



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:52 PM UTC

PDB ID : 1EFX / pdb_00001efx
Title : STRUCTURE OF A COMPLEX BETWEEN THE HUMAN NATURAL KILLER CELL RECEPTOR KIR2DL2 AND A CLASS I MHC LIGAND HLA-CW3
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Deposited on : 2000-02-10
Resolution : 3.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

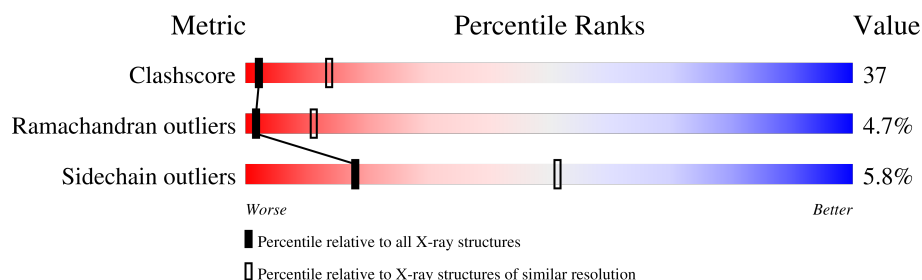
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	278	36% 55% 8% .
2	B	100	47% 45% 8%
3	C	9	44% 33% 22%
4	D	200	44% 46% 8% ..
4	E	200	42% 48% 8% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6411 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 HLA-CW3 (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	0	0
			2274	1420	413	435	6			

- Molecule 2 is a protein called BETA-2-MICROGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	SEE REMARK 999	UNP P61769

- Molecule 3 is a protein called PEPTIDE FROM IMPORTIN ALPHA-2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			61	40	9	12			

- Molecule 4 is a protein called NATURAL KILLER CELL RECEPTOR KIR2DL2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	197	Total	C	N	O	S	0	0	0
			1527	962	266	291	8			
4	E	197	Total	C	N	O	S	0	0	0
			1527	962	266	291	8			

- Molecule 5 is water.

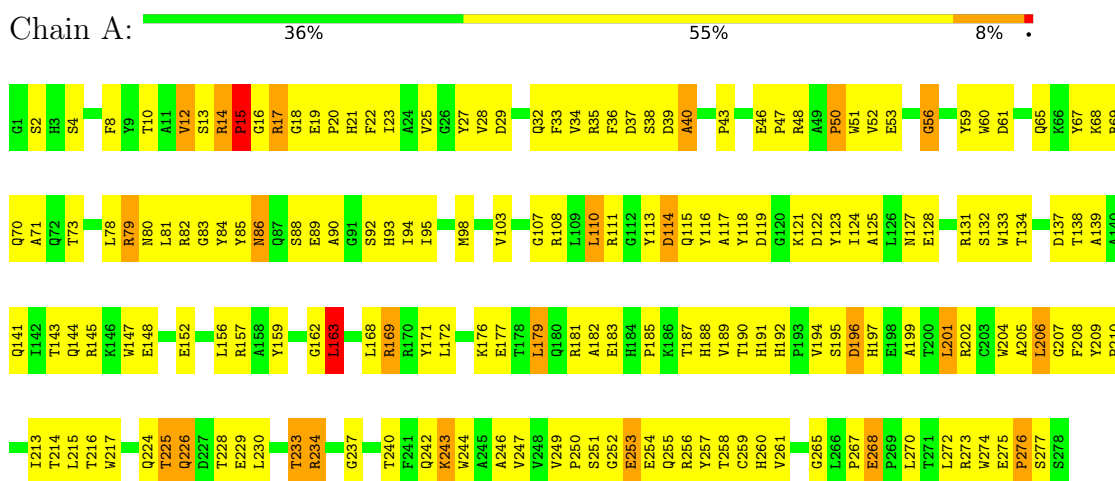
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	75	Total 75	O 75	0	0
5	B	28	Total 28	O 28	0	0
5	C	5	Total 5	O 5	0	0
5	D	46	Total 46	O 46	0	0
5	E	31	Total 31	O 31	0	0

3 Residue-property plots

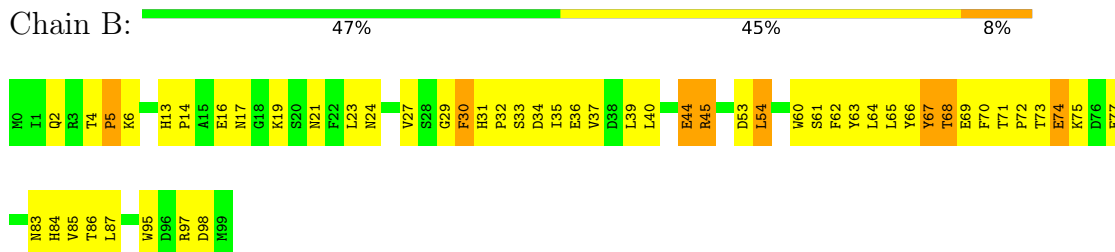
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

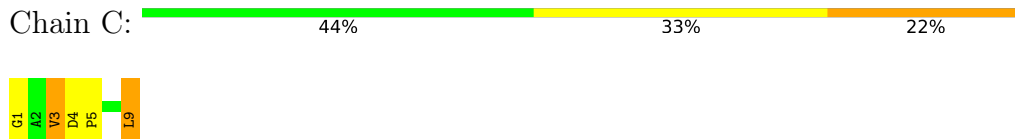
• Molecule 1: HLA-CW3 (HEAVY CHAIN)



• Molecule 2: BETA-2-MICROGLOBULIN

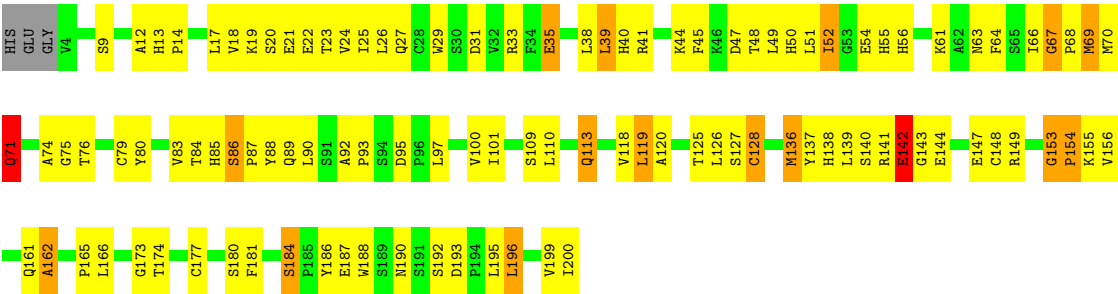


• Molecule 3: PEPTIDE FROM IMPORTIN ALPHA-2

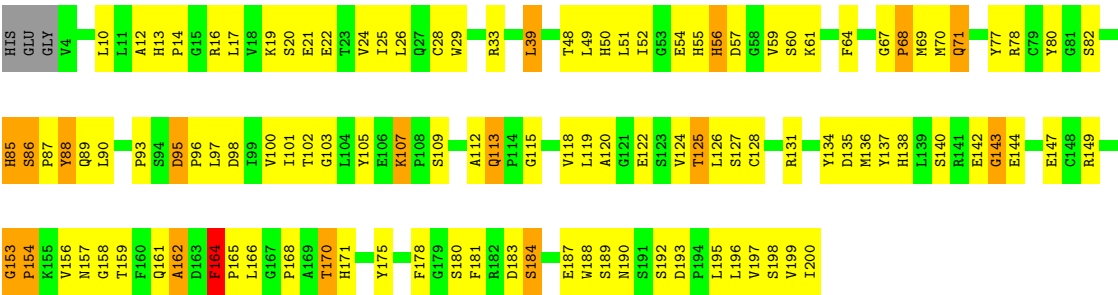


• Molecule 4: NATURAL KILLER CELL RECEPTOR KIR2DL2





● Molecule 4: NATURAL KILLER CELL RECEPTOR KIR2DL2



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.54Å 90.33Å 207.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 3.00	Depositor
% Data completeness (in resolution range)	88.5 (10.00-3.00)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.231 , 0.294	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6411	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.52	0/2340	1.00	14/3181 (0.4%)
2	B	0.51	0/860	1.05	6/1162 (0.5%)
3	C	0.69	0/61	0.82	0/82
4	D	0.60	0/1572	1.07	8/2133 (0.4%)
4	E	0.50	0/1572	1.07	13/2133 (0.6%)
All	All	0.54	0/6405	1.04	41/8691 (0.5%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	275	GLU	CA-C-N	10.58	133.06	119.84
1	A	275	GLU	C-N-CA	10.58	133.06	119.84
2	B	44	GLU	N-CA-C	7.88	121.05	109.07
1	A	56	GLY	CA-C-N	7.50	129.22	119.84
1	A	56	GLY	C-N-CA	7.50	129.22	119.84
4	D	184	SER	CA-C-N	6.81	126.95	119.87
4	D	184	SER	C-N-CA	6.81	126.95	119.87
4	E	95	ASP	CA-C-N	6.71	126.67	119.76
4	E	95	ASP	C-N-CA	6.71	126.67	119.76
2	B	6	LYS	N-CA-C	-6.29	101.62	110.50
1	A	14	ARG	CA-C-N	6.27	127.67	119.84
1	A	14	ARG	C-N-CA	6.27	127.67	119.84
4	D	153	GLY	CA-C-N	5.96	127.28	119.84
4	D	153	GLY	C-N-CA	5.96	127.28	119.84
4	E	184	SER	CA-C-N	5.84	125.54	119.82
4	E	184	SER	C-N-CA	5.84	125.54	119.82
1	A	226	GLN	N-CA-C	5.71	118.45	111.82
1	A	163	LEU	N-CA-C	5.71	118.44	111.82
1	A	89	GLU	N-CA-C	-5.70	104.68	111.69
4	D	31	ASP	N-CA-C	-5.59	105.96	112.89
1	A	110	LEU	N-CA-C	-5.50	104.20	111.96
4	E	159	THR	N-CA-C	5.50	117.42	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	69	MET	N-CA-C	5.49	118.44	110.42
4	E	131	ARG	N-CA-C	-5.42	105.45	111.36
2	B	5	PRO	N-CA-C	5.40	119.06	110.21
2	B	74	GLU	N-CA-C	-5.38	105.98	112.54
4	E	164	PHE	CB-CA-C	-5.36	105.62	110.44
1	A	107	GLY	N-CA-C	-5.33	100.56	113.18
2	B	35	ILE	N-CA-C	5.32	114.87	108.06
4	E	107	LYS	CA-C-N	-5.26	113.27	119.84
4	E	107	LYS	C-N-CA	-5.26	113.27	119.84
4	E	153	GLY	CA-C-N	5.22	126.36	119.84
4	E	153	GLY	C-N-CA	5.22	126.36	119.84
2	B	45	ARG	N-CA-C	5.21	117.64	110.35
1	A	169	ARG	N-CA-C	-5.20	105.31	110.97
4	E	69	MET	N-CA-C	5.16	117.58	110.35
4	D	95	ASP	CA-C-N	5.14	124.90	119.76
4	D	95	ASP	C-N-CA	5.14	124.90	119.76
4	E	158	GLY	N-CA-C	5.06	122.29	115.00
1	A	268	GLU	CA-C-N	5.05	125.05	119.90
1	A	268	GLU	C-N-CA	5.05	125.05	119.90

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	2274	0	2119	172	0
2	B	837	0	803	56	0
3	C	61	0	68	10	0
4	D	1527	0	1452	117	0
4	E	1527	0	1452	110	0
5	A	75	0	0	9	0
5	B	28	0	0	3	0
5	C	5	0	0	0	0
5	D	46	0	0	3	0
5	E	31	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6411	0	5894	447	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 37.

All (447) 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)
4:D:140:SER:HB2	4:D:147:GLU:OE1	1.58	1.02
4:D:52:ILE:HD12	4:D:52:ILE:H	1.26	0.96
4:D:86:SER:OG	4:D:87:PRO:HD3	1.67	0.93
4:E:86:SER:HB3	4:E:87:PRO:HD3	1.52	0.92
2:B:4:THR:HG23	2:B:5:PRO:HD2	1.59	0.84
4:E:102:THR:HG22	4:E:188:TRP:HB2	1.60	0.83
4:D:110:LEU:HD13	4:D:177:CYS:SG	2.21	0.81
4:E:19:LYS:HB2	4:E:22:GLU:HG3	1.62	0.81
4:E:136:MET:HG2	4:E:149:ARG:HH21	1.44	0.81
4:E:86:SER:CB	4:E:87:PRO:HD3	2.12	0.80
4:D:136:MET:HE3	4:D:180:SER:OG	1.82	0.78
4:D:125:THR:HG22	4:D:165:PRO:HA	1.66	0.78
1:A:147:TRP:CZ2	3:C:9:LEU:HD13	2.19	0.77
2:B:40:LEU:HD22	2:B:44:GLU:HA	1.66	0.77
1:A:159:TYR:CE2	3:C:3:VAL:HG23	2.20	0.75
2:B:29:GLY:HA2	2:B:61:SER:HB2	1.67	0.75
4:D:126:LEU:HD12	4:D:139:LEU:HD21	1.67	0.75
1:A:15:PRO:HB3	1:A:90:ALA:HA	1.69	0.74
2:B:73:THR:HG22	2:B:75:LYS:H	1.52	0.74
1:A:237:GLY:HA3	5:A:314:HOH:O	1.86	0.74
4:E:136:MET:CG	4:E:149:ARG:HH21	2.01	0.73
4:D:17:LEU:HD13	4:D:100:VAL:HB	1.69	0.73
1:A:35:ARG:HD3	1:A:48:ARG:CZ	2.19	0.72
2:B:84:HIS:CD2	2:B:85:VAL:H	2.07	0.72
4:D:56:HIS:NE2	4:D:61:LYS:HE3	2.05	0.72
4:E:52:ILE:H	4:E:52:ILE:HD12	1.55	0.71
4:D:17:LEU:CD1	4:D:100:VAL:HB	2.20	0.71
2:B:30:PHE:H	2:B:30:PHE:HD1	1.39	0.71
1:A:43:PRO:HG2	1:A:68:LYS:NZ	2.07	0.70
1:A:258:THR:HG22	1:A:273:ARG:HG2	1.74	0.69
1:A:12:VAL:HG22	1:A:94:ILE:HG12	1.74	0.69
4:D:119:LEU:CD2	4:D:200:ILE:HG22	2.23	0.69
1:A:182:ALA:HB1	1:A:265:GLY:HA2	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:149:ARG:HD3	4:D:149:ARG:C	2.19	0.66
1:A:163:LEU:HD12	1:A:163:LEU:H	1.59	0.66
4:D:35:GLU:HG2	4:D:85:HIS:NE2	2.11	0.66
1:A:82:ARG:HB2	5:A:299:HOH:O	1.96	0.66
4:E:17:LEU:HD23	4:E:100:VAL:HB	1.77	0.65
4:E:156:VAL:HG11	5:E:222:HOH:O	1.96	0.65
1:A:2:SER:HB2	1:A:103:VAL:O	1.96	0.65
4:D:49:LEU:N	4:D:49:LEU:HD12	2.10	0.65
4:D:70:MET:HA	4:D:101:ILE:HD12	1.78	0.65
4:E:136:MET:HE3	4:E:180:SER:OG	1.97	0.65
1:A:43:PRO:HD2	5:A:321:HOH:O	1.96	0.65
4:D:136:MET:CG	4:D:149:ARG:HH21	2.10	0.65
1:A:98:MET:HG3	5:A:346:HOH:O	1.96	0.64
1:A:33:PHE:HB2	1:A:52:VAL:HG11	1.79	0.64
1:A:207:GLY:HA2	1:A:240:THR:HB	1.78	0.64
2:B:97:ARG:HG2	5:B:106:HOH:O	1.98	0.64
4:D:75:GLY:HA2	4:D:186:TYR:CD2	2.32	0.64
1:A:111:ARG:NH1	1:A:128:GLU:HG3	2.13	0.63
4:E:136:MET:HG2	4:E:149:ARG:NH2	2.12	0.63
1:A:181:ARG:HG2	1:A:182:ALA:H	1.62	0.63
4:E:149:ARG:HD3	4:E:149:ARG:C	2.22	0.63
4:D:119:LEU:HD23	4:D:200:ILE:HG22	1.80	0.63
4:D:87:PRO:O	4:D:88:TYR:HB2	1.98	0.63
1:A:35:ARG:HD2	2:B:53:ASP:OD1	1.98	0.63
1:A:213:ILE:HG22	5:A:312:HOH:O	1.99	0.63
4:E:135:ASP:OD2	4:E:181:PHE:HA	1.99	0.63
4:D:44:LYS:HG2	4:D:45:PHE:CD2	2.33	0.63
4:D:18:VAL:HB	4:D:69:MET:HG2	1.81	0.62
2:B:73:THR:HG22	2:B:74:GLU:N	2.13	0.62
4:D:86:SER:CB	4:D:87:PRO:HD3	2.29	0.62
1:A:194:VAL:HG23	1:A:195:SER:H	1.64	0.62
4:D:51:LEU:HD22	4:D:64:PHE:CD1	2.34	0.62
4:E:17:LEU:CD2	4:E:100:VAL:HB	2.30	0.62
4:E:140:SER:OG	4:E:143:GLY:HA2	1.99	0.62
4:E:149:ARG:HD3	4:E:149:ARG:O	2.00	0.62
4:D:54:GLU:HA	4:D:54:GLU:OE1	2.00	0.62
2:B:39:LEU:HD13	2:B:68:THR:HG22	1.83	0.61
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.82	0.61
2:B:84:HIS:HD2	2:B:85:VAL:H	1.49	0.61
4:E:29:TRP:HB3	4:E:61:LYS:HG2	1.83	0.61
2:B:4:THR:CG2	2:B:5:PRO:HD2	2.30	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:19:LYS:O	4:D:22:GLU:HB2	2.01	0.60
4:E:200:ILE:HG22	4:E:200:ILE:O	2.01	0.60
2:B:23:LEU:O	2:B:67:TYR:HA	2.02	0.60
4:E:12:ALA:HB2	4:E:26:LEU:HD22	1.82	0.60
4:E:137:TYR:HD2	4:E:162:ALA:HB2	1.65	0.60
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.37	0.60
4:D:18:VAL:HG21	4:D:69:MET:HE2	1.84	0.60
1:A:163:LEU:HD12	1:A:163:LEU:N	2.17	0.60
4:D:200:ILE:HG23	4:E:48:THR:OG1	2.02	0.60
4:E:16:ARG:HG3	4:E:98:ASP:O	2.01	0.60
4:E:70:MET:HA	4:E:101:ILE:CD1	2.32	0.60
1:A:28:VAL:O	1:A:29:ASP:HB2	2.02	0.59
2:B:4:THR:HG23	2:B:87:LEU:HD21	1.83	0.59
1:A:114:ASP:OD2	1:A:156:LEU:HD13	2.03	0.59
1:A:138:THR:HA	1:A:141:GLN:HG3	1.84	0.59
2:B:29:GLY:HA2	2:B:61:SER:CB	2.33	0.59
4:D:88:TYR:HB3	4:D:90:LEU:CD2	2.32	0.59
4:E:156:VAL:O	4:E:157:ASN:HB2	2.02	0.59
4:D:184:SER:HB3	4:D:187:GLU:CD	2.28	0.58
4:E:89:GLN:HB2	5:E:216:HOH:O	2.03	0.58
1:A:205:ALA:O	1:A:208:PHE:HE2	1.86	0.58
1:A:207:GLY:HA2	1:A:240:THR:CB	2.33	0.58
1:A:234:ARG:CD	1:A:242:GLN:HB2	2.33	0.58
4:D:100:VAL:HG11	4:D:188:TRP:CE2	2.38	0.58
4:E:161:GLN:HG2	4:E:162:ALA:H	1.68	0.58
2:B:30:PHE:HE1	2:B:63:TYR:HA	1.68	0.58
4:E:55:HIS:O	4:E:56:HIS:HB3	2.03	0.58
4:E:24:VAL:HG22	4:E:25:ILE:N	2.19	0.58
4:D:24:VAL:HG22	4:D:25:ILE:N	2.18	0.57
1:A:204:TRP:HB3	1:A:206:LEU:HD11	1.85	0.57
1:A:206:LEU:N	1:A:206:LEU:HD12	2.19	0.57
1:A:81:LEU:HD23	1:A:118:TYR:CD1	2.39	0.57
1:A:189:VAL:HA	1:A:202:ARG:O	2.03	0.57
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.40	0.57
4:E:51:LEU:HB2	4:E:64:PHE:CE1	2.40	0.57
4:E:137:TYR:HD2	4:E:162:ALA:CB	2.16	0.56
4:E:109:SER:O	4:E:128:CYS:HA	2.05	0.56
4:E:142:GLU:O	4:E:144:GLU:N	2.38	0.56
1:A:213:ILE:HG12	1:A:214:THR:H	1.71	0.56
1:A:234:ARG:HD3	1:A:242:GLN:HB2	1.88	0.56
4:D:174:THR:HA	4:D:195:LEU:O	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ARG:HD2	2:B:53:ASP:CG	2.30	0.56
4:D:50:HIS:CD2	4:D:88:TYR:OH	2.59	0.56
4:D:100:VAL:HG11	4:D:188:TRP:CD2	2.41	0.56
4:E:87:PRO:O	4:E:88:TYR:O	2.23	0.56
4:E:138:HIS:HB2	4:E:178:PHE:HB2	1.87	0.56
4:D:70:MET:HA	4:D:101:ILE:CD1	2.35	0.56
1:A:224:GLN:O	1:A:226:GLN:N	2.39	0.56
4:D:200:ILE:CG2	4:E:48:THR:HG21	2.34	0.56
4:E:54:GLU:OE1	4:E:54:GLU:HA	2.06	0.56
4:E:125:THR:HB	4:E:165:PRO:HA	1.88	0.56
1:A:79:ARG:HG3	1:A:79:ARG:HH11	1.69	0.56
1:A:201:LEU:HD23	1:A:201:LEU:H	1.70	0.56
4:D:126:LEU:HG	4:D:166:LEU:HD11	1.88	0.55
4:D:149:ARG:HD3	4:D:149:ARG:O	2.07	0.55
4:D:195:LEU:HD23	4:D:196:LEU:N	2.21	0.55
2:B:27:VAL:HG23	2:B:27:VAL:O	2.07	0.55
4:D:138:HIS:ND1	4:D:188:TRP:CZ2	2.75	0.55
4:D:44:LYS:HG2	4:D:45:PHE:CE2	2.41	0.55
1:A:21:HIS:NE2	1:A:23:ILE:HD11	2.22	0.55
4:D:141:ARG:HD3	4:D:144:GLU:HG3	1.88	0.55
1:A:217:TRP:HB2	1:A:228:THR:HG23	1.89	0.55
4:D:83:VAL:HG23	5:D:244:HOH:O	2.07	0.55
1:A:213:ILE:HG12	1:A:214:THR:N	2.22	0.55
4:D:80:TYR:CE1	4:D:93:PRO:HA	2.42	0.55
4:E:10:LEU:HG	4:E:97:LEU:HD22	1.89	0.54
1:A:131:ARG:NH2	1:A:157:ARG:HH21	2.05	0.54
1:A:133:TRP:HB2	1:A:144:GLN:NE2	2.22	0.54
4:E:70:MET:HA	4:E:101:ILE:HD12	1.88	0.54
4:E:156:VAL:HG21	5:E:222:HOH:O	2.07	0.54
1:A:123:TYR:HD2	1:A:124:ILE:HG22	1.71	0.54
1:A:27:TYR:CE2	1:A:32:GLN:HB2	2.41	0.54
1:A:243:LYS:HG2	1:A:244:TRP:N	2.22	0.54
4:D:27:GLN:HG3	4:D:63:ASN:ND2	2.23	0.54
4:E:86:SER:CB	4:E:87:PRO:CD	2.85	0.54
2:B:33:SER:O	2:B:34:ASP:C	2.51	0.54
4:E:59:VAL:HG12	4:E:60:SER:N	2.23	0.54
4:D:80:TYR:HE1	4:D:93:PRO:HB3	1.73	0.54
1:A:124:ILE:HG12	1:A:125:ALA:N	2.22	0.53
1:A:217:TRP:CD1	1:A:247:VAL:HG13	2.43	0.53
1:A:8:PHE:HB2	1:A:25:VAL:HG22	1.89	0.53
4:D:74:ALA:HB1	4:D:186:TYR:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:136:MET:HE2	4:D:181:PHE:C	2.34	0.53
1:A:156:LEU:HD11	5:A:307:HOH:O	2.08	0.53
2:B:84:HIS:CD2	2:B:85:VAL:N	2.75	0.53
4:E:184:SER:HB3	4:E:187:GLU:HG3	1.91	0.53
4:E:142:GLU:C	4:E:144:GLU:H	2.15	0.53
4:E:33:ARG:NH1	4:E:55:HIS:ND1	2.57	0.53
4:E:80:TYR:HE1	4:E:93:PRO:HB3	1.74	0.53
4:D:141:ARG:NH1	4:D:173:GLY:HA3	2.24	0.53
1:A:217:TRP:CD1	1:A:228:THR:HG23	2.44	0.53
4:D:101:ILE:HG22	4:D:186:TYR:O	2.09	0.52
4:E:80:TYR:HB2	4:E:90:LEU:HD12	1.91	0.52
1:A:217:TRP:HD1	1:A:228:THR:HG23	1.74	0.52
1:A:35:ARG:HH12	1:A:37:ASP:HB2	1.74	0.52
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.92	0.52
1:A:217:TRP:CZ3	1:A:258:THR:C	2.87	0.52
4:D:141:ARG:O	4:D:142:GLU:C	2.50	0.52
4:E:80:TYR:CE1	4:E:93:PRO:HB3	2.44	0.52
1:A:273:ARG:HG3	1:A:273:ARG:HH11	1.73	0.52
2:B:19:LYS:O	2:B:71:THR:HG23	2.10	0.52
4:D:136:MET:HB3	4:D:149:ARG:HH21	1.75	0.52
1:A:81:LEU:HD22	1:A:95:ILE:HD11	1.92	0.52
1:A:32:GLN:NE2	1:A:48:ARG:HG2	2.25	0.51
1:A:47:PRO:HB3	1:A:60:TRP:CZ2	2.45	0.51
4:E:57:ASP:HB3	5:E:204:HOH:O	2.10	0.51
4:D:52:ILE:H	4:D:52:ILE:CD1	1.97	0.51
1:A:182:ALA:HB1	1:A:265:GLY:CA	2.40	0.51
1:A:233:THR:OG1	1:A:243:LYS:HE2	2.11	0.51
2:B:67:TYR:N	2:B:67:TYR:CD1	2.79	0.51
1:A:163:LEU:H	1:A:163:LEU:CD1	2.23	0.51
2:B:16:GLU:O	2:B:16:GLU:HG3	2.10	0.51
1:A:202:ARG:HH11	1:A:202:ARG:HG2	1.74	0.51
4:D:52:ILE:HD12	4:D:52:ILE:N	2.10	0.51
2:B:21:ASN:O	2:B:70:PHE:N	2.41	0.51
2:B:37:VAL:HB	2:B:66:TYR:CZ	2.47	0.51
1:A:33:PHE:O	1:A:48:ARG:N	2.44	0.50
1:A:80:ASN:O	1:A:84:TYR:HD1	1.93	0.50
1:A:196:ASP:O	1:A:197:HIS:CB	2.58	0.50
4:D:80:TYR:CE1	4:D:93:PRO:HB3	2.45	0.50
1:A:13:SER:HB3	1:A:78:LEU:HD13	1.93	0.50
1:A:202:ARG:HG2	1:A:246:ALA:HB2	1.93	0.50
1:A:51:TRP:CZ3	1:A:171:TYR:HB3	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:84:THR:HG21	4:D:89:GLN:OE1	2.11	0.50
4:E:126:LEU:CD1	4:E:166:LEU:HD11	2.42	0.50
1:A:224:GLN:O	1:A:225:THR:C	2.55	0.50
4:E:52:ILE:HD12	4:E:52:ILE:N	2.23	0.50
4:E:118:VAL:O	4:E:199:VAL:HA	2.11	0.50
4:D:44:LYS:HB3	5:D:207:HOH:O	2.11	0.49
4:D:156:VAL:O	4:D:156:VAL:HG13	2.12	0.49
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.48	0.49
2:B:36:GLU:HG3	5:B:102:HOH:O	2.12	0.49
2:B:67:TYR:N	2:B:67:TYR:HD1	2.10	0.49
1:A:81:LEU:C	1:A:83:GLY:N	2.68	0.49
1:A:182:ALA:O	1:A:209:TYR:O	2.31	0.49
4:D:138:HIS:O	4:D:177:CYS:HA	2.12	0.49
4:D:154:PRO:HB3	5:D:203:HOH:O	2.12	0.49
4:E:39:LEU:HD21	4:E:77:TYR:HB3	1.93	0.49
1:A:70:GLN:O	1:A:73:THR:N	2.42	0.49
1:A:127:ASN:ND2	1:A:134:THR:OG1	2.36	0.49
4:D:51:LEU:HD22	4:D:64:PHE:CE1	2.48	0.49
4:E:147:GLU:OE1	4:E:147:GLU:HA	2.12	0.49
1:A:209:TYR:CD1	1:A:209:TYR:C	2.91	0.49
4:D:118:VAL:O	4:D:199:VAL:HA	2.12	0.49
4:E:195:LEU:C	4:E:195:LEU:HD23	2.38	0.49
1:A:93:HIS:HD1	1:A:119:ASP:CG	2.19	0.49
4:E:105:TYR:HB3	4:E:134:TYR:OH	2.13	0.49
1:A:35:ARG:NH1	1:A:37:ASP:HB2	2.28	0.49
1:A:185:PRO:HB3	1:A:208:PHE:CD2	2.47	0.49
1:A:273:ARG:HG3	1:A:273:ARG:NH1	2.27	0.48
1:A:117:ALA:HB2	2:B:60:TRP:CZ2	2.48	0.48
2:B:23:LEU:C	2:B:24:ASN:HD22	2.22	0.48
4:D:195:LEU:HD23	4:D:195:LEU:C	2.38	0.48
4:E:39:LEU:HD21	4:E:77:TYR:HD1	1.79	0.48
4:D:184:SER:HB3	4:D:187:GLU:OE2	2.14	0.48
1:A:190:THR:OG1	1:A:192:HIS:HE1	1.97	0.48
1:A:234:ARG:CG	1:A:234:ARG:HH11	2.27	0.48
2:B:74:GLU:HA	2:B:97:ARG:HH22	1.79	0.48
4:D:33:ARG:HG3	4:D:33:ARG:HH11	1.79	0.48
4:D:85:HIS:O	4:D:87:PRO:N	2.47	0.48
1:A:14:ARG:CZ	1:A:21:HIS:HB2	2.44	0.48
1:A:267:PRO:HG2	1:A:268:GLU:H	1.78	0.47
4:E:113:GLN:HE21	4:E:113:GLN:CA	2.25	0.47
1:A:98:MET:HE1	1:A:113:TYR:CE1	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:TYR:HA	1:A:210:PRO:C	2.40	0.47
2:B:2:GLN:HB3	2:B:86:THR:HG22	1.96	0.47
1:A:187:THR:O	1:A:188:HIS:HB3	2.14	0.47
4:D:80:TYR:HB2	4:D:90:LEU:CD1	2.44	0.47
4:E:71:GLN:NE2	4:E:187:GLU:OE2	2.42	0.47
4:D:66:ILE:O	4:D:67:GLY:O	2.33	0.47
4:E:103:GLY:HA2	4:E:190:ASN:CG	2.40	0.47
4:E:183:ASP:OD1	4:E:184:SER:N	2.47	0.47
2:B:39:LEU:CD1	2:B:68:THR:HG22	2.44	0.47
4:D:17:LEU:N	4:D:17:LEU:HD22	2.29	0.47
1:A:19:GLU:HB3	1:A:20:PRO:HD2	1.96	0.47
1:A:132:SER:HA	5:A:305:HOH:O	2.15	0.47
1:A:143:THR:CG2	3:C:9:LEU:HD12	2.45	0.47
4:E:82:SER:HB2	4:E:89:GLN:H	1.79	0.47
4:E:85:HIS:O	4:E:87:PRO:N	2.47	0.47
1:A:253:GLU:OE1	1:A:256:ARG:NH1	2.47	0.47
1:A:274:TRP:O	1:A:276:PRO:HD3	2.13	0.47
4:D:23:THR:HG22	4:D:24:VAL:N	2.29	0.47
2:B:54:LEU:HA	2:B:64:LEU:HD21	1.97	0.47
4:D:109:SER:O	4:D:128:CYS:HA	2.15	0.47
4:E:68:PRO:O	4:E:70:MET:HG2	2.15	0.47
4:E:107:LYS:NZ	5:E:201:HOH:O	2.48	0.47
4:E:119:LEU:O	4:E:122:GLU:HB3	2.15	0.46
1:A:61:ASP:O	1:A:65:GLN:HG2	2.15	0.46
2:B:54:LEU:HA	2:B:64:LEU:CD2	2.46	0.46
4:D:136:MET:HB3	4:D:149:ARG:HE	1.79	0.46
1:A:249:VAL:HG13	1:A:250:PRO:HD2	1.97	0.46
1:A:183:GLU:OE1	1:A:183:GLU:HA	2.14	0.46
1:A:22:PHE:H	1:A:38:SER:HB3	1.80	0.46
1:A:171:TYR:OH	3:C:1:GLY:N	2.49	0.46
2:B:24:ASN:HD22	2:B:24:ASN:N	2.13	0.46
4:D:50:HIS:CG	4:D:51:LEU:N	2.83	0.46
1:A:22:PHE:HB3	1:A:38:SER:HB2	1.97	0.46
2:B:73:THR:CG2	2:B:74:GLU:N	2.77	0.46
4:D:12:ALA:HB2	4:D:26:LEU:HD22	1.98	0.46
1:A:78:LEU:O	1:A:79:ARG:C	2.57	0.46
1:A:176:LYS:HG3	1:A:177:GLU:N	2.29	0.46
4:D:85:HIS:O	4:D:86:SER:C	2.58	0.46
1:A:215:LEU:CD2	1:A:261:VAL:HG22	2.46	0.46
2:B:33:SER:HB3	2:B:62:PHE:CE2	2.51	0.46
4:D:136:MET:CB	4:D:149:ARG:HH21	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:75:GLY:HA2	4:D:186:TYR:CE2	2.50	0.46
4:E:50:HIS:CG	4:E:51:LEU:N	2.84	0.46
4:E:55:HIS:O	4:E:56:HIS:CB	2.63	0.46
1:A:111:ARG:HH11	1:A:128:GLU:HG3	1.80	0.45
4:D:9:SER:HB3	4:D:29:TRP:CE2	2.51	0.45
4:E:101:ILE:HG23	4:E:101:ILE:O	2.16	0.45
4:E:112:ALA:HB2	4:E:126:LEU:CD2	2.46	0.45
4:E:137:TYR:CD2	4:E:162:ALA:CB	2.99	0.45
1:A:190:THR:OG1	1:A:192:HIS:CE1	2.69	0.45
1:A:252:GLY:C	1:A:254:GLU:H	2.24	0.45
1:A:249:VAL:HG22	1:A:257:TYR:CE2	2.51	0.45
1:A:141:GLN:HE21	1:A:141:GLN:HB3	1.59	0.45
1:A:267:PRO:HG2	1:A:268:GLU:N	2.31	0.45
4:E:49:LEU:HD12	4:E:49:LEU:N	2.32	0.45
4:D:22:GLU:O	4:D:69:MET:HB2	2.16	0.45
1:A:121:LYS:O	1:A:122:ASP:C	2.60	0.45
1:A:144:GLN:HG3	1:A:148:GLU:HG3	1.99	0.45
2:B:13:HIS:O	2:B:14:PRO:C	2.60	0.45
2:B:30:PHE:CD1	2:B:30:PHE:N	2.77	0.45
4:D:192:SER:O	4:D:193:ASP:C	2.60	0.45
4:E:170:THR:HB	4:E:171:HIS:CE1	2.51	0.45
4:E:39:LEU:CD2	4:E:77:TYR:HD1	2.30	0.45
1:A:39:ASP:O	1:A:40:ALA:C	2.59	0.45
4:D:80:TYR:CE1	4:D:93:PRO:CA	3.00	0.45
4:E:196:LEU:HD23	4:E:197:VAL:N	2.31	0.45
1:A:137:ASP:OD1	1:A:139:ALA:N	2.50	0.44
4:E:70:MET:O	4:E:71:GLN:C	2.60	0.44
4:E:118:VAL:HG21	4:E:124:VAL:CG2	2.47	0.44
1:A:4:SER:HA	1:A:168:LEU:HD21	1.99	0.44
1:A:51:TRP:H	1:A:51:TRP:CD1	2.35	0.44
4:D:26:LEU:HD12	4:D:39:LEU:HD11	1.98	0.44
4:E:153:GLY:O	4:E:154:PRO:O	2.35	0.44
4:D:136:MET:HB3	4:D:149:ARG:NH2	2.31	0.44
2:B:85:VAL:HG23	2:B:86:THR:N	2.33	0.44
4:D:138:HIS:CD2	4:D:138:HIS:N	2.84	0.44
4:D:97:LEU:HD12	4:D:97:LEU:HA	1.87	0.44
4:D:153:GLY:HA2	4:D:154:PRO:HD2	1.78	0.44
4:E:10:LEU:HD13	4:E:28:CYS:SG	2.58	0.44
4:E:29:TRP:CB	4:E:61:LYS:HG2	2.47	0.44
1:A:159:TYR:CD2	3:C:3:VAL:HG23	2.52	0.44
1:A:181:ARG:HG2	1:A:182:ALA:N	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:CYS:HB3	1:A:272:LEU:HG	2.00	0.44
4:D:17:LEU:HD21	4:D:147:GLU:OE2	2.17	0.44
2:B:69:GLU:O	2:B:70:PHE:HB3	2.16	0.44
2:B:72:PRO:HD2	5:B:119:HOH:O	2.18	0.44
4:D:41:ARG:HA	4:D:76:THR:O	2.17	0.44
4:E:51:LEU:HD12	4:E:51:LEU:HA	1.85	0.44
4:E:154:PRO:O	4:E:156:VAL:HG13	2.17	0.44
1:A:168:LEU:HD12	1:A:172:LEU:HG	1.99	0.44
2:B:29:GLY:CA	2:B:61:SER:HB2	2.44	0.44
1:A:61:ASP:HA	5:A:344:HOH:O	2.18	0.43
4:E:192:SER:O	4:E:193:ASP:C	2.61	0.43
1:A:254:GLU:O	1:A:256:ARG:N	2.51	0.43
4:E:161:GLN:CG	4:E:162:ALA:H	2.30	0.43
1:A:28:VAL:HG23	1:A:33:PHE:CE1	2.53	0.43
1:A:36:PHE:CD2	1:A:67:TYR:HB3	2.52	0.43
1:A:116:TYR:N	1:A:116:TYR:CD1	2.85	0.43
1:A:207:GLY:CA	1:A:240:THR:HB	2.47	0.43
2:B:85:VAL:CG2	2:B:86:THR:N	2.81	0.43
4:E:19:LYS:O	4:E:22:GLU:HB2	2.18	0.43
4:E:20:SER:O	4:E:21:GLU:HB2	2.17	0.43
4:E:142:GLU:C	4:E:144:GLU:N	2.77	0.43
1:A:98:MET:HE3	1:A:115:GLN:NE2	2.33	0.43
1:A:230:LEU:HD23	1:A:230:LEU:N	2.33	0.43
4:E:70:MET:CA	4:E:101:ILE:HD12	2.48	0.43
4:E:142:GLU:O	4:E:144:GLU:HG3	2.17	0.43
1:A:141:GLN:O	1:A:145:ARG:HG3	2.19	0.43
1:A:144:GLN:O	1:A:145:ARG:C	2.59	0.43
1:A:182:ALA:CB	1:A:265:GLY:HA2	2.46	0.43
4:D:50:HIS:HD2	4:D:88:TYR:OH	2.00	0.43
1:A:138:THR:O	1:A:141:GLN:HB2	2.17	0.43
4:E:17:LEU:HD11	4:E:188:TRP:HZ3	1.83	0.43
2:B:39:LEU:HD13	2:B:68:THR:CG2	2.47	0.43
4:D:23:THR:CG2	4:D:24:VAL:N	2.82	0.43
4:D:38:LEU:HD13	4:D:50:HIS:HA	2.00	0.43
1:A:10:THR:HG21	2:B:54:LEU:HD21	2.01	0.43
1:A:78:LEU:O	1:A:80:ASN:N	2.51	0.43
4:D:113:GLN:CA	4:D:113:GLN:HE21	2.32	0.43
1:A:133:TRP:HH2	1:A:156:LEU:HD12	1.84	0.43
1:A:204:TRP:HZ2	2:B:98:ASP:O	2.02	0.43
4:D:20:SER:O	4:D:21:GLU:HB2	2.18	0.43
4:E:13:HIS:HA	4:E:14:PRO:C	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:ARG:NH2	4:D:70:MET:HE1	2.33	0.43
1:A:159:TYR:CD2	3:C:3:VAL:CG2	3.02	0.43
1:A:187:THR:HA	1:A:204:TRP:O	2.19	0.43
4:E:10:LEU:HD11	4:E:26:LEU:HB3	2.00	0.43
4:E:39:LEU:O	4:E:48:THR:HA	2.19	0.43
1:A:53:GLU:O	1:A:53:GLU:HG2	2.19	0.42
1:A:56:GLY:O	1:A:59:TYR:HB3	2.19	0.42
1:A:276:PRO:O	1:A:277:SER:HB3	2.19	0.42
3:C:3:VAL:CG1	3:C:4:ASP:N	2.79	0.42
4:D:40:HIS:ND1	4:D:47:ASP:O	2.51	0.42
1:A:14:ARG:HB3	1:A:17:ARG:HH21	1.85	0.42
1:A:252:GLY:N	1:A:254:GLU:OE2	2.53	0.42
2:B:95:TRP:CZ2	2:B:97:ARG:HB3	2.54	0.42
4:D:48:THR:C	4:D:49:LEU:HD12	2.44	0.42
2:B:31:HIS:ND1	2:B:32:PRO:HA	2.35	0.42
2:B:67:TYR:O	2:B:68:THR:HB	2.18	0.42
4:D:38:LEU:O	4:D:79:CYS:HA	2.19	0.42
4:D:84:THR:CG2	4:D:89:GLN:HB2	2.49	0.42
1:A:147:TRP:CG	1:A:152:GLU:HG3	2.55	0.42
4:D:127:SER:HB2	4:D:161:GLN:NE2	2.34	0.42
1:A:113:TYR:O	1:A:114:ASP:HB2	2.20	0.42
4:E:77:TYR:C	4:E:78:ARG:HG3	2.45	0.42
1:A:79:ARG:HG3	1:A:79:ARG:NH1	2.35	0.42
4:E:105:TYR:HB2	4:E:189:SER:CB	2.50	0.42
4:E:59:VAL:CG1	4:E:60:SER:N	2.83	0.42
4:E:164:PHE:N	4:E:164:PHE:CD2	2.87	0.42
4:D:92:ALA:HB1	4:D:93:PRO:HD2	2.02	0.42
1:A:43:PRO:HG2	1:A:68:LYS:HZ3	1.84	0.41
1:A:201:LEU:HD23	1:A:247:VAL:O	2.20	0.41
4:D:125:THR:CG2	4:D:165:PRO:HA	2.43	0.41
4:D:137:TYR:HD2	4:D:162:ALA:HB2	1.85	0.41
4:E:120:ALA:HA	4:E:199:VAL:CG1	2.49	0.41
4:E:138:HIS:CE1	4:E:149:ARG:HG3	2.55	0.41
4:E:184:SER:CB	4:E:187:GLU:OE1	2.68	0.41
1:A:16:GLY:C	1:A:18:GLY:H	2.27	0.41
4:D:51:LEU:HB2	4:D:64:PHE:CE1	2.54	0.41
4:D:142:GLU:HB2	4:D:143:GLY:H	1.50	0.41
4:E:112:ALA:HB2	4:E:126:LEU:HD23	2.02	0.41
1:A:70:GLN:NE2	3:C:5:PRO:HB3	2.35	0.41
1:A:254:GLU:C	1:A:256:ARG:H	2.29	0.41
4:E:112:ALA:O	4:E:115:GLY:CA	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:196:LEU:HD23	4:E:196:LEU:C	2.45	0.41
1:A:124:ILE:CG1	1:A:125:ALA:N	2.84	0.41
4:D:54:GLU:OE1	4:D:54:GLU:CA	2.67	0.41
1:A:70:GLN:HE21	3:C:5:PRO:HB3	1.84	0.41
2:B:24:ASN:N	2:B:24:ASN:ND2	2.68	0.41
2:B:40:LEU:HD23	2:B:45:ARG:N	2.36	0.41
4:D:13:HIS:HA	4:D:14:PRO:C	2.45	0.41
4:E:51:LEU:HD22	4:E:64:PHE:CD1	2.56	0.41
4:D:136:MET:HB3	4:D:149:ARG:NE	2.35	0.41
1:A:85:TYR:O	1:A:86:ASN:C	2.61	0.41
4:D:70:MET:CA	4:D:101:ILE:HD12	2.49	0.41
4:D:120:ALA:HA	4:D:199:VAL:CG1	2.51	0.41
1:A:34:VAL:HG11	5:A:279:HOH:O	2.20	0.41
1:A:103:VAL:HA	1:A:110:LEU:HD13	2.01	0.41
4:D:70:MET:HB3	4:D:70:MET:HE2	1.62	0.41
4:E:39:LEU:HD12	4:E:64:PHE:CG	2.55	0.41
4:E:134:TYR:HA	4:E:181:PHE:CE1	2.55	0.41
4:E:175:TYR:N	4:E:175:TYR:CD1	2.89	0.41
1:A:16:GLY:O	1:A:18:GLY:N	2.53	0.41
1:A:70:GLN:O	1:A:71:ALA:C	2.64	0.41
1:A:80:ASN:O	1:A:84:TYR:CD1	2.73	0.41
1:A:168:LEU:O	1:A:172:LEU:HG	2.21	0.41
1:A:179:LEU:HD23	1:A:179:LEU:N	2.36	0.41
1:A:216:THR:HG22	1:A:217:TRP:N	2.36	0.41
1:A:260:HIS:HA	1:A:270:LEU:O	2.21	0.41
2:B:31:HIS:HA	2:B:32:PRO:C	2.46	0.41
4:D:27:GLN:HE21	4:D:27:GLN:HB3	1.58	0.41
4:D:141:ARG:HH12	4:D:173:GLY:HA3	1.84	0.41
4:E:95:ASP:HA	4:E:96:PRO:HD3	1.87	0.41
1:A:217:TRP:CE3	1:A:258:THR:O	2.74	0.41
4:D:86:SER:OG	4:D:87:PRO:CD	2.53	0.41
4:E:77:TYR:C	4:E:78:ARG:CG	2.93	0.41
1:A:143:THR:HG21	3:C:9:LEU:HD12	2.02	0.40
4:D:86:SER:CB	4:D:87:PRO:CD	2.98	0.40
4:E:52:ILE:H	4:E:52:ILE:CD1	2.30	0.40
1:A:33:PHE:HB2	1:A:52:VAL:CG1	2.51	0.40
1:A:33:PHE:HD2	1:A:34:VAL:HG13	1.84	0.40
1:A:195:SER:O	1:A:196:ASP:C	2.64	0.40
1:A:217:TRP:HZ3	1:A:258:THR:C	2.29	0.40
1:A:191:HIS:CE1	1:A:199:ALA:HB1	2.57	0.40
4:D:24:VAL:CG2	4:D:25:ILE:N	2.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:71:GLN:NE2	4:D:187:GLU:CD	2.79	0.40
1:A:50:PRO:O	1:A:51:TRP:C	2.64	0.40
1:A:229:GLU:O	1:A:246:ALA:N	2.49	0.40
4:D:55:HIS:O	4:D:56:HIS:HB3	2.21	0.40

There are no symmetry-related clashes.

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	276/278 (99%)	225 (82%)	34 (12%)	17 (6%)	1	6
2	B	98/100 (98%)	79 (81%)	17 (17%)	2 (2%)	6	28
3	C	7/9 (78%)	5 (71%)	2 (29%)	0	100	100
4	D	195/200 (98%)	159 (82%)	28 (14%)	8 (4%)	2	13
4	E	195/200 (98%)	160 (82%)	26 (13%)	9 (5%)	2	11
All	All	771/787 (98%)	628 (82%)	107 (14%)	36 (5%)	2	11

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
4	D	71	GLN
4	D	86	SER
4	D	142	GLU
4	D	154	PRO
4	E	56	HIS
4	E	71	GLN
4	E	86	SER
4	E	88	TYR
4	E	154	PRO

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Mol	Chain	Res	Type
1	A	92	SER
1	A	108	ARG
1	A	114	ASP
1	A	196	ASP
1	A	225	THR
1	A	255	GLN
2	B	17	ASN
4	D	67	GLY
4	E	67	GLY
4	E	143	GLY
1	A	15	PRO
1	A	86	ASN
1	A	251	SER
1	A	276	PRO
2	B	68	THR
4	E	68	PRO
1	A	243	LYS
1	A	253	GLU
4	D	155	LYS
4	D	162	ALA
4	E	162	ALA
1	A	40	ALA
1	A	50	PRO
4	D	68	PRO
1	A	79	ARG
1	A	162	GLY

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	233/233 (100%)	222 (95%)	11 (5%)	23	58
2	B	95/95 (100%)	90 (95%)	5 (5%)	20	54
3	C	6/6 (100%)	4 (67%)	2 (33%)	0	1
4	D	169/172 (98%)	157 (93%)	12 (7%)	13	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	E	169/172 (98%)	160 (95%)	9 (5%)	20	54
All	All	672/678 (99%)	633 (94%)	39 (6%)	18	51

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	15	PRO
1	A	46	GLU
1	A	88	SER
1	A	163	LEU
1	A	169	ARG
1	A	179	LEU
1	A	201	LEU
1	A	206	LEU
1	A	233	THR
1	A	234	ARG
2	B	30	PHE
2	B	54	LEU
2	B	67	TYR
2	B	77	GLU
2	B	83	ASN
3	C	3	VAL
3	C	9	LEU
4	D	35	GLU
4	D	39	LEU
4	D	52	ILE
4	D	71	GLN
4	D	113	GLN
4	D	119	LEU
4	D	128	CYS
4	D	136	MET
4	D	142	GLU
4	D	148	CYS
4	D	190	ASN
4	D	196	LEU
4	E	39	LEU
4	E	85	HIS
4	E	113	GLN
4	E	125	THR
4	E	127	SER
4	E	164	PHE

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Mol	Chain	Res	Type
4	E	168	PRO
4	E	170	THR
4	E	198	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	115	GLN
1	A	127	ASN
1	A	141	GLN
1	A	144	GLN
1	A	155	GLN
1	A	180	GLN
1	A	191	HIS
1	A	192	HIS
1	A	262	GLN
1	A	263	HIS
2	B	24	ASN
2	B	83	ASN
2	B	84	HIS
4	D	27	GLN
4	D	50	HIS
4	D	63	ASN
4	D	71	GLN
4	D	113	GLN
4	D	161	GLN
4	E	113	GLN
4	E	138	HIS

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

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.