



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:22 PM UTC

PDB ID : 3O3W / pdb_00003o3w
Title : Crystal Structure of BH2092 protein (residues 14-131) from *Bacillus halodurans*, Northeast Structural Genomics Consortium Target BhR228A
Authors : Forouhar, F.; Neely, H.; Seetharaman, J.; Sahdev, S.; Xiao, R.; Ciccocanti, C.; Lee, D.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2010-07-26
Resolution : 2.91 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

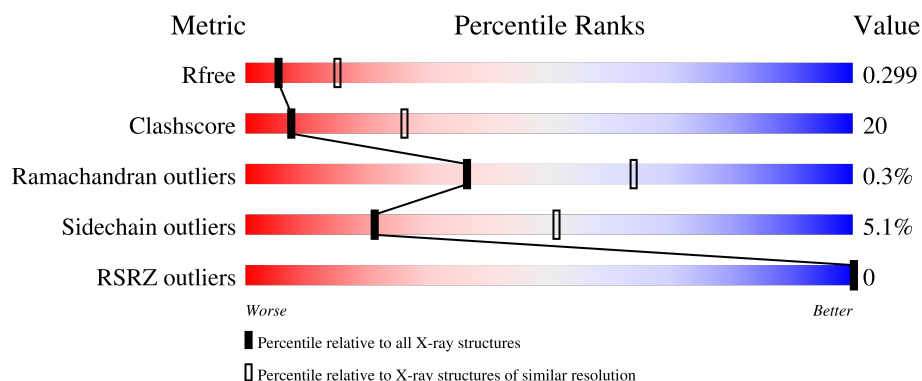
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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	126	 60% 31% • 6%
1	B	126	 60% 31% • 7%
1	C	126	 58% 31% 5% 6%
1	D	126	 60% 28% • 10%

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Mol	Chain	Length	Quality of chain
1	E	126	52%37%5%6%
1	F	126	60%29%5%6%
1	G	126	56%32%10%
1	H	126	64%25%8%
1	I	126	63%25%10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8353 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 BH2092 protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	118	Total	C	N	O	S	Se	0	0	0
			928	591	159	174	3	1			
1	B	117	Total	C	N	O	S	Se	0	0	0
			928	591	158	175	3	1			
1	C	118	Total	C	N	O	S	Se	0	0	0
			933	594	159	176	3	1			
1	D	114	Total	C	N	O	S	Se	0	0	0
			907	578	155	170	3	1			
1	E	118	Total	C	N	O	S	Se	0	0	0
			933	594	159	176	3	1			
1	F	118	Total	C	N	O	S	Se	0	0	0
			929	591	161	173	3	1			
1	G	114	Total	C	N	O	S	Se	0	0	0
			907	578	155	170	3	1			
1	H	116	Total	C	N	O	S	Se	0	0	0
			919	586	157	172	3	1			
1	I	113	Total	C	N	O	S	Se	0	0	0
			890	567	153	166	3	1			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	132	LEU	-	expression tag	UNP Q9KB42
A	133	GLU	-	expression tag	UNP Q9KB42
A	134	HIS	-	expression tag	UNP Q9KB42
A	135	HIS	-	expression tag	UNP Q9KB42
A	136	HIS	-	expression tag	UNP Q9KB42
A	137	HIS	-	expression tag	UNP Q9KB42
A	138	HIS	-	expression tag	UNP Q9KB42
A	139	HIS	-	expression tag	UNP Q9KB42
B	132	LEU	-	expression tag	UNP Q9KB42
B	133	GLU	-	expression tag	UNP Q9KB42
B	134	HIS	-	expression tag	UNP Q9KB42

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Chain	Residue	Modelled	Actual	Comment	Reference
B	135	HIS	-	expression tag	UNP Q9KB42
B	136	HIS	-	expression tag	UNP Q9KB42
B	137	HIS	-	expression tag	UNP Q9KB42
B	138	HIS	-	expression tag	UNP Q9KB42
B	139	HIS	-	expression tag	UNP Q9KB42
C	132	LEU	-	expression tag	UNP Q9KB42
C	133	GLU	-	expression tag	UNP Q9KB42
C	134	HIS	-	expression tag	UNP Q9KB42
C	135	HIS	-	expression tag	UNP Q9KB42
C	136	HIS	-	expression tag	UNP Q9KB42
C	137	HIS	-	expression tag	UNP Q9KB42
C	138	HIS	-	expression tag	UNP Q9KB42
C	139	HIS	-	expression tag	UNP Q9KB42
D	132	LEU	-	expression tag	UNP Q9KB42
D	133	GLU	-	expression tag	UNP Q9KB42
D	134	HIS	-	expression tag	UNP Q9KB42
D	135	HIS	-	expression tag	UNP Q9KB42
D	136	HIS	-	expression tag	UNP Q9KB42
D	137	HIS	-	expression tag	UNP Q9KB42
D	138	HIS	-	expression tag	UNP Q9KB42
D	139	HIS	-	expression tag	UNP Q9KB42
E	132	LEU	-	expression tag	UNP Q9KB42
E	133	GLU	-	expression tag	UNP Q9KB42
E	134	HIS	-	expression tag	UNP Q9KB42
E	135	HIS	-	expression tag	UNP Q9KB42
E	136	HIS	-	expression tag	UNP Q9KB42
E	137	HIS	-	expression tag	UNP Q9KB42
E	138	HIS	-	expression tag	UNP Q9KB42
E	139	HIS	-	expression tag	UNP Q9KB42
F	132	LEU	-	expression tag	UNP Q9KB42
F	133	GLU	-	expression tag	UNP Q9KB42
F	134	HIS	-	expression tag	UNP Q9KB42
F	135	HIS	-	expression tag	UNP Q9KB42
F	136	HIS	-	expression tag	UNP Q9KB42
F	137	HIS	-	expression tag	UNP Q9KB42
F	138	HIS	-	expression tag	UNP Q9KB42
F	139	HIS	-	expression tag	UNP Q9KB42
G	132	LEU	-	expression tag	UNP Q9KB42
G	133	GLU	-	expression tag	UNP Q9KB42
G	134	HIS	-	expression tag	UNP Q9KB42
G	135	HIS	-	expression tag	UNP Q9KB42
G	136	HIS	-	expression tag	UNP Q9KB42

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Chain	Residue	Modelled	Actual	Comment	Reference
G	137	HIS	-	expression tag	UNP Q9KB42
G	138	HIS	-	expression tag	UNP Q9KB42
G	139	HIS	-	expression tag	UNP Q9KB42
H	132	LEU	-	expression tag	UNP Q9KB42
H	133	GLU	-	expression tag	UNP Q9KB42
H	134	HIS	-	expression tag	UNP Q9KB42
H	135	HIS	-	expression tag	UNP Q9KB42
H	136	HIS	-	expression tag	UNP Q9KB42
H	137	HIS	-	expression tag	UNP Q9KB42
H	138	HIS	-	expression tag	UNP Q9KB42
H	139	HIS	-	expression tag	UNP Q9KB42
I	132	LEU	-	expression tag	UNP Q9KB42
I	133	GLU	-	expression tag	UNP Q9KB42
I	134	HIS	-	expression tag	UNP Q9KB42
I	135	HIS	-	expression tag	UNP Q9KB42
I	136	HIS	-	expression tag	UNP Q9KB42
I	137	HIS	-	expression tag	UNP Q9KB42
I	138	HIS	-	expression tag	UNP Q9KB42
I	139	HIS	-	expression tag	UNP Q9KB42

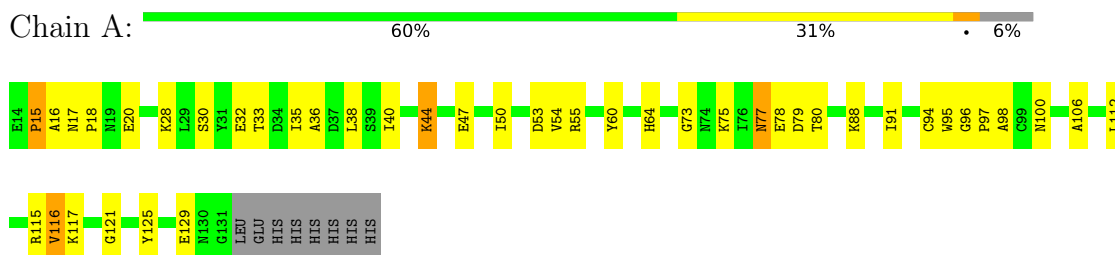
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	14	Total O 14 14	0	0
2	B	9	Total O 9 9	0	0
2	C	6	Total O 6 6	0	0
2	D	10	Total O 10 10	0	0
2	E	11	Total O 11 11	0	0
2	F	7	Total O 7 7	0	0
2	G	10	Total O 10 10	0	0
2	H	2	Total O 2 2	0	0
2	I	10	Total O 10 10	0	0

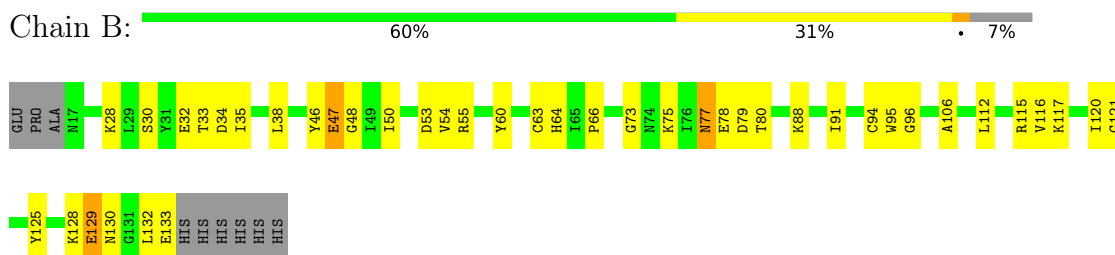
3 Residue-property plots

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.

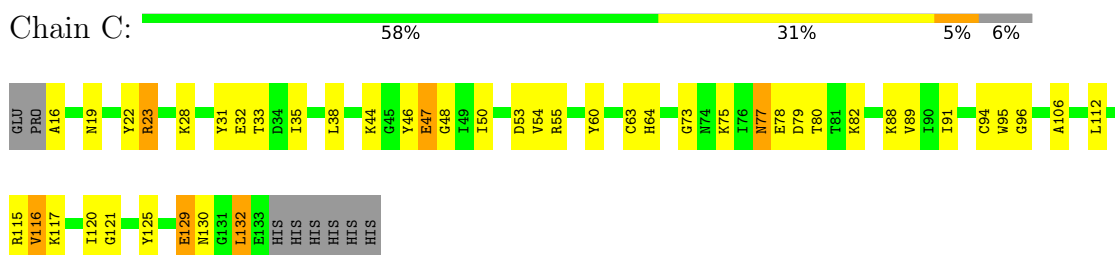
• Molecule 1: BH2092 protein



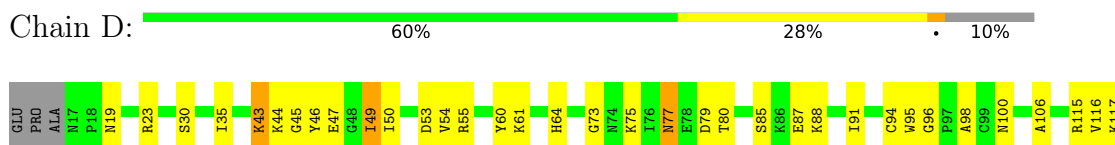
• Molecule 1: BH2092 protein

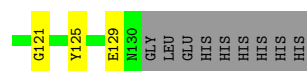


• Molecule 1: BH2092 protein



• Molecule 1: BH2092 protein





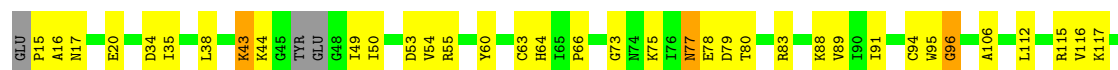
• Molecule 1: BH2092 protein

Chain E: 52% 37% 5% 6%



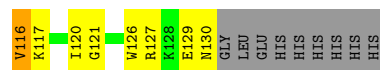
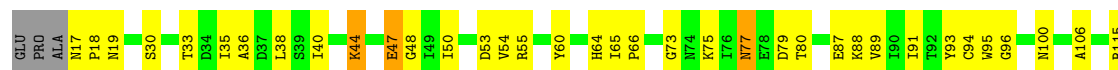
• Molecule 1: BH2092 protein

Chain F: 60% 29% 5% 6%



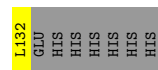
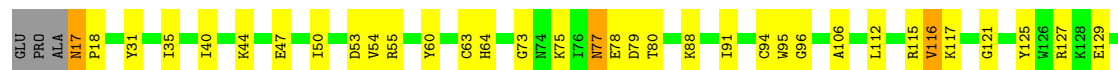
• Molecule 1: BH2092 protein

Chain G: 56% 32% 10% 2%



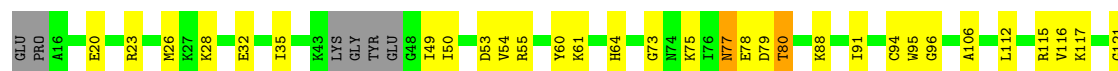
• Molecule 1: BH2092 protein

Chain H: 64% 25% 8% 3%



• Molecule 1: BH2092 protein

Chain I: 63% 25% 10% 2%



R127	R128	E129	L132	GLU	HIS	HIS	HIS	HIS	HIS	HIS
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4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	114.24Å 114.24Å 85.45Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.79 – 2.91 19.79 – 2.91	Depositor EDS
% Data completeness (in resolution range)	93.9 (19.79-2.91) 98.5 (19.79-2.91)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	0.17	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.90Å)	Xtriage
Refinement program	CNS 1.2, REFMAC	Depositor
R, R_{free}	0.250 , 0.292 0.261 , 0.299	Depositor DCC
R_{free} test set	5410 reflections (9.88%)	wwPDB-VP
Wilson B-factor (Å ²)	33.1	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 4.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.098 for -h,-k,l 0.099 for h,-h-k,-l 0.286 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8353	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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	0.56	0/946	0.90	1/1275 (0.1%)
1	B	0.57	0/945	0.86	0/1272
1	C	0.58	1/950 (0.1%)	1.12	4/1279 (0.3%)
1	D	0.55	0/924	0.85	1/1244 (0.1%)
1	E	0.58	0/950	0.90	5/1279 (0.4%)
1	F	0.55	0/946	0.89	2/1272 (0.2%)
1	G	0.50	0/924	0.91	3/1244 (0.2%)
1	H	0.47	0/936	0.87	3/1260 (0.2%)
1	I	0.50	0/905	0.84	0/1218
All	All	0.54	1/8426 (0.0%)	0.91	19/11343 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	23	ARG	CD-NE	-5.21	1.39	1.46

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	23	ARG	NE-CZ-NH2	15.46	133.12	119.20
1	C	23	ARG	NE-CZ-NH1	-14.90	106.60	121.50
1	C	23	ARG	CD-NE-CZ	12.42	141.78	124.40
1	G	17	ASN	CA-C-N	8.76	130.78	119.84
1	G	17	ASN	C-N-CA	8.76	130.78	119.84
1	D	44	LYS	N-CA-C	-6.04	105.89	113.20
1	A	116	VAL	N-CA-C	5.87	116.45	110.05
1	E	116	VAL	N-CA-C	5.60	116.15	110.05
1	E	17	ASN	CA-C-N	-5.31	113.72	119.24
1	E	17	ASN	C-N-CA	-5.31	113.72	119.24
1	G	116	VAL	N-CA-C	5.29	115.81	110.05
1	F	96	GLY	CA-C-N	5.22	125.02	119.28
1	F	96	GLY	C-N-CA	5.22	125.02	119.28
1	H	17	ASN	CA-C-N	5.19	126.33	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	17	ASN	C-N-CA	5.19	126.33	119.84
1	C	116	VAL	N-CA-C	5.13	115.64	110.05
1	E	96	GLY	CA-C-N	5.06	124.72	119.56
1	E	96	GLY	C-N-CA	5.06	124.72	119.56
1	H	116	VAL	N-CA-C	5.00	115.50	110.05

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	928	0	927	45	0
1	B	928	0	930	35	0
1	C	933	0	935	60	1
1	D	907	0	910	44	0
1	E	933	0	935	63	0
1	F	929	0	934	47	1
1	G	907	0	910	36	0
1	H	919	0	924	29	0
1	I	890	0	897	42	0
2	A	14	0	0	1	0
2	B	9	0	0	2	0
2	C	6	0	0	1	0
2	D	10	0	0	1	0
2	E	11	0	0	4	0
2	F	7	0	0	2	0
2	G	10	0	0	4	0
2	H	2	0	0	0	0
2	I	10	0	0	6	0
All	All	8353	0	8302	337	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 20.

All (337) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:78:GLU:OE1	1:E:26:MSE:HE2	1.37	1.24
1:I:49:ILE:HD12	1:I:91:ILE:HD12	1.37	1.07
1:E:115:ARG:HH12	1:F:15:PRO:HB2	1.17	1.05
1:D:23:ARG:NH2	1:I:20:GLU:HB2	1.73	1.03
1:G:35:ILE:HG21	1:G:129:GLU:HG3	1.42	0.98
1:B:28:LYS:HD2	2:B:144:HOH:O	1.66	0.96
1:A:16:ALA:N	1:C:115:ARG:HH12	1.65	0.95
1:D:23:ARG:NH2	1:I:20:GLU:CB	2.31	0.92
1:A:15:PRO:C	1:C:115:ARG:HH12	1.79	0.91
1:E:115:ARG:HH22	1:F:15:PRO:C	1.81	0.89
1:D:23:ARG:HH21	1:I:20:GLU:CB	1.87	0.86
1:G:100:ASN:HB2	2:G:143:HOH:O	1.75	0.86
1:E:115:ARG:HH12	1:F:15:PRO:CB	1.91	0.84
1:F:43:LYS:HD3	1:F:44:LYS:N	1.91	0.83
1:B:77:ASN:HD22	1:B:79:ASP:H	1.26	0.83
1:E:94:CYS:HB3	2:E:9:HOH:O	1.79	0.83
1:D:77:ASN:HD22	1:D:79:ASP:H	1.26	0.83
1:C:77:ASN:HD22	1:C:79:ASP:H	1.28	0.82
1:C:35:ILE:HG21	1:C:129:GLU:HG3	1.59	0.82
1:I:77:ASN:HD22	1:I:79:ASP:H	1.28	0.82
1:C:78:GLU:CD	1:E:26:MSE:HE2	2.05	0.80
1:F:77:ASN:HD22	1:F:79:ASP:H	1.30	0.80
1:B:35:ILE:HG21	1:B:129:GLU:HG3	1.64	0.80
1:E:77:ASN:HD22	1:E:79:ASP:H	1.29	0.79
1:I:35:ILE:HG21	1:I:129:GLU:HG3	1.64	0.79
1:B:129:GLU:HB2	1:B:130:ASN:HD22	1.47	0.78
1:E:35:ILE:HG21	1:E:129:GLU:HG3	1.63	0.78
1:G:77:ASN:HD22	1:G:79:ASP:H	1.30	0.78
1:H:77:ASN:HD22	1:H:79:ASP:H	1.30	0.77
1:A:16:ALA:N	1:C:115:ARG:NH1	2.32	0.77
1:A:77:ASN:HD22	1:A:79:ASP:H	1.31	0.77
1:E:115:ARG:NH1	1:F:15:PRO:HB2	1.98	0.77
1:D:23:ARG:NH2	1:I:20:GLU:CG	2.47	0.76
1:I:26:MSE:HE3	2:I:148:HOH:O	1.84	0.76
1:A:15:PRO:C	1:C:115:ARG:HH22	1.94	0.74
1:H:35:ILE:HG21	1:H:129:GLU:HG3	1.68	0.74
1:C:82:LYS:NZ	1:E:19:ASN:HD22	1.85	0.73
1:F:129:GLU:HA	2:F:140:HOH:O	1.89	0.73
1:C:78:GLU:OE1	1:E:26:MSE:CE	2.29	0.72
1:D:35:ILE:HG21	1:D:129:GLU:HG3	1.71	0.72
1:E:63:CYS:SG	1:E:132:LEU:HD13	2.29	0.71
1:F:35:ILE:HG21	1:F:129:GLU:HG3	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:CYS:SG	1:B:132:LEU:HD13	2.31	0.71
1:A:35:ILE:HG21	1:A:129:GLU:HG3	1.72	0.70
1:H:17:ASN:HD22	1:H:18:PRO:HD2	1.55	0.70
1:I:127:ARG:HG2	1:I:132:LEU:HD21	1.74	0.68
1:H:17:ASN:HD22	1:H:18:PRO:CD	2.07	0.68
1:E:30:SER:O	1:F:15:PRO:HB3	1.92	0.68
1:A:106:ALA:HA	1:A:116:VAL:HG21	1.76	0.68
1:E:44:LYS:HE3	1:E:44:LYS:HA	1.75	0.68
1:C:79:ASP:OD2	1:E:23:ARG:HG2	1.94	0.68
1:I:106:ALA:HA	1:I:116:VAL:HG21	1.76	0.67
1:D:23:ARG:HH21	1:I:20:GLU:HB2	1.51	0.67
1:G:106:ALA:HA	1:G:116:VAL:HG21	1.77	0.67
1:H:47:GLU:O	1:H:88:LYS:HE3	1.95	0.67
1:A:15:PRO:CB	1:C:115:ARG:HH12	2.09	0.66
1:D:106:ALA:HA	1:D:116:VAL:HG21	1.76	0.66
1:E:47:GLU:O	1:E:88:LYS:HE3	1.95	0.66
1:B:106:ALA:HA	1:B:116:VAL:HG21	1.77	0.66
1:H:106:ALA:HA	1:H:116:VAL:HG21	1.77	0.66
1:E:106:ALA:HA	1:E:116:VAL:HG21	1.78	0.66
1:G:47:GLU:O	1:G:88:LYS:HE3	1.95	0.65
1:A:15:PRO:C	1:C:115:ARG:NH1	2.53	0.65
1:A:15:PRO:HB2	1:C:115:ARG:HH12	1.61	0.65
1:D:23:ARG:NH2	1:I:20:GLU:HG3	2.12	0.65
1:F:106:ALA:HA	1:F:116:VAL:HG21	1.78	0.65
1:B:77:ASN:HD22	1:B:79:ASP:N	1.95	0.64
1:G:18:PRO:HB2	2:G:147:HOH:O	1.97	0.64
1:A:15:PRO:HB2	1:C:115:ARG:NH1	2.13	0.64
1:A:15:PRO:HG2	1:C:115:ARG:NH1	2.13	0.64
1:A:47:GLU:O	1:A:88:LYS:HE3	1.97	0.64
1:D:77:ASN:HD22	1:D:79:ASP:N	1.95	0.64
1:I:94:CYS:SG	1:I:95:TRP:N	2.70	0.64
1:B:47:GLU:O	1:B:88:LYS:HE3	1.97	0.64
1:C:106:ALA:HA	1:C:116:VAL:HG21	1.80	0.64
1:G:87:GLU:HB2	2:G:142:HOH:O	1.98	0.63
1:C:82:LYS:NZ	1:E:19:ASN:ND2	2.47	0.63
1:C:77:ASN:HD22	1:C:79:ASP:N	1.97	0.62
1:H:77:ASN:HD22	1:H:79:ASP:N	1.98	0.62
1:G:44:LYS:HB2	1:G:44:LYS:NZ	2.15	0.62
1:E:77:ASN:HD22	1:E:79:ASP:N	1.98	0.62
1:E:70:SER:HB3	2:E:141:HOH:O	1.99	0.61
1:C:132:LEU:HD12	1:C:132:LEU:C	2.26	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:PRO:HA	1:B:133:GLU:OE1	2.01	0.61
1:C:82:LYS:HZ1	1:E:19:ASN:HD22	1.47	0.61
1:C:82:LYS:NZ	1:E:23:ARG:NH2	2.49	0.60
1:G:35:ILE:HG21	1:G:129:GLU:CG	2.27	0.60
1:E:66:PRO:HG3	1:E:133:GLU:HB2	1.82	0.60
1:D:77:ASN:ND2	1:D:79:ASP:HB2	2.17	0.60
1:I:77:ASN:HD22	1:I:79:ASP:N	1.98	0.60
1:G:77:ASN:HD22	1:G:79:ASP:N	1.98	0.60
1:E:33:THR:O	1:E:120:ILE:HG12	2.01	0.60
1:A:15:PRO:CB	1:C:115:ARG:NH1	2.65	0.60
1:D:30:SER:O	1:D:117:LYS:HE3	2.02	0.59
1:G:94:CYS:SG	1:G:95:TRP:N	2.74	0.59
1:D:23:ARG:HH21	1:I:20:GLU:CA	2.15	0.59
1:F:77:ASN:HD22	1:F:79:ASP:N	1.99	0.59
1:C:130:ASN:OD1	2:C:140:HOH:O	2.16	0.59
1:A:77:ASN:HD22	1:A:79:ASP:N	1.99	0.58
1:I:23:ARG:NH1	2:I:146:HOH:O	2.37	0.58
1:A:15:PRO:O	1:C:115:ARG:NH2	2.22	0.57
1:G:77:ASN:ND2	1:G:79:ASP:HB2	2.19	0.57
1:A:77:ASN:ND2	1:A:79:ASP:HB2	2.20	0.57
1:H:77:ASN:ND2	1:H:79:ASP:HB2	2.18	0.57
1:A:30:SER:O	1:A:117:LYS:HE3	2.03	0.57
1:C:33:THR:O	1:C:120:ILE:HG12	2.05	0.56
1:G:44:LYS:HB2	1:G:44:LYS:HZ2	1.70	0.56
1:F:17:ASN:ND2	1:F:20:GLU:HB2	2.20	0.56
1:F:66:PRO:HA	1:F:133:GLU:OE2	2.06	0.56
1:I:77:ASN:ND2	1:I:79:ASP:HB2	2.21	0.56
1:H:31:TYR:OH	1:I:20:GLU:HG2	2.05	0.56
1:H:17:ASN:HD22	1:H:18:PRO:N	2.04	0.56
1:C:47:GLU:O	1:C:88:LYS:HE3	2.06	0.56
1:D:23:ARG:HH22	1:I:20:GLU:CG	2.17	0.55
1:B:77:ASN:ND2	1:B:79:ASP:HB2	2.20	0.55
1:H:40:ILE:HG22	1:H:44:LYS:HD2	1.89	0.55
1:H:63:CYS:SG	1:H:132:LEU:HD13	2.46	0.55
1:F:77:ASN:ND2	1:F:79:ASP:HB2	2.21	0.55
1:F:49:ILE:HG22	1:F:89:VAL:HB	1.89	0.55
1:E:77:ASN:ND2	1:E:79:ASP:HB2	2.22	0.54
1:G:35:ILE:CG2	1:G:129:GLU:HG3	2.27	0.54
1:H:40:ILE:O	1:H:44:LYS:HG3	2.07	0.54
1:D:23:ARG:CZ	1:I:20:GLU:HB2	2.36	0.54
1:A:54:VAL:HG12	1:A:54:VAL:O	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:ASN:ND2	1:C:79:ASP:HB2	2.23	0.53
1:E:115:ARG:HH22	1:F:15:PRO:N	2.07	0.53
1:E:115:ARG:NH2	1:F:15:PRO:C	2.59	0.53
1:D:77:ASN:HD21	1:D:79:ASP:HB2	1.72	0.52
1:E:82:LYS:HB2	2:E:146:HOH:O	2.09	0.52
1:H:96:GLY:HA2	1:H:121:GLY:O	2.09	0.52
1:B:128:LYS:HA	2:B:143:HOH:O	2.10	0.52
1:B:77:ASN:HD21	1:B:79:ASP:HB2	1.74	0.52
1:F:116:VAL:HG22	1:F:117:LYS:N	2.24	0.52
1:G:30:SER:O	1:G:117:LYS:HE3	2.10	0.52
1:I:80:THR:CG2	2:I:142:HOH:O	2.57	0.52
1:C:94:CYS:SG	1:C:95:TRP:N	2.81	0.52
1:A:116:VAL:HG22	1:A:117:LYS:N	2.25	0.52
1:A:15:PRO:C	1:C:115:ARG:NH2	2.65	0.52
1:D:23:ARG:HH22	1:I:20:GLU:HG3	1.73	0.52
1:F:17:ASN:HD21	1:F:20:GLU:HB2	1.75	0.51
1:I:77:ASN:HD21	1:I:79:ASP:HB2	1.75	0.51
1:B:33:THR:O	1:B:120:ILE:HG12	2.10	0.51
1:H:17:ASN:ND2	1:H:18:PRO:HD2	2.24	0.51
1:A:98:ALA:O	1:B:34:ASP:HB2	2.10	0.51
1:H:116:VAL:HG22	1:H:117:LYS:N	2.25	0.51
1:A:77:ASN:HD21	1:A:79:ASP:HB2	1.75	0.51
1:C:54:VAL:HG12	1:C:54:VAL:O	2.09	0.51
1:G:96:GLY:HA2	1:G:121:GLY:O	2.10	0.51
1:H:17:ASN:HD22	1:H:17:ASN:C	2.17	0.51
1:H:77:ASN:HD21	1:H:79:ASP:HB2	1.74	0.51
1:H:94:CYS:SG	1:H:95:TRP:N	2.76	0.51
1:C:60:TYR:O	1:C:64:HIS:HD2	1.94	0.51
1:D:47:GLU:O	1:D:88:LYS:HE3	2.10	0.51
1:E:117:LYS:NZ	1:F:15:PRO:HG3	2.26	0.51
1:F:77:ASN:HD21	1:F:79:ASP:HB2	1.75	0.51
1:G:53:ASP:OD1	1:G:55:ARG:HD3	2.11	0.51
1:E:53:ASP:OD1	1:E:55:ARG:HD3	2.11	0.50
1:E:115:ARG:HH12	1:F:16:ALA:H	1.58	0.50
1:B:116:VAL:HG22	1:B:117:LYS:N	2.27	0.50
1:I:116:VAL:HG22	1:I:117:LYS:N	2.26	0.50
1:F:53:ASP:OD1	1:F:55:ARG:HD3	2.12	0.50
1:I:53:ASP:OD1	1:I:55:ARG:HD3	2.12	0.50
1:F:94:CYS:SG	1:F:95:TRP:N	2.75	0.49
1:C:82:LYS:HZ2	1:E:19:ASN:ND2	2.09	0.49
1:G:33:THR:O	1:G:120:ILE:HG12	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:26:MSE:CE	2:I:148:HOH:O	2.51	0.49
1:C:77:ASN:HD21	1:C:79:ASP:HB2	1.77	0.49
1:C:116:VAL:HG22	1:C:117:LYS:N	2.28	0.49
1:D:94:CYS:SG	1:D:95:TRP:N	2.80	0.49
1:D:96:GLY:HA2	1:D:121:GLY:O	2.13	0.49
1:G:77:ASN:HD21	1:G:79:ASP:HB2	1.77	0.49
1:A:15:PRO:CG	1:C:115:ARG:NH1	2.75	0.49
1:E:115:ARG:NH1	1:F:16:ALA:H	2.11	0.49
1:G:50:ILE:HD11	1:G:88:LYS:HD2	1.95	0.48
1:G:116:VAL:HG22	1:G:117:LYS:N	2.28	0.48
1:E:60:TYR:O	1:E:64:HIS:HD2	1.96	0.48
1:E:115:ARG:NH1	1:F:16:ALA:N	2.61	0.48
1:D:43:LYS:NZ	1:D:43:LYS:HB3	2.28	0.48
1:C:64:HIS:O	1:C:132:LEU:HB2	2.13	0.48
1:C:78:GLU:HG2	1:C:112:LEU:HD22	1.96	0.48
1:D:60:TYR:O	1:D:64:HIS:HD2	1.95	0.48
1:B:60:TYR:O	1:B:64:HIS:HD2	1.96	0.48
1:E:115:ARG:HH22	1:F:15:PRO:CA	2.26	0.48
1:H:53:ASP:OD1	1:H:55:ARG:HD3	2.13	0.48
1:E:54:VAL:HG12	1:E:54:VAL:O	2.13	0.48
1:H:60:TYR:O	1:H:64:HIS:HD2	1.96	0.48
1:I:96:GLY:HA2	1:I:121:GLY:O	2.14	0.48
1:E:116:VAL:HG22	1:E:117:LYS:N	2.29	0.47
1:F:60:TYR:O	1:F:64:HIS:HD2	1.97	0.47
1:A:15:PRO:CA	1:C:115:ARG:HH12	2.27	0.47
1:A:40:ILE:O	1:A:44:LYS:HG3	2.14	0.47
1:I:54:VAL:O	1:I:54:VAL:HG12	2.13	0.47
1:E:77:ASN:HD21	1:E:79:ASP:HB2	1.78	0.47
1:F:43:LYS:HD3	1:F:44:LYS:H	1.76	0.47
1:F:50:ILE:HD11	1:F:88:LYS:HD2	1.95	0.47
1:D:116:VAL:HG22	1:D:117:LYS:N	2.30	0.47
1:G:129:GLU:O	1:G:130:ASN:C	2.55	0.47
1:E:30:SER:HB2	1:F:15:PRO:HA	1.96	0.47
1:E:98:ALA:O	1:F:34:ASP:HB2	2.15	0.47
1:E:115:ARG:HH12	1:F:16:ALA:N	2.12	0.47
1:B:30:SER:O	1:B:117:LYS:HE3	2.14	0.47
1:E:73:GLY:C	1:E:75:LYS:H	2.23	0.47
1:B:54:VAL:O	1:B:54:VAL:HG12	2.14	0.47
1:G:54:VAL:HG12	1:G:54:VAL:O	2.15	0.47
1:A:60:TYR:O	1:A:64:HIS:HD2	1.97	0.47
1:B:115:ARG:HG3	1:B:115:ARG:HH11	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:50:ILE:HD11	1:E:88:LYS:HD2	1.97	0.47
1:F:94:CYS:HB3	2:F:142:HOH:O	2.15	0.46
1:G:91:ILE:HA	1:G:117:LYS:O	2.16	0.46
1:I:60:TYR:O	1:I:64:HIS:HD2	1.98	0.46
1:D:19:ASN:HB2	2:D:6:HOH:O	2.15	0.46
1:F:73:GLY:C	1:F:75:LYS:H	2.23	0.46
1:B:66:PRO:HG3	1:B:133:GLU:HB2	1.98	0.46
1:G:73:GLY:C	1:G:75:LYS:H	2.23	0.46
1:I:115:ARG:HG3	1:I:115:ARG:HH11	1.80	0.46
1:A:53:ASP:OD1	1:A:55:ARG:HD3	2.16	0.46
1:A:16:ALA:O	1:C:115:ARG:NH1	2.49	0.46
1:C:46:TYR:CZ	1:C:48:GLY:HA3	2.51	0.46
1:D:49:ILE:HD12	1:D:49:ILE:O	2.16	0.45
1:E:115:ARG:NH2	1:F:15:PRO:N	2.64	0.45
1:F:54:VAL:O	1:F:54:VAL:HG12	2.15	0.45
1:G:60:TYR:O	1:G:64:HIS:HD2	1.98	0.45
1:B:129:GLU:CB	1:B:130:ASN:HD22	2.24	0.45
1:C:115:ARG:HG3	1:C:115:ARG:HH11	1.80	0.45
1:C:121:GLY:HA3	1:C:125:TYR:CD2	2.52	0.45
1:C:91:ILE:HA	1:C:117:LYS:O	2.16	0.45
1:A:94:CYS:SG	1:A:95:TRP:N	2.81	0.45
1:B:73:GLY:C	1:B:75:LYS:H	2.25	0.45
1:D:115:ARG:HH11	1:D:115:ARG:HG3	1.81	0.45
1:D:53:ASP:OD1	1:D:55:ARG:HD3	2.17	0.45
1:D:73:GLY:C	1:D:75:LYS:H	2.25	0.45
1:E:17:ASN:H	1:E:18:PRO:CD	2.29	0.45
1:I:78:GLU:HG2	1:I:112:LEU:HD22	1.99	0.45
1:E:36:ALA:N	2:E:143:HOH:O	2.45	0.45
1:I:73:GLY:C	1:I:75:LYS:H	2.24	0.45
1:C:73:GLY:C	1:C:75:LYS:H	2.24	0.44
1:F:115:ARG:HG3	1:F:115:ARG:HH11	1.82	0.44
1:A:17:ASN:O	1:A:18:PRO:C	2.59	0.44
1:A:115:ARG:HG3	1:A:115:ARG:HH11	1.81	0.44
1:C:63:CYS:HB2	1:C:132:LEU:CD1	2.47	0.44
1:A:91:ILE:HA	1:A:117:LYS:O	2.18	0.44
1:C:96:GLY:HA2	1:C:121:GLY:O	2.16	0.44
1:D:50:ILE:HD11	1:D:88:LYS:HD2	1.99	0.44
1:F:78:GLU:HG2	1:F:112:LEU:HD22	2.00	0.44
1:D:121:GLY:HA3	1:D:125:TYR:CD2	2.53	0.44
1:G:48:GLY:HA2	1:G:88:LYS:HG2	2.00	0.44
1:H:54:VAL:O	1:H:54:VAL:HG12	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:115:ARG:HG3	1:G:115:ARG:HH11	1.82	0.44
1:B:96:GLY:HA2	1:B:121:GLY:O	2.18	0.44
1:A:50:ILE:HD11	1:A:88:LYS:HD2	1.98	0.44
1:H:73:GLY:C	1:H:75:LYS:H	2.26	0.44
1:A:96:GLY:HA2	1:A:121:GLY:O	2.17	0.44
1:H:91:ILE:HA	1:H:117:LYS:O	2.17	0.44
1:H:115:ARG:HG3	1:H:115:ARG:HH11	1.83	0.44
1:B:77:ASN:ND2	1:B:79:ASP:H	2.05	0.44
1:B:115:ARG:NH1	1:C:16:ALA:N	2.66	0.44
1:D:91:ILE:HA	1:D:117:LYS:O	2.18	0.44
1:E:28:LYS:O	1:E:32:GLU:HG2	2.17	0.44
1:E:78:GLU:HG2	1:E:112:LEU:HD22	1.99	0.43
1:F:96:GLY:HA2	1:F:121:GLY:O	2.18	0.43
1:G:38:LEU:HD12	1:G:126:TRP:CZ3	2.53	0.43
1:B:94:CYS:SG	1:B:95:TRP:N	2.81	0.43
1:F:43:LYS:HB2	1:F:43:LYS:HE2	1.69	0.43
1:A:28:LYS:O	1:A:32:GLU:HG2	2.17	0.43
1:B:46:TYR:CE2	1:B:48:GLY:HA3	2.54	0.43
1:D:43:LYS:NZ	1:D:43:LYS:CB	2.81	0.43
1:B:28:LYS:O	1:B:32:GLU:HG2	2.18	0.43
1:B:91:ILE:HA	1:B:117:LYS:O	2.18	0.43
1:C:53:ASP:OD1	1:C:55:ARG:HD3	2.19	0.43
1:C:89:VAL:HG13	1:C:115:ARG:O	2.18	0.43
1:C:125:TYR:O	1:C:129:GLU:HG2	2.19	0.43
1:D:115:ARG:NH1	1:E:17:ASN:OD1	2.52	0.43
1:G:89:VAL:HG13	1:G:115:ARG:O	2.19	0.43
1:C:50:ILE:HD11	1:C:88:LYS:HD2	2.00	0.43
1:I:77:ASN:ND2	1:I:79:ASP:H	2.07	0.43
1:D:98:ALA:O	1:E:34:ASP:HB2	2.19	0.42
1:A:73:GLY:C	1:A:75:LYS:H	2.26	0.42
1:E:62:GLU:O	1:E:63:CYS:HB3	2.19	0.42
1:B:50:ILE:HD11	1:B:88:LYS:HD2	2.01	0.42
1:F:91:ILE:HA	1:F:117:LYS:O	2.19	0.42
1:D:23:ARG:HH21	1:I:20:GLU:HA	1.84	0.42
1:A:97:PRO:HB2	1:B:125:TYR:CE2	2.55	0.42
1:D:43:LYS:C	1:D:45:GLY:H	2.28	0.42
1:E:30:SER:HB2	1:F:15:PRO:CA	2.49	0.42
1:E:115:ARG:HG3	1:E:115:ARG:HH11	1.85	0.42
1:A:125:TYR:O	1:A:129:GLU:HG2	2.18	0.42
1:B:115:ARG:HH12	1:C:16:ALA:N	2.17	0.42
1:D:54:VAL:O	1:D:54:VAL:HG12	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:63:CYS:SG	1:F:132:LEU:HD22	2.59	0.42
1:D:77:ASN:ND2	1:D:79:ASP:H	2.05	0.42
1:D:115:ARG:CZ	1:E:17:ASN:OD1	2.68	0.42
1:E:91:ILE:HA	1:E:117:LYS:O	2.19	0.42
1:H:50:ILE:HD11	1:H:88:LYS:HD2	2.01	0.42
1:I:50:ILE:HD11	1:I:88:LYS:HD2	2.01	0.42
1:A:54:VAL:O	1:A:54:VAL:CG1	2.68	0.42
1:A:100:ASN:HD22	1:A:100:ASN:HA	1.62	0.42
1:B:121:GLY:HA3	1:B:125:TYR:CD2	2.55	0.42
1:C:19:ASN:HB3	1:C:23:ARG:NH1	2.35	0.42
1:I:91:ILE:HA	1:I:117:LYS:O	2.20	0.42
1:G:44:LYS:NZ	1:G:44:LYS:CB	2.82	0.41
1:G:127:ARG:O	1:G:130:ASN:HB2	2.20	0.41
1:B:78:GLU:HG2	1:B:112:LEU:HD22	2.02	0.41
1:D:46:TYR:HD2	1:D:49:ILE:HG13	1.85	0.41
1:D:125:TYR:O	1:D:129:GLU:HG2	2.20	0.41
1:G:36:ALA:O	1:G:40:ILE:HG12	2.20	0.41
1:I:61:LYS:HB2	1:I:61:LYS:HE3	1.87	0.41
1:D:85:SER:C	1:D:87:GLU:H	2.28	0.41
1:H:127:ARG:HG2	1:H:132:LEU:HD11	2.02	0.41
1:I:75:LYS:HE3	2:I:142:HOH:O	2.21	0.41
1:F:121:GLY:HA3	1:F:125:TYR:CD2	2.56	0.41
1:G:19:ASN:OD1	2:G:147:HOH:O	2.21	0.41
1:I:28:LYS:O	1:I:32:GLU:HG2	2.21	0.41
1:A:36:ALA:O	1:A:40:ILE:HG12	2.21	0.41
1:C:77:ASN:ND2	1:C:79:ASP:H	2.06	0.41
1:D:61:LYS:HE3	1:D:61:LYS:HB2	1.90	0.41
1:E:121:GLY:HA3	1:E:125:TYR:CD2	2.56	0.41
1:A:20:GLU:HG2	1:C:31:TYR:OH	2.21	0.41
1:C:28:LYS:O	1:C:32:GLU:HG2	2.21	0.41
1:C:79:ASP:CG	1:E:23:ARG:HG2	2.46	0.41
1:E:96:GLY:HA2	1:E:121:GLY:O	2.21	0.41
1:I:54:VAL:HG13	1:I:73:GLY:CA	2.51	0.41
1:H:121:GLY:HA3	1:H:125:TYR:CD2	2.56	0.41
1:I:80:THR:HG23	2:I:142:HOH:O	2.20	0.41
1:A:78:GLU:HG2	1:A:112:LEU:HD22	2.02	0.40
1:C:19:ASN:HB3	1:C:23:ARG:HH12	1.86	0.40
1:H:78:GLU:HG2	1:H:112:LEU:HD22	2.03	0.40
1:G:65:ILE:HA	1:G:66:PRO:HD3	1.99	0.40
1:I:53:ASP:OD2	1:I:55:ARG:NH1	2.53	0.40
1:A:33:THR:HA	2:A:11:HOH:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:94:CYS:SG	1:E:95:TRP:N	2.78	0.40
1:F:77:ASN:ND2	1:F:79:ASP:H	2.09	0.40
1:E:115:ARG:NH2	1:F:16:ALA:N	2.70	0.40
1:B:53:ASP:OD1	1:B:55:ARG:HD3	2.21	0.40
1:C:82:LYS:HZ1	1:E:23:ARG:NH2	2.19	0.40
1:D:100:ASN:HD22	1:D:100:ASN:HA	1.64	0.40
1:E:45:GLY:O	1:E:46:TYR:C	2.65	0.40
1:G:93:TYR:CD1	1:G:93:TYR:C	3.00	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:C:22:TYR:OH	1:F:83:ARG:NH1[2_455]	2.11	0.09

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	116/126 (92%)	108 (93%)	7 (6%)	1 (1%)	14	40
1	B	115/126 (91%)	109 (95%)	6 (5%)	0	100	100
1	C	116/126 (92%)	108 (93%)	7 (6%)	1 (1%)	14	40
1	D	112/126 (89%)	108 (96%)	4 (4%)	0	100	100
1	E	116/126 (92%)	109 (94%)	6 (5%)	1 (1%)	14	40
1	F	114/126 (90%)	106 (93%)	8 (7%)	0	100	100
1	G	112/126 (89%)	103 (92%)	9 (8%)	0	100	100
1	H	114/126 (90%)	105 (92%)	9 (8%)	0	100	100
1	I	109/126 (86%)	102 (94%)	7 (6%)	0	100	100
All	All	1024/1134 (90%)	958 (94%)	63 (6%)	3 (0%)	36	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	17	ASN
1	A	15	PRO
1	C	132	LEU

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	96/104 (92%)	92 (96%)	4 (4%)	26	59
1	B	97/104 (93%)	92 (95%)	5 (5%)	21	50
1	C	97/104 (93%)	91 (94%)	6 (6%)	16	44
1	D	95/104 (91%)	91 (96%)	4 (4%)	26	59
1	E	97/104 (93%)	89 (92%)	8 (8%)	10	31
1	F	97/104 (93%)	89 (92%)	8 (8%)	10	31
1	G	95/104 (91%)	91 (96%)	4 (4%)	26	59
1	H	96/104 (92%)	94 (98%)	2 (2%)	47	76
1	I	93/104 (89%)	90 (97%)	3 (3%)	34	66
All	All	863/936 (92%)	819 (95%)	44 (5%)	21	51

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	LEU
1	A	44	LYS
1	A	77	ASN
1	A	80	THR
1	B	38	LEU
1	B	47	GLU
1	B	77	ASN
1	B	80	THR
1	B	129	GLU
1	C	38	LEU

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Mol	Chain	Res	Type
1	C	44	LYS
1	C	47	GLU
1	C	77	ASN
1	C	80	THR
1	C	129	GLU
1	D	43	LYS
1	D	49	ILE
1	D	77	ASN
1	D	80	THR
1	E	38	LEU
1	E	43	LYS
1	E	44	LYS
1	E	77	ASN
1	E	80	THR
1	E	129	GLU
1	E	130	ASN
1	E	133	GLU
1	F	38	LEU
1	F	43	LYS
1	F	77	ASN
1	F	80	THR
1	F	129	GLU
1	F	130	ASN
1	F	132	LEU
1	F	133	GLU
1	G	44	LYS
1	G	47	GLU
1	G	77	ASN
1	G	80	THR
1	H	77	ASN
1	H	80	THR
1	I	77	ASN
1	I	80	THR
1	I	129	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	ASN
1	A	100	ASN
1	A	111	GLN
1	B	17	ASN

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Mol	Chain	Res	Type
1	B	77	ASN
1	B	100	ASN
1	B	111	GLN
1	B	130	ASN
1	C	17	ASN
1	C	77	ASN
1	C	100	ASN
1	C	111	GLN
1	D	77	ASN
1	D	100	ASN
1	D	111	GLN
1	E	19	ASN
1	E	77	ASN
1	E	100	ASN
1	E	111	GLN
1	F	77	ASN
1	F	100	ASN
1	F	111	GLN
1	F	134	HIS
1	G	77	ASN
1	G	100	ASN
1	G	111	GLN
1	H	17	ASN
1	H	77	ASN
1	H	100	ASN
1	H	111	GLN
1	I	17	ASN
1	I	77	ASN
1	I	100	ASN
1	I	111	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	117/126 (92%)	-1.63	0 100 100	12, 25, 46, 55	0
1	B	116/126 (92%)	-1.67	0 100 100	6, 25, 47, 52	0
1	C	117/126 (92%)	-1.66	0 100 100	11, 24, 44, 52	0
1	D	113/126 (89%)	-1.66	0 100 100	10, 27, 47, 54	0
1	E	117/126 (92%)	-1.63	0 100 100	5, 23, 46, 55	0
1	F	117/126 (92%)	-1.61	0 100 100	9, 26, 46, 52	0
1	G	113/126 (89%)	-1.43	0 100 100	22, 38, 57, 66	0
1	H	115/126 (91%)	-1.51	0 100 100	24, 37, 55, 69	0
1	I	112/126 (88%)	-1.47	0 100 100	25, 38, 55, 62	0
All	All	1037/1134 (91%)	-1.59	0 100 100	5, 31, 51, 69	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.