



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:38 am GMT

PDB ID : 1HRT
Title : THE STRUCTURE OF A COMPLEX OF BOVINE ALPHA-THROMBIN
AND RECOMBINANT HIRUDIN AT 2.8 ANGSTROMS RESOLUTION
Authors : Vitali, J.; Edwards, B.F.P.
Deposited on : 1993-02-25
Resolution : 2.80 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

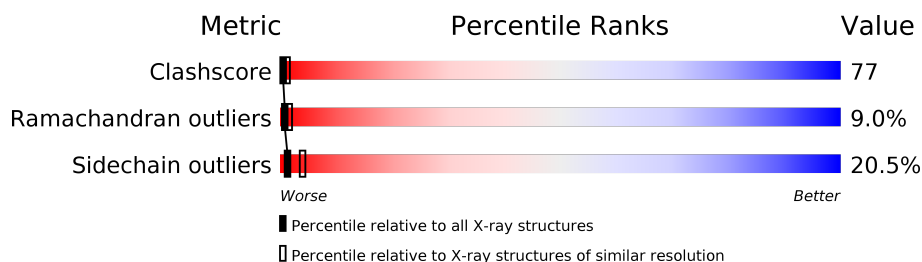
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	49	
2	H	259	
3	I	65	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2996 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 THROMBIN (SMALL SUBUNIT).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	36	Total	C	N	O	S	0	0	0
			290	181	48	60	1			

- Molecule 2 is a protein called THROMBIN (LARGE SUBUNIT).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	259	Total	C	N	O	S	0	0	0
			2094	1337	376	369	12			

- Molecule 3 is a protein called HIRUDIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	65	Total	C	N	O	S	0	0	0
			483	287	80	110	6			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	79	Total	O	0	0
			79	79		
4	I	25	Total	O	0	0
			25	25		
4	L	25	Total	O	0	0
			25	25		

4 Data and refinement statistics

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

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	59.11Å 102.62Å 143.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.80	Depositor
% Data completeness (in resolution range)	(Not available) (7.00-2.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT	Depositor
R, R_{free}	0.155 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2996	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	L	1.16	1/294 (0.3%)	1.93	5/390 (1.3%)
2	H	1.14	2/2148 (0.1%)	2.12	73/2905 (2.5%)
3	I	1.13	0/489	2.05	10/659 (1.5%)
All	All	1.14	3/2931 (0.1%)	2.09	88/3954 (2.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	L	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	171	SER	CB-OG	-5.88	1.34	1.42
2	H	42	CYS	CB-SG	-5.84	1.72	1.81
1	L	14(M)	GLY	N-CA	-5.41	1.38	1.46

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	110	ARG	CD-NE-CZ	18.33	149.26	123.60
2	H	144	ARG	CD-NE-CZ	18.11	148.96	123.60
2	H	187	ARG	NE-CZ-NH1	14.19	127.39	120.30
2	H	113	GLU	CA-CB-CG	13.56	143.23	113.40
2	H	137	ARG	CD-NE-CZ	12.35	140.89	123.60
2	H	187	ARG	CD-NE-CZ	11.94	140.32	123.60
2	H	130	LEU	CA-CB-CG	10.53	139.51	115.30
3	I	62	GLU	C-N-CA	10.41	147.72	121.70
2	H	110	ARG	NE-CZ-NH1	10.38	125.49	120.30
2	H	144	ARG	NE-CZ-NH2	10.37	125.48	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	21	ASP	CB-CG-OD1	9.54	126.89	118.30
2	H	67	ARG	NE-CZ-NH2	9.16	124.88	120.30
3	I	3	TYR	CA-CB-CG	9.11	130.70	113.40
2	H	123	LEU	CB-CA-C	8.73	126.79	110.20
2	H	146	GLU	CA-CB-CG	8.48	132.06	113.40
2	H	217	GLU	CA-CB-CG	8.26	131.56	113.40
3	I	33	ASP	CB-CG-OD1	-8.19	110.93	118.30
2	H	204(B)	ASN	CB-CA-C	7.99	126.38	110.40
1	L	14(M)	GLY	N-CA-C	7.97	133.01	113.10
2	H	182	CYS	CA-CB-SG	7.91	128.24	114.00
2	H	77(A)	ARG	NE-CZ-NH1	7.87	124.23	120.30
2	H	62	ASP	CB-CG-OD1	7.85	125.36	118.30
2	H	145	ARG	NE-CZ-NH2	7.75	124.18	120.30
2	H	40	LEU	CA-CB-CG	7.68	132.97	115.30
2	H	194	ASP	CB-CG-OD2	7.63	125.17	118.30
2	H	199	PHE	N-CA-CB	7.52	124.14	110.60
2	H	149(B)	SER	N-CA-C	7.16	130.32	111.00
2	H	194	ASP	CB-CG-OD1	-7.09	111.92	118.30
2	H	149(D)	ALA	N-CA-C	-7.02	92.04	111.00
2	H	206	ARG	NE-CZ-NH2	-6.99	116.80	120.30
2	H	244	ARG	C-N-CA	6.92	139.01	121.70
2	H	129(B)	LYS	CA-CB-CG	6.91	128.59	113.40
2	H	185	LYS	CB-CA-C	6.85	124.09	110.40
2	H	97(A)	GLU	CA-CB-CG	6.82	128.40	113.40
2	H	204(B)	ASN	CA-C-N	6.71	131.96	117.20
2	H	129(B)	LYS	CB-CG-CD	6.71	129.04	111.60
2	H	73	ARG	NE-CZ-NH1	-6.53	117.04	120.30
2	H	204(B)	ASN	CA-C-O	-6.47	106.51	120.10
2	H	101	ARG	CD-NE-CZ	6.41	132.57	123.60
2	H	93	ARG	NE-CZ-NH2	6.40	123.50	120.30
2	H	55	ALA	N-CA-CB	6.38	119.04	110.10
2	H	149(B)	SER	CA-C-O	6.36	133.45	120.10
2	H	149(D)	ALA	N-CA-CB	6.31	118.93	110.10
2	H	233	ARG	NE-CZ-NH2	6.28	123.44	120.30
2	H	187	ARG	N-CA-CB	-6.19	99.45	110.60
2	H	189	ASP	CB-CG-OD2	6.18	123.86	118.30
2	H	37	PRO	N-CA-C	6.18	128.16	112.10
2	H	33	LEU	CA-CB-CG	6.15	129.45	115.30
2	H	162	LEU	CB-CA-C	6.03	121.66	110.20
2	H	173	ARG	NE-CZ-NH2	6.02	123.31	120.30
3	I	62	GLU	N-CA-CB	5.99	121.39	110.60
3	I	15	LEU	CA-CB-CG	5.97	129.02	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	116	ASP	CB-CG-OD2	5.90	123.61	118.30
2	H	145	ARG	CA-CB-CG	5.85	126.26	113.40
2	H	60(A)	TYR	CB-CG-CD1	-5.83	117.50	121.00
2	H	60(H)	PHE	CA-CB-CG	5.78	127.77	113.90
1	L	3	LEU	CA-CB-CG	5.77	128.57	115.30
2	H	21	ASP	CA-CB-CG	5.74	126.03	113.40
2	H	76	TYR	N-CA-CB	5.60	120.68	110.60
1	L	1(C)	GLU	CG-CD-OE2	5.59	129.48	118.30
2	H	244	ARG	NE-CZ-NH1	5.58	123.09	120.30
3	I	43	GLU	CA-CB-CG	5.58	125.67	113.40
2	H	116	ASP	CB-CG-OD1	-5.51	113.34	118.30
2	H	146	GLU	OE1-CD-OE2	-5.47	116.73	123.30
2	H	206	ARG	CB-CA-C	-5.45	99.50	110.40
2	H	35	ARG	CD-NE-CZ	5.42	131.19	123.60
1	L	14(H)	GLU	CA-CB-CG	5.39	125.26	113.40
2	H	46	LEU	CB-CA-C	5.35	120.37	110.20
2	H	158	VAL	O-C-N	5.32	131.21	122.70
3	I	6	CYS	O-C-N	5.31	131.20	122.70
2	H	141	TRP	CA-CB-CG	5.30	123.77	113.70
2	H	144	ARG	NE-CZ-NH1	-5.29	117.65	120.30
2	H	241	VAL	CA-CB-CG2	5.27	118.81	110.90
2	H	189	ASP	CB-CG-OD1	-5.23	113.60	118.30
2	H	129(B)	LYS	N-CA-CB	5.22	120.00	110.60
3	I	63	TYR	N-CA-CB	5.22	119.99	110.60
2	H	172	THR	N-CA-CB	-5.20	100.42	110.30
1	L	14(L)	GLU	C-N-CA	5.17	133.16	122.30
2	H	65	LEU	CA-CB-CG	5.16	127.16	115.30
3	I	22	CYS	N-CA-C	-5.14	97.13	111.00
2	H	187	ARG	CA-CB-CG	5.12	124.67	113.40
2	H	77(A)	ARG	CD-NE-CZ	5.11	130.75	123.60
2	H	191	CYS	CA-CB-SG	5.09	123.16	114.00
2	H	149(D)	ALA	O-C-N	5.08	130.84	122.70
2	H	94	TYR	C-N-CA	5.05	134.34	121.70
2	H	145	ARG	CD-NE-CZ	5.05	130.67	123.60
2	H	111	PRO	N-CD-CG	-5.04	95.64	103.20
3	I	18	GLY	N-CA-C	-5.03	100.53	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	4	ARG	Sidechain

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	L	290	0	276	61	0
2	H	2094	0	2096	344	0
3	I	483	0	430	84	0
4	H	79	0	0	42	0
4	I	25	0	0	3	0
4	L	25	0	0	10	0
All	All	2996	0	2802	434	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 77.

All (434) 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:H:143:ASN:HA	2:H:150:VAL:O	1.53	1.09
3:I:49:GLN:OE1	3:I:49:GLN:HA	1.48	1.09
1:L:1(G):PHE:HB2	2:H:51:TRP:NE1	1.73	1.03
2:H:147:THR:HA	2:H:149(B):SER:OG	1.58	1.03
2:H:204:PRO:HD2	2:H:204(A):TYR:HD1	1.22	1.02
1:L:1(G):PHE:HB2	2:H:51:TRP:HE1	1.21	0.97
2:H:149(C):VAL:HG13	4:H:527:HOH:O	1.67	0.93
2:H:19:GLY:HA2	2:H:158:VAL:HG13	1.52	0.92
2:H:174:ILE:HD11	3:I:21:VAL:CG1	2.00	0.92
1:L:14(D):LYS:H	1:L:14(D):LYS:HE2	1.36	0.90
2:H:60(I):THR:HG22	2:H:62:ASP:H	1.38	0.89
3:I:60:PRO:HB2	3:I:63:TYR:CE2	2.08	0.88
2:H:137:ARG:HD2	2:H:157:VAL:CG2	2.04	0.86
2:H:204:PRO:HD2	2:H:204(A):TYR:CD1	2.11	0.86
2:H:186:PRO:HG3	2:H:223:GLY:HA2	1.59	0.85
1:L:9:LYS:HE3	4:L:522:HOH:O	1.77	0.84
2:H:31:VAL:HG21	4:H:565:HOH:O	1.76	0.84
2:H:136:GLY:HA3	2:H:199:PHE:CZ	2.14	0.83
3:I:48:PRO:O	3:I:49:GLN:HB2	1.79	0.83
2:H:127:GLN:HE21	2:H:127:GLN:HA	1.44	0.82
2:H:31:VAL:HG13	2:H:66:VAL:HG13	1.61	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:244:ARG:HA	2:H:244:ARG:NE	1.93	0.82
2:H:98:ASN:HA	4:H:487:HOH:O	1.77	0.82
2:H:236:LYS:HD2	2:H:239:GLN:OE1	1.79	0.81
3:I:47:LYS:HD2	3:I:48:PRO:HD2	1.61	0.81
2:H:196:GLY:HA2	2:H:212:ILE:HG22	1.63	0.81
2:H:174:ILE:HD11	3:I:21:VAL:HG12	1.61	0.80
2:H:105:LEU:HD13	2:H:242:ILE:HD11	1.61	0.80
2:H:210:MET:O	2:H:231:VAL:HG23	1.80	0.80
2:H:173:ARG:NH1	4:H:490:HOH:O	2.11	0.79
2:H:165:ARG:NH2	2:H:178:ASP:HA	1.97	0.79
2:H:32:MET:CE	2:H:70:LYS:HD3	2.13	0.79
2:H:122:CYS:SG	2:H:206:ARG:HG2	2.22	0.79
2:H:46:LEU:O	2:H:120:PRO:HA	1.83	0.79
2:H:16:ILE:HG23	2:H:158:VAL:HG22	1.64	0.79
2:H:41:LEU:HD21	2:H:60(H):PHE:CE2	2.17	0.78
2:H:16:ILE:CG2	2:H:158:VAL:HG22	2.14	0.78
2:H:173:ARG:HB2	4:I:554:HOH:O	1.83	0.78
2:H:149(C):VAL:O	4:H:581:HOH:O	2.01	0.77
2:H:164:GLU:HG2	2:H:167:VAL:HG13	1.66	0.77
1:L:14(D):LYS:H	1:L:14(D):LYS:CE	1.97	0.77
2:H:174:ILE:HD11	3:I:21:VAL:HG11	1.67	0.76
2:H:36:LYS:HG2	2:H:36:LYS:O	1.85	0.76
3:I:12:ASN:HD21	3:I:24:GLN:HA	1.48	0.76
2:H:50:ARG:NH1	4:H:558:HOH:O	2.05	0.75
2:H:16:ILE:N	2:H:194:ASP:OD2	2.19	0.75
2:H:46:LEU:HD23	2:H:51:TRP:O	1.85	0.75
1:L:1(G):PHE:CB	2:H:51:TRP:HE1	1.97	0.75
2:H:131:HIS:HA	4:H:511:HOH:O	1.87	0.74
3:I:27:LYS:HE3	3:I:42:GLY:HA3	1.69	0.74
2:H:149(B):SER:O	4:H:523:HOH:O	2.05	0.74
2:H:148:TRP:CD1	3:I:4:THR:HG21	2.23	0.74
2:H:57:HIS:HE1	2:H:214:SER:O	1.70	0.74
2:H:100:ASP:CG	2:H:177:THR:HG21	2.09	0.73
1:L:14(H):GLU:O	1:L:14(L):GLU:HG2	1.89	0.72
2:H:143:ASN:HD22	2:H:192:GLU:HG3	1.53	0.72
1:L:8:GLU:O	1:L:11:GLN:N	2.15	0.72
3:I:11:GLN:HA	3:I:45:THR:O	1.89	0.72
2:H:244:ARG:CA	2:H:244:ARG:NE	2.53	0.71
1:L:6:LEU:HA	1:L:10:LYS:HD3	1.72	0.71
2:H:146:GLU:HA	4:H:507:HOH:O	1.89	0.71
2:H:146:GLU:CA	4:H:507:HOH:O	2.37	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:36:LYS:HD2	3:I:64:LEU:HA	1.73	0.71
1:L:15:ARG:NH1	4:L:583:HOH:O	2.24	0.70
2:H:31:VAL:CG1	2:H:66:VAL:HG13	2.21	0.70
3:I:49:GLN:OE1	3:I:49:GLN:CA	2.34	0.70
2:H:146:GLU:HB3	2:H:148:TRP:HE1	1.55	0.70
2:H:137:ARG:HD2	2:H:157:VAL:HG21	1.74	0.70
2:H:30:GLN:HG2	2:H:155:LEU:HD11	1.73	0.70
1:L:6:LEU:O	1:L:10:LYS:HB2	1.91	0.70
2:H:73:ARG:NE	2:H:152:PRO:O	2.24	0.70
2:H:48:SER:HB2	2:H:51:TRP:HD1	1.57	0.69
2:H:29:TRP:CH2	2:H:207:TRP:HB2	2.27	0.69
2:H:126:LYS:HG3	2:H:232:PHE:HZ	1.57	0.69
2:H:19:GLY:CA	2:H:158:VAL:HG13	2.22	0.69
2:H:31:VAL:HG11	4:H:565:HOH:O	1.92	0.69
1:L:1(G):PHE:HB2	2:H:51:TRP:CE2	2.28	0.69
3:I:60:PRO:HB2	3:I:63:TYR:HE2	1.57	0.69
2:H:192:GLU:OE2	4:H:508:HOH:O	2.09	0.69
2:H:47:ILE:O	2:H:120:PRO:HB3	1.93	0.68
1:L:3:LEU:HB3	4:L:522:HOH:O	1.94	0.68
2:H:146:GLU:HB3	2:H:148:TRP:NE1	2.09	0.68
1:L:1(H):THR:OG1	2:H:245:LEU:HD11	1.94	0.68
2:H:136:GLY:N	2:H:160:LEU:O	2.21	0.68
2:H:60(I):THR:HG22	2:H:61:VAL:N	2.09	0.68
2:H:244:ARG:HE	2:H:244:ARG:CA	2.05	0.68
2:H:138:VAL:HG22	2:H:199:PHE:HD2	1.59	0.67
2:H:186:PRO:CG	2:H:223:GLY:HA2	2.24	0.67
2:H:69:GLY:N	4:H:512:HOH:O	2.21	0.67
3:I:59:ILE:HG12	3:I:60:PRO:HD2	1.76	0.67
2:H:245:LEU:HD12	2:H:246:GLY:N	2.10	0.66
2:H:197:GLY:O	2:H:213:VAL:HG23	1.96	0.66
2:H:204:PRO:O	4:H:586:HOH:O	2.13	0.66
1:L:3:LEU:CB	4:L:522:HOH:O	2.44	0.66
2:H:65:LEU:HD13	2:H:82:ILE:HG21	1.77	0.66
2:H:186(D):LYS:HD2	4:H:533:HOH:O	1.95	0.65
3:I:53:ASP:CB	3:I:56:PHE:HE2	2.08	0.65
3:I:47:LYS:HB3	3:I:48:PRO:HD2	1.77	0.65
2:H:192:GLU:HG3	4:H:508:HOH:O	1.96	0.65
2:H:240:LYS:HD2	2:H:244:ARG:NH1	2.12	0.65
2:H:62:ASP:OD1	4:H:541:HOH:O	2.14	0.65
3:I:49:GLN:HG3	3:I:50:SER:H	1.60	0.65
2:H:46:LEU:C	2:H:46:LEU:HD22	2.17	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:14(G):PHE:HA	1:L:14(J):TYR:CD2	2.31	0.65
2:H:141:TRP:HH2	4:H:578:HOH:O	1.79	0.64
2:H:49:ASP:HB2	4:H:503:HOH:O	1.97	0.64
1:L:1(G):PHE:CB	2:H:51:TRP:NE1	2.56	0.64
2:H:147:THR:OG1	2:H:149(B):SER:HB2	1.97	0.64
2:H:163:VAL:HG13	2:H:167:VAL:CG2	2.27	0.64
2:H:163:VAL:HG13	2:H:167:VAL:HG22	1.78	0.64
2:H:129(A):ALA:O	4:H:572:HOH:O	2.14	0.64
2:H:211:GLY:HA2	2:H:231:VAL:HG23	1.80	0.64
2:H:60(A):TYR:HB3	2:H:60(F):LYS:HD3	1.79	0.64
2:H:66:VAL:CG2	2:H:85:LEU:HD11	2.27	0.64
2:H:127:GLN:NE2	2:H:127:GLN:HA	2.11	0.63
2:H:137:ARG:HD3	2:H:159:ASN:ND2	2.13	0.63
3:I:29:ILE:HD12	3:I:40:VAL:CG1	2.28	0.63
2:H:32:MET:HE3	2:H:70:LYS:HD3	1.80	0.63
2:H:125:ASP:O	2:H:127:GLN:N	2.32	0.62
1:L:4:ARG:NH1	1:L:8:GLU:OE1	2.32	0.62
2:H:109:LYS:HB3	2:H:110:ARG:NH1	2.14	0.62
2:H:148:TRP:NE1	3:I:4:THR:CG2	2.62	0.62
2:H:135:LYS:HA	2:H:160:LEU:O	1.98	0.62
2:H:236:LYS:HA	2:H:239:GLN:OE1	1.99	0.62
3:I:13:LEU:HD12	3:I:46:PRO:HB3	1.81	0.62
2:H:70:LYS:HB3	4:H:578:HOH:O	1.98	0.62
1:L:14(E):GLU:OE1	4:L:529:HOH:O	2.16	0.62
3:I:52:ASN:O	3:I:53:ASP:HB2	2.00	0.62
2:H:148:TRP:CZ3	3:I:5:ASP:HB2	2.34	0.62
2:H:40:LEU:HD13	2:H:40:LEU:O	2.00	0.61
1:L:1(E):ALA:HA	2:H:48:SER:CA	2.30	0.61
2:H:73:ARG:O	3:I:56:PHE:HE1	1.82	0.61
2:H:82:ILE:HD11	3:I:59:ILE:HD11	1.82	0.61
2:H:82:ILE:CD1	3:I:59:ILE:CD1	2.78	0.61
3:I:38:GLN:HB2	4:I:486:HOH:O	2.01	0.60
2:H:110:ARG:N	2:H:110:ARG:HD2	2.16	0.60
2:H:59:LEU:N	2:H:59:LEU:HD13	2.16	0.60
3:I:22:CYS:SG	3:I:26:ASN:HB3	2.42	0.60
3:I:49:GLN:HG3	3:I:50:SER:N	2.17	0.60
1:L:5:PRO:HA	1:L:9:LYS:HG3	1.83	0.60
2:H:243:ASP:O	2:H:245:LEU:HG	2.01	0.60
2:H:126:LYS:HG3	2:H:232:PHE:CZ	2.37	0.60
3:I:12:ASN:HB3	3:I:46:PRO:HA	1.84	0.60
1:L:14:ASP:OD2	2:H:26:LEU:HG	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:216:GLY:O	3:I:3:TYR:HD1	1.84	0.59
3:I:53:ASP:HB3	3:I:56:PHE:HE2	1.67	0.59
2:H:87:LYS:HA	2:H:87:LYS:HZ2	1.67	0.59
2:H:148:TRP:CE3	3:I:5:ASP:HB2	2.38	0.59
2:H:122:CYS:SG	2:H:206:ARG:CG	2.90	0.59
2:H:20:GLN:NE2	2:H:157:VAL:O	2.35	0.59
2:H:147:THR:CB	2:H:149(B):SER:HB2	2.33	0.59
2:H:147:THR:C	2:H:149:THR:H	2.05	0.59
1:L:4:ARG:NH1	2:H:207:TRP:HE1	2.01	0.59
3:I:60:PRO:CB	3:I:63:TYR:HE2	2.16	0.58
2:H:190:ALA:O	2:H:220:CYS:HB3	2.02	0.58
2:H:221(A):ARG:HB3	2:H:224:LYS:HG3	1.84	0.58
2:H:41:LEU:HD21	2:H:60(H):PHE:CZ	2.38	0.58
2:H:66:VAL:HG23	2:H:85:LEU:HD11	1.85	0.58
2:H:233:ARG:NE	4:H:549:HOH:O	2.35	0.58
2:H:52:VAL:CG2	4:H:565:HOH:O	2.51	0.58
2:H:243:ASP:OD2	2:H:244:ARG:NH2	2.37	0.58
2:H:151:GLN:HG3	4:H:579:HOH:O	2.04	0.57
1:L:14(D):LYS:HZ3	1:L:14(D):LYS:N	2.01	0.57
2:H:137:ARG:O	2:H:199:PHE:HA	2.03	0.57
2:H:105:LEU:CD1	2:H:242:ILE:HD11	2.32	0.57
2:H:243:ASP:O	2:H:244:ARG:C	2.42	0.57
3:I:13:LEU:HD11	3:I:24:GLN:HE21	1.69	0.57
1:L:4:ARG:HH22	1:L:14:ASP:CB	2.17	0.57
2:H:60(A):TYR:CB	2:H:60(F):LYS:HD3	2.34	0.57
2:H:82:ILE:HD13	3:I:59:ILE:HD13	1.86	0.57
3:I:60:PRO:HG2	3:I:63:TYR:HE2	1.69	0.57
2:H:240:LYS:HD3	2:H:244:ARG:HH22	1.69	0.57
3:I:29:ILE:HD12	3:I:40:VAL:HG11	1.87	0.57
2:H:143:ASN:HD22	2:H:192:GLU:HB2	1.70	0.57
3:I:60:PRO:CG	3:I:63:TYR:HE2	2.18	0.57
2:H:30:GLN:HG2	2:H:155:LEU:CD1	2.34	0.56
2:H:52:VAL:HG22	4:H:565:HOH:O	2.05	0.56
3:I:13:LEU:HA	3:I:22:CYS:O	2.04	0.56
3:I:47:LYS:CD	3:I:48:PRO:HD2	2.31	0.56
2:H:177:THR:HG23	2:H:180:MET:SD	2.46	0.56
2:H:109:LYS:C	2:H:110:ARG:HD2	2.26	0.56
2:H:98:ASN:O	2:H:99:LEU:HB2	2.06	0.56
1:L:4:ARG:HH22	1:L:14:ASP:HB3	1.70	0.56
2:H:141:TRP:NE1	2:H:155:LEU:HB2	2.21	0.56
2:H:36:LYS:O	2:H:36(A):SER:OG	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:77(A):ARG:HD2	4:H:520:HOH:O	2.05	0.56
1:L:4:ARG:NH2	1:L:14:ASP:CB	2.68	0.56
2:H:51:TRP:CH2	2:H:107:LYS:HD3	2.41	0.56
1:L:1(E):ALA:HA	2:H:48:SER:HA	1.88	0.56
2:H:33:LEU:CD1	2:H:44:ALA:HB2	2.36	0.55
1:L:14(K):ILE:C	1:L:14(L):GLU:OE1	2.43	0.55
2:H:165:ARG:O	2:H:169:LYS:HD2	2.07	0.55
2:H:31:VAL:CG1	2:H:32:MET:N	2.68	0.55
2:H:82:ILE:CD1	3:I:59:ILE:HD11	2.36	0.55
2:H:102:ASP:OD2	2:H:214:SER:OG	2.18	0.55
2:H:109:LYS:HG3	2:H:109:LYS:O	2.06	0.55
2:H:204(B):ASN:OD1	2:H:206:ARG:HD3	2.07	0.55
2:H:148:TRP:CD1	3:I:4:THR:CG2	2.90	0.55
2:H:91:HIS:O	2:H:93:ARG:N	2.40	0.55
1:L:1(H):THR:HG22	4:L:488:HOH:O	2.06	0.55
2:H:176:ILE:HG13	2:H:180:MET:CE	2.37	0.55
2:H:221(A):ARG:NH1	3:I:15:LEU:HD12	2.21	0.55
3:I:47:LYS:HD2	3:I:48:PRO:CD	2.35	0.54
2:H:144:ARG:N	2:H:150:VAL:O	2.40	0.54
2:H:146:GLU:HB3	4:H:507:HOH:O	2.07	0.54
2:H:48:SER:HB2	2:H:51:TRP:CD1	2.39	0.54
1:L:8:GLU:O	1:L:10:LYS:N	2.40	0.54
2:H:49:ASP:O	2:H:112:ILE:HD13	2.07	0.54
2:H:137:ARG:CD	2:H:159:ASN:ND2	2.71	0.54
3:I:12:ASN:O	3:I:22:CYS:O	2.26	0.54
3:I:25:GLY:HA2	3:I:43:GLU:OE1	2.07	0.54
1:L:9:LYS:HB3	4:L:526:HOH:O	2.07	0.54
2:H:140:GLY:HA2	4:H:577:HOH:O	2.08	0.54
2:H:186(D):LYS:CD	4:H:533:HOH:O	2.54	0.54
2:H:200:VAL:HG23	2:H:207:TRP:CE3	2.43	0.54
2:H:105:LEU:HD12	2:H:241:VAL:HG22	1.89	0.53
2:H:163:VAL:HG12	2:H:168:CYS:SG	2.48	0.53
2:H:51:TRP:CE2	2:H:242:ILE:HD12	2.44	0.53
1:L:14(G):PHE:CZ	2:H:202:LYS:HD3	2.44	0.53
2:H:60(D):TRP:CZ3	3:I:49:GLN:HG2	2.44	0.53
2:H:91:HIS:CE1	2:H:101:ARG:HD3	2.43	0.53
2:H:244:ARG:HE	2:H:244:ARG:N	2.06	0.53
2:H:109:LYS:O	2:H:110:ARG:HB3	2.09	0.52
2:H:141:TRP:CE2	2:H:155:LEU:HB2	2.44	0.52
2:H:54:THR:CG2	2:H:59:LEU:HD21	2.39	0.52
2:H:82:ILE:CD1	3:I:59:ILE:HD13	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:52:VAL:HG13	4:H:565:HOH:O	2.09	0.52
2:H:184(A):TYR:OH	2:H:186(D):LYS:HE3	2.08	0.52
2:H:125:ASP:C	2:H:127:GLN:N	2.61	0.52
2:H:114:LEU:CD2	4:H:503:HOH:O	2.56	0.52
2:H:138:VAL:HG22	2:H:199:PHE:CD2	2.42	0.52
2:H:204(B):ASN:ND2	2:H:208:TYR:OH	2.42	0.52
2:H:31:VAL:HG12	2:H:32:MET:N	2.24	0.52
2:H:38:GLN:NE2	4:H:569:HOH:O	2.43	0.52
2:H:60(A):TYR:HD2	2:H:60(D):TRP:HB2	1.75	0.52
2:H:36:LYS:O	2:H:36:LYS:CG	2.56	0.52
3:I:59:ILE:HG12	3:I:60:PRO:CD	2.40	0.52
3:I:60:PRO:HB2	3:I:63:TYR:CD2	2.45	0.52
1:L:14(H):GLU:HG3	4:L:591:HOH:O	2.10	0.52
2:H:221(A):ARG:HB3	2:H:224:LYS:HB2	1.92	0.52
2:H:130:LEU:C	2:H:131:HIS:HD2	2.14	0.52
2:H:220:CYS:SG	3:I:2:VAL:HG11	2.50	0.51
2:H:224:LYS:HE2	4:H:501:HOH:O	2.11	0.51
2:H:244:ARG:HE	2:H:244:ARG:H	1.57	0.51
2:H:26:LEU:O	2:H:26:LEU:HD23	2.09	0.51
2:H:36:LYS:CD	3:I:64:LEU:HA	2.40	0.51
1:L:14(G):PHE:CE1	2:H:204:PRO:HB3	2.44	0.51
2:H:105:LEU:HD12	2:H:241:VAL:CG2	2.40	0.51
3:I:47:LYS:HB3	3:I:48:PRO:CD	2.41	0.51
1:L:8:GLU:O	1:L:9:LYS:C	2.48	0.51
2:H:109:LYS:CB	2:H:110:ARG:NH1	2.74	0.51
2:H:147:THR:N	4:H:507:HOH:O	2.43	0.51
2:H:235:LYS:HE2	2:H:239:GLN:CD	2.31	0.51
2:H:240:LYS:HD3	2:H:244:ARG:NH2	2.26	0.51
2:H:149:THR:O	2:H:149(A):THR:HG23	2.11	0.51
2:H:109:LYS:CG	2:H:110:ARG:NH1	2.74	0.50
2:H:52:VAL:CG1	4:H:565:HOH:O	2.59	0.50
2:H:176:ILE:HG13	2:H:180:MET:HE2	1.93	0.50
2:H:200:VAL:HG23	2:H:207:TRP:HE3	1.77	0.50
2:H:165:ARG:HH21	2:H:178:ASP:HA	1.76	0.50
2:H:245:LEU:C	2:H:245:LEU:HD12	2.32	0.50
2:H:125:ASP:C	2:H:127:GLN:H	2.15	0.50
2:H:212:ILE:O	2:H:228:TYR:HA	2.11	0.50
2:H:60(I):THR:CG2	2:H:61:VAL:N	2.75	0.50
3:I:53:ASP:HB3	3:I:56:PHE:CE2	2.45	0.50
1:L:14(B):THR:OG1	2:H:159:ASN:OD1	2.29	0.50
2:H:200:VAL:CG2	2:H:207:TRP:HE3	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:14(M):GLY:O	1:L:15:ARG:HG2	2.12	0.50
2:H:93:ARG:HB2	2:H:101:ARG:HD2	1.94	0.50
2:H:165:ARG:CB	2:H:166:PRO:CD	2.90	0.50
2:H:145:ARG:N	2:H:150:VAL:HG12	2.27	0.50
2:H:217:GLU:HB3	4:H:499:HOH:O	2.12	0.50
2:H:35:ARG:NH2	2:H:39:GLU:OE2	2.45	0.50
1:L:1:CYS:O	2:H:206:ARG:HG3	2.12	0.50
2:H:50:ARG:HH11	2:H:50:ARG:HG2	1.77	0.50
2:H:60(A):TYR:CZ	3:I:1:VAL:HG21	2.47	0.50
2:H:60(D):TRP:CH2	3:I:49:GLN:HG2	2.47	0.49
2:H:108:LEU:C	2:H:110:ARG:H	2.16	0.49
3:I:49:GLN:CG	3:I:50:SER:H	2.24	0.49
2:H:60(A):TYR:HB3	2:H:60(F):LYS:CD	2.42	0.49
2:H:164:GLU:HG2	2:H:167:VAL:CG1	2.39	0.49
2:H:56:ALA:O	2:H:59:LEU:HD22	2.13	0.49
2:H:143:ASN:HD22	2:H:192:GLU:CG	2.24	0.49
2:H:164:GLU:H	2:H:164:GLU:CD	2.16	0.49
2:H:17:VAL:HG23	2:H:191:CYS:HB2	1.95	0.49
1:L:4:ARG:HH11	2:H:207:TRP:HE1	1.60	0.49
2:H:127:GLN:O	2:H:129(B):LYS:HG2	2.13	0.49
2:H:148:TRP:NE1	3:I:4:THR:HG21	2.26	0.49
2:H:66:VAL:HG21	2:H:85:LEU:HD11	1.94	0.49
1:L:5:PRO:HA	1:L:9:LYS:CG	2.43	0.49
2:H:105:LEU:HD13	2:H:242:ILE:CD1	2.36	0.49
2:H:60(B):PRO:N	2:H:60(C):PRO:CD	2.76	0.49
2:H:137:ARG:HD2	2:H:157:VAL:HG23	1.93	0.48
1:L:1(G):PHE:HB3	2:H:50:ARG:HD2	1.94	0.48
2:H:108:LEU:HD13	2:H:112:ILE:CG2	2.43	0.48
2:H:143:ASN:CA	2:H:150:VAL:O	2.44	0.48
1:L:10:LYS:O	1:L:11:GLN:HB2	2.13	0.48
2:H:114:LEU:HG	4:H:503:HOH:O	2.13	0.48
2:H:91:HIS:HA	2:H:92:PRO:HD2	1.68	0.48
2:H:149(D):ALA:O	2:H:150:VAL:N	2.47	0.48
2:H:202:LYS:HE3	2:H:205:ASN:O	2.14	0.48
3:I:47:LYS:CB	3:I:48:PRO:HD2	2.41	0.48
2:H:58:CYS:C	2:H:59:LEU:HD13	2.34	0.48
2:H:147:THR:HA	2:H:149(B):SER:CB	2.44	0.48
2:H:165:ARG:N	2:H:166:PRO:HD2	2.28	0.48
2:H:20:GLN:O	2:H:20:GLN:HG2	2.14	0.48
2:H:210:MET:O	2:H:231:VAL:CG2	2.58	0.48
2:H:60(B):PRO:N	2:H:60(C):PRO:HD3	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:36:LYS:O	2:H:36(A):SER:CB	2.61	0.47
1:L:5:PRO:HA	1:L:9:LYS:CB	2.44	0.47
2:H:192:GLU:CG	4:H:508:HOH:O	2.60	0.47
1:L:14(D):LYS:N	1:L:14(D):LYS:CE	2.73	0.47
1:L:3:LEU:HG	4:L:522:HOH:O	2.15	0.47
2:H:60(B):PRO:HG2	2:H:96:TRP:CE2	2.49	0.47
3:I:48:PRO:O	3:I:49:GLN:CB	2.53	0.47
2:H:16:ILE:HG23	2:H:158:VAL:CG2	2.39	0.47
2:H:109:LYS:CB	2:H:110:ARG:HH11	2.26	0.47
2:H:39:GLU:HB2	3:I:52:ASN:O	2.15	0.47
2:H:87:LYS:HZ1	2:H:88:ILE:H	1.63	0.47
1:L:4:ARG:NH1	2:H:207:TRP:NE1	2.63	0.47
2:H:46:LEU:HD22	2:H:47:ILE:N	2.30	0.47
3:I:14:CYS:O	3:I:22:CYS:N	2.47	0.47
1:L:1(A):ASP:OD1	1:L:3:LEU:HD23	2.15	0.47
1:L:3:LEU:CG	4:L:522:HOH:O	2.62	0.47
2:H:146:GLU:CB	2:H:148:TRP:HE1	2.24	0.47
2:H:25:GLY:O	2:H:28:PRO:HD3	2.14	0.47
2:H:136:GLY:O	2:H:159:ASN:HA	2.15	0.47
2:H:204:PRO:HG2	2:H:204(A):TYR:CE1	2.50	0.47
2:H:234:LEU:O	2:H:235:LYS:C	2.52	0.47
1:L:1(F):GLY:C	1:L:1(D):GLY:H	2.18	0.47
2:H:215:TRP:HZ3	2:H:217:GLU:HB2	1.80	0.46
2:H:50:ARG:NH1	2:H:50:ARG:HG2	2.30	0.46
2:H:57:HIS:O	2:H:60(A):TYR:HB2	2.15	0.46
1:L:14(D):LYS:HZ2	1:L:14(D):LYS:HB3	1.80	0.46
2:H:200:VAL:CG2	2:H:207:TRP:CE3	2.99	0.46
2:H:29:TRP:CH2	2:H:207:TRP:CB	2.97	0.46
3:I:12:ASN:HB3	3:I:46:PRO:CA	2.45	0.46
2:H:33:LEU:HD23	2:H:64:LEU:HD13	1.96	0.46
2:H:51:TRP:CZ3	2:H:107:LYS:HB2	2.51	0.46
2:H:97(A):GLU:HB3	2:H:98:ASN:H	1.44	0.46
2:H:52:VAL:HB	2:H:106:LEU:HB2	1.98	0.46
2:H:50:ARG:HA	2:H:108:LEU:HB2	1.98	0.46
2:H:143:ASN:HB3	2:H:150:VAL:HB	1.96	0.46
2:H:137:ARG:CD	2:H:159:ASN:HD21	2.29	0.46
2:H:187:ARG:NH2	2:H:222:ASP:OD2	2.43	0.46
2:H:240:LYS:CD	2:H:244:ARG:NH1	2.79	0.46
1:L:1(A):ASP:O	2:H:119:HIS:NE2	2.49	0.46
2:H:149(A):THR:CG2	4:H:513:HOH:O	2.64	0.46
2:H:58:CYS:HB2	2:H:59:LEU:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:230:HIS:O	2:H:234:LEU:HD12	2.15	0.45
2:H:108:LEU:HD13	2:H:112:ILE:HG23	1.98	0.45
3:I:61:GLU:O	3:I:64:LEU:HD12	2.17	0.45
1:L:3:LEU:O	2:H:119:HIS:CD2	2.69	0.45
1:L:3:LEU:O	1:L:5:PRO:HD3	2.16	0.45
2:H:204:PRO:CD	2:H:204(A):TYR:CD1	2.93	0.45
3:I:53:ASP:CB	3:I:56:PHE:CE2	2.95	0.45
1:L:14(G):PHE:CE2	2:H:202:LYS:HD3	2.51	0.45
2:H:110:ARG:HG2	2:H:110:ARG:O	2.13	0.45
1:L:14(G):PHE:CD1	2:H:204:PRO:HB3	2.52	0.45
2:H:240:LYS:NZ	2:H:244:ARG:HH12	2.14	0.45
2:H:16:ILE:HD12	2:H:194:ASP:CG	2.37	0.44
2:H:200:VAL:HB	2:H:208:TYR:O	2.17	0.44
2:H:245:LEU:CD1	2:H:245:LEU:C	2.86	0.44
3:I:15:LEU:HD13	3:I:19:SER:HA	1.99	0.44
2:H:217:GLU:CB	4:H:499:HOH:O	2.65	0.44
3:I:49:GLN:CG	3:I:50:SER:N	2.79	0.44
2:H:16:ILE:HG21	2:H:158:VAL:HG22	1.96	0.44
2:H:148:TRP:NE1	3:I:4:THR:HG22	2.32	0.44
1:L:14(D):LYS:NZ	1:L:14(D):LYS:N	2.65	0.44
2:H:234:LEU:O	2:H:236:LYS:N	2.51	0.44
2:H:67:ARG:C	2:H:68:ILE:HG12	2.38	0.44
3:I:38:GLN:CB	4:I:489:HOH:O	2.65	0.44
2:H:51:TRP:HE3	2:H:106:LEU:O	2.00	0.44
2:H:95:ASN:OD1	2:H:97(A):GLU:HB2	2.18	0.44
1:L:5:PRO:HA	1:L:9:LYS:HB2	1.99	0.44
2:H:17:VAL:HG21	2:H:220:CYS:HB2	1.99	0.43
2:H:160:LEU:HA	2:H:161:PRO:HD3	1.82	0.43
2:H:142:GLY:HA3	2:H:192:GLU:O	2.18	0.43
2:H:235:LYS:HA	2:H:238:ILE:HD12	1.99	0.43
2:H:60(D):TRP:CZ3	3:I:49:GLN:CG	3.02	0.43
2:H:76:TYR:HE1	4:H:510:HOH:O	1.96	0.43
2:H:149(C):VAL:O	2:H:149(D):ALA:HB3	2.18	0.43
2:H:240:LYS:CD	2:H:244:ARG:HH12	2.32	0.43
2:H:129(A):ALA:O	2:H:131:HIS:NE2	2.45	0.43
1:L:14(B):THR:O	1:L:14(E):GLU:HB2	2.18	0.43
2:H:164:GLU:C	2:H:166:PRO:HD2	2.39	0.43
2:H:164:GLU:OE1	2:H:164:GLU:N	2.52	0.43
3:I:58:GLU:H	3:I:58:GLU:CD	2.22	0.43
2:H:164:GLU:O	2:H:167:VAL:N	2.52	0.42
2:H:143:ASN:ND2	2:H:192:GLU:HG3	2.28	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:137:ARG:HG3	2:H:138:VAL:N	2.34	0.42
2:H:184(A):TYR:C	2:H:185:LYS:O	2.56	0.42
3:I:47:LYS:CB	3:I:48:PRO:CD	2.96	0.42
2:H:139:THR:HA	2:H:156:GLN:O	2.19	0.42
2:H:201:MET:SD	2:H:210:MET:HG2	2.59	0.42
2:H:235:LYS:HE2	2:H:239:GLN:OE1	2.19	0.42
2:H:48:SER:HB3	2:H:50:ARG:H	1.84	0.42
3:I:45:THR:HA	3:I:46:PRO:HD3	1.88	0.42
3:I:32:SER:HB3	3:I:33:ASP:H	1.71	0.42
2:H:60(D):TRP:CH2	3:I:1:VAL:CG1	3.03	0.42
2:H:60(D):TRP:CH2	3:I:1:VAL:HG13	2.55	0.42
2:H:109:LYS:CG	2:H:110:ARG:HH11	2.33	0.42
2:H:191:CYS:O	2:H:194:ASP:HB2	2.20	0.41
2:H:197:GLY:H	2:H:213:VAL:HB	1.85	0.41
2:H:89:TYR:O	2:H:104:ALA:HB1	2.20	0.41
2:H:140:GLY:HA3	2:H:194:ASP:OD1	2.19	0.41
2:H:146:GLU:O	2:H:149:THR:OG1	2.38	0.41
2:H:221(A):ARG:NH1	3:I:15:LEU:CD1	2.83	0.41
1:L:14(I):SER:OG	2:H:134:PHE:HA	2.20	0.41
2:H:53:LEU:HD21	2:H:103:ILE:HG13	2.02	0.41
2:H:163:VAL:CG1	2:H:167:VAL:CG2	2.98	0.41
2:H:60(D):TRP:CD2	3:I:49:GLN:HB3	2.55	0.41
3:I:29:ILE:HD12	3:I:40:VAL:HG13	2.01	0.41
2:H:87:LYS:HD2	2:H:87:LYS:HA	1.84	0.41
2:H:177:THR:OG1	2:H:178:ASP:N	2.54	0.41
2:H:66:VAL:HG23	2:H:85:LEU:CD1	2.50	0.41
2:H:73:ARG:HD2	3:I:53:ASP:OD1	2.20	0.41
2:H:165:ARG:O	2:H:166:PRO:C	2.59	0.41
2:H:59:LEU:CD1	2:H:59:LEU:N	2.82	0.41
2:H:67:ARG:NH1	2:H:70:LYS:HZ2	2.18	0.41
2:H:73:ARG:O	3:I:56:PHE:CE1	2.70	0.41
2:H:196:GLY:HA2	2:H:212:ILE:CG2	2.43	0.41
2:H:234:LEU:C	2:H:236:LYS:N	2.69	0.41
2:H:36:LYS:HD2	3:I:63:TYR:O	2.21	0.41
3:I:60:PRO:CB	3:I:63:TYR:CE2	2.89	0.41
2:H:145:ARG:HB2	2:H:150:VAL:CG1	2.50	0.41
2:H:56:ALA:O	2:H:58:CYS:N	2.53	0.41
1:L:1(G):PHE:CD1	1:L:1(G):PHE:N	2.89	0.41
2:H:212:ILE:O	2:H:229:THR:N	2.50	0.40
2:H:57:HIS:CE1	2:H:214:SER:O	2.61	0.40
2:H:51:TRP:C	2:H:52:VAL:HG23	2.42	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:71:HIS:CD2	2:H:154:VAL:HG22	2.56	0.40
2:H:87:LYS:NZ	2:H:88:ILE:H	2.19	0.40
2:H:150:VAL:CG2	4:H:515:HOH:O	2.69	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	L	34/49 (69%)	22 (65%)	8 (24%)	4 (12%)	0	1
2	H	257/259 (99%)	200 (78%)	36 (14%)	21 (8%)	1	2
3	I	63/65 (97%)	49 (78%)	7 (11%)	7 (11%)	0	1
All	All	354/373 (95%)	271 (77%)	51 (14%)	32 (9%)	1	1

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	36(A)	SER
2	H	77(A)	ARG
2	H	110	ARG
2	H	150	VAL
2	H	245	LEU
3	I	49	GLN
3	I	50	SER
3	I	61	GLU
1	L	9	LYS
2	H	48	SER
2	H	57	HIS
2	H	97(A)	GLU
2	H	126	LYS
2	H	240	LYS

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Mol	Chain	Res	Type
2	H	243	ASP
2	H	244	ARG
3	I	53	ASP
1	L	7	PHE
1	L	8	GLU
1	L	14(M)	GLY
2	H	56	ALA
2	H	149(D)	ALA
2	H	149(E)	GLU
2	H	165	ARG
2	H	179	ASN
2	H	36	LYS
2	H	92	PRO
2	H	129	ALA
3	I	60	PRO
3	I	52	ASN
3	I	63	TYR
2	H	37	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	L	30/43 (70%)	22 (73%)	8 (27%)	0	1
2	H	226/226 (100%)	181 (80%)	45 (20%)	1	4
3	I	56/56 (100%)	45 (80%)	11 (20%)	1	4
All	All	312/325 (96%)	248 (80%)	64 (20%)	1	4

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	3	LEU
1	L	8	GLU
1	L	14	ASP
1	L	14(B)	THR

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Mol	Chain	Res	Type
1	L	14(D)	LYS
1	L	14(I)	SER
1	L	14(L)	GLU
1	L	15	ARG
2	H	20	GLN
2	H	21	ASP
2	H	27	SER
2	H	37	PRO
2	H	38	GLN
2	H	39	GLU
2	H	40	LEU
2	H	46	LEU
2	H	50	ARG
2	H	53	LEU
2	H	59	LEU
2	H	64	LEU
2	H	68	ILE
2	H	72	SER
2	H	75	ARG
2	H	81	LYS
2	H	102	ASP
2	H	110	ARG
2	H	112	ILE
2	H	127	GLN
2	H	129(B)	LYS
2	H	129(C)	LEU
2	H	139	THR
2	H	149	THR
2	H	149(A)	THR
2	H	151	GLN
2	H	155	LEU
2	H	158	VAL
2	H	167	VAL
2	H	172	THR
2	H	177	THR
2	H	182	CYS
2	H	192	GLU
2	H	203	SER
2	H	204(B)	ASN
2	H	210	MET
2	H	220	CYS
2	H	221(A)	ARG

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Mol	Chain	Res	Type
2	H	222	ASP
2	H	224	LYS
2	H	233	ARG
2	H	235	LYS
2	H	243	ASP
2	H	244	ARG
2	H	245	LEU
3	I	8	GLU
3	I	9	SER
3	I	19	SER
3	I	27	LYS
3	I	41	THR
3	I	43	GLU
3	I	45	THR
3	I	47	LYS
3	I	53	ASP
3	I	59	ILE
3	I	65	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	11	GLN
1	L	14(A)	GLN
2	H	20	GLN
2	H	57	HIS
2	H	60(G)	ASN
2	H	71	HIS
2	H	127	GLN
2	H	143	ASN
2	H	156	GLN
2	H	159	ASN
2	H	230	HIS
3	I	12	ASN
3	I	24	GLN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.