3%)	3 (1%)	9	17
1	B	207/209 (99%)	196 (95%)	10 (5%)	1 (0%)	24	43
1	C	207/209 (99%)	196 (95%)	7 (3%)	4 (2%)	6	11
1	D	207/209 (99%)	192 (93%)	9 (4%)	6 (3%)	3	5
1	E	207/209 (99%)	199 (96%)	5 (2%)	3 (1%)	9	17
1	F	207/209 (99%)	199 (96%)	4 (2%)	4 (2%)	6	11
1	G	207/209 (99%)	198 (96%)	6 (3%)	3 (1%)	9	17
1	H	207/209 (99%)	197 (95%)	6 (3%)	4 (2%)	6	11

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1656/1672 (99%)	1575 (95%)	53 (3%)	28 (2%)	7	13

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	141	THR
1	D	48	LEU
1	F	35	VAL
1	G	42	SER
1	H	141	THR
1	C	145	GLY
1	A	110	ASN
1	D	40	GLU
1	D	141	THR
1	E	110	ASN
1	F	110	ASN
1	B	110	ASN
1	C	110	ASN
1	D	49	TYR
1	E	64	GLN
1	F	141	THR
1	G	64	GLN
1	G	110	ASN
1	H	64	GLN
1	A	64	GLN
1	C	64	GLN
1	D	64	GLN
1	F	64	GLN
1	H	110	ASN
1	D	187	PRO
1	E	187	PRO
1	H	187	PRO
1	A	187	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	175/175 (100%)	160 (91%)	15 (9%)	10	21
1	B	175/175 (100%)	162 (93%)	13 (7%)	13	27
1	C	175/175 (100%)	163 (93%)	12 (7%)	14	30
1	D	175/175 (100%)	161 (92%)	14 (8%)	11	24
1	E	175/175 (100%)	162 (93%)	13 (7%)	13	27
1	F	175/175 (100%)	161 (92%)	14 (8%)	11	24
1	G	175/175 (100%)	158 (90%)	17 (10%)	8	17
1	H	175/175 (100%)	157 (90%)	18 (10%)	7	14
All	All	1400/1400 (100%)	1284 (92%)	116 (8%)	10	22

All (116) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	VAL
1	A	36	GLU
1	A	43	LEU
1	A	62	LEU
1	A	72	LEU
1	A	76	LEU
1	A	109	THR
1	A	110	ASN
1	A	122	LEU
1	A	143	ILE
1	A	146	ASP
1	A	147	GLN
1	A	170	LEU
1	A	175	LEU
1	A	189	LEU
1	B	5	VAL
1	B	36	GLU
1	B	43	LEU
1	B	62	LEU
1	B	72	LEU
1	B	76	LEU
1	B	109	THR
1	B	110	ASN
1	B	122	LEU
1	B	143	ILE
1	B	170	LEU
1	B	175	LEU

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Mol	Chain	Res	Type
1	B	189	LEU
1	C	5	VAL
1	C	43	LEU
1	C	62	LEU
1	C	72	LEU
1	C	76	LEU
1	C	109	THR
1	C	110	ASN
1	C	122	LEU
1	C	143	ILE
1	C	170	LEU
1	C	175	LEU
1	C	189	LEU
1	D	5	VAL
1	D	35	VAL
1	D	36	GLU
1	D	40	GLU
1	D	43	LEU
1	D	62	LEU
1	D	72	LEU
1	D	76	LEU
1	D	109	THR
1	D	110	ASN
1	D	122	LEU
1	D	170	LEU
1	D	175	LEU
1	D	189	LEU
1	E	5	VAL
1	E	31	GLU
1	E	43	LEU
1	E	62	LEU
1	E	72	LEU
1	E	76	LEU
1	E	109	THR
1	E	110	ASN
1	E	122	LEU
1	E	133	LEU
1	E	170	LEU
1	E	175	LEU
1	E	189	LEU
1	F	5	VAL
1	F	36	GLU

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Mol	Chain	Res	Type
1	F	37	THR
1	F	43	LEU
1	F	62	LEU
1	F	72	LEU
1	F	76	LEU
1	F	109	THR
1	F	110	ASN
1	F	122	LEU
1	F	141	THR
1	F	170	LEU
1	F	175	LEU
1	F	189	LEU
1	G	5	VAL
1	G	34	THR
1	G	43	LEU
1	G	60	LEU
1	G	62	LEU
1	G	72	LEU
1	G	76	LEU
1	G	91	MET
1	G	109	THR
1	G	110	ASN
1	G	122	LEU
1	G	133	LEU
1	G	146	ASP
1	G	147	GLN
1	G	170	LEU
1	G	175	LEU
1	G	189	LEU
1	H	5	VAL
1	H	36	GLU
1	H	39	GLN
1	H	43	LEU
1	H	60	LEU
1	H	62	LEU
1	H	72	LEU
1	H	76	LEU
1	H	91	MET
1	H	109	THR
1	H	110	ASN
1	H	122	LEU
1	H	141	THR

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Mol	Chain	Res	Type
1	H	146	ASP
1	H	147	GLN
1	H	170	LEU
1	H	175	LEU
1	H	189	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (79) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	39	GLN
1	A	56	GLN
1	A	66	ASN
1	A	83	GLN
1	A	93	ASN
1	A	110	ASN
1	A	125	GLN
1	A	136	ASN
1	A	147	GLN
1	A	154	ASN
1	B	24	GLN
1	B	56	GLN
1	B	66	ASN
1	B	83	GLN
1	B	93	ASN
1	B	110	ASN
1	B	125	GLN
1	B	136	ASN
1	B	154	ASN
1	C	24	GLN
1	C	56	GLN
1	C	66	ASN
1	C	93	ASN
1	C	110	ASN
1	C	125	GLN
1	C	136	ASN
1	C	154	ASN
1	C	204	ASN
1	D	24	GLN
1	D	56	GLN
1	D	66	ASN
1	D	93	ASN

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Mol	Chain	Res	Type
1	D	110	ASN
1	D	125	GLN
1	D	154	ASN
1	E	24	GLN
1	E	56	GLN
1	E	66	ASN
1	E	83	GLN
1	E	93	ASN
1	E	110	ASN
1	E	125	GLN
1	E	136	ASN
1	E	147	GLN
1	E	154	ASN
1	F	24	GLN
1	F	51	GLN
1	F	56	GLN
1	F	66	ASN
1	F	93	ASN
1	F	110	ASN
1	F	136	ASN
1	F	147	GLN
1	F	154	ASN
1	F	200	ASN
1	F	209	GLN
1	G	24	GLN
1	G	56	GLN
1	G	66	ASN
1	G	93	ASN
1	G	110	ASN
1	G	135	GLN
1	G	147	GLN
1	G	154	ASN
1	G	209	GLN
1	H	24	GLN
1	H	39	GLN
1	H	56	GLN
1	H	66	ASN
1	H	83	GLN
1	H	93	ASN
1	H	110	ASN
1	H	125	GLN
1	H	136	ASN

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Mol	Chain	Res	Type
1	H	137	GLN
1	H	154	ASN
1	H	204	ASN
1	H	209	GLN

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	209/209 (100%)	-1.57	0 100 100	11, 27, 45, 58	0
1	B	209/209 (100%)	-1.57	0 100 100	9, 27, 39, 50	0
1	C	209/209 (100%)	-1.44	0 100 100	17, 30, 65, 79	0
1	D	209/209 (100%)	-1.47	0 100 100	15, 29, 66, 75	0
1	E	209/209 (100%)	-1.53	0 100 100	10, 28, 52, 69	0
1	F	209/209 (100%)	-1.53	0 100 100	13, 27, 45, 58	0
1	G	209/209 (100%)	-1.44	0 100 100	14, 29, 67, 76	0
1	H	209/209 (100%)	-1.47	0 100 100	14, 31, 62, 81	0
All	All	1672/1672 (100%)	-1.50	0 100 100	9, 28, 56, 81	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.