



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2021 – 04:29 AM EDT

PDB ID : 1KTL
Title : The human non-classical major histocompatibility complex molecule HLA-E
Authors : Holmes, M.A.; Strong, R.K.
Deposited on : 2002-01-16
Resolution : 3.10 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

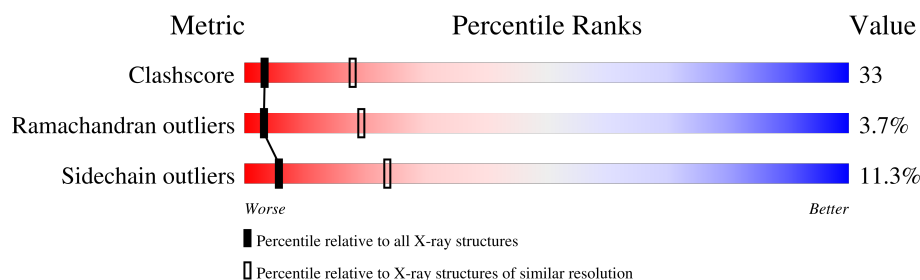
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)

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	274	<div> <div>46%</div> <div>45%</div> <div>8%</div> <div>.</div> </div>
1	C	274	<div> <div>49%</div> <div>36%</div> <div>12%</div> <div>.</div> </div>
2	B	100	<div> <div>56%</div> <div>37%</div> <div>6%</div> <div>.</div> </div>
2	D	100	<div> <div>42%</div> <div>46%</div> <div>10%</div> <div>.</div> </div>
3	P	9	<div> <div>33%</div> <div>44%</div> <div>11%</div> <div>11%</div> </div>
3	Q	9	<div> <div>33%</div> <div>44%</div> <div>11%</div> <div>11%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6211 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 CLASS I HISTOCOMPATIBILITY ANTIGEN, ALPHA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2200	1376	389	428	7			
1	C	274	Total	C	N	O	S	0	0	0
			2212	1383	394	428	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	GLY	ARG	engineered mutation	UNP P13747
A	256	ALA	ARG	conflict	UNP P13747
C	107	GLY	ARG	engineered mutation	UNP P13747
C	256	ALA	ARG	conflict	UNP P13747

- 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
			828	526	139	159	4			
2	D	100	Total	C	N	O	S	0	0	0
			828	526	139	159	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	cloning artifact	UNP P01884
D	1	MET	-	cloning artifact	UNP P01884

- Molecule 3 is a protein called Peptide VTAPRTLTL.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	9	Total	C	N	O	0	0	0
			69	45	12	12			
3	Q	9	Total	C	N	O	0	0	0
			69	45	12	12			

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



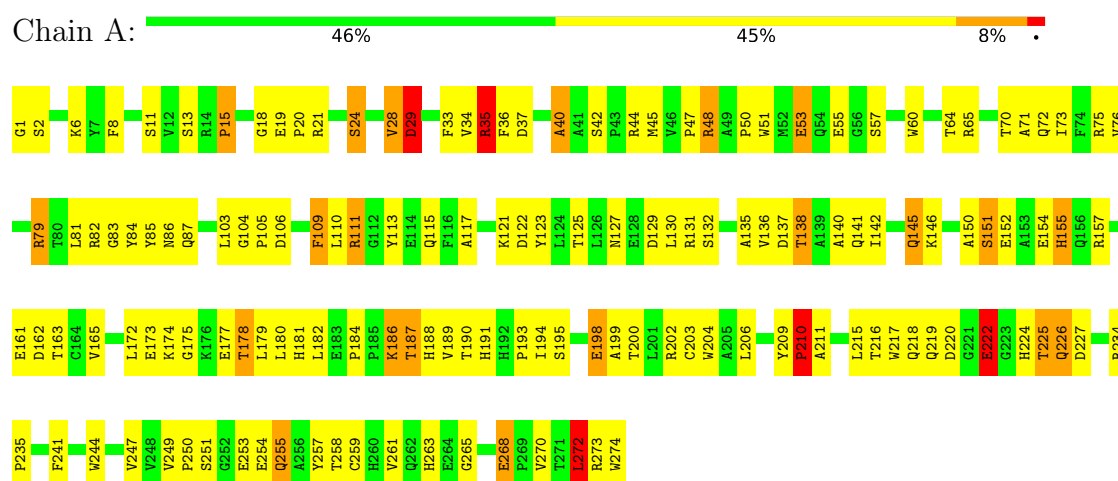
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

3 Residue-property plots

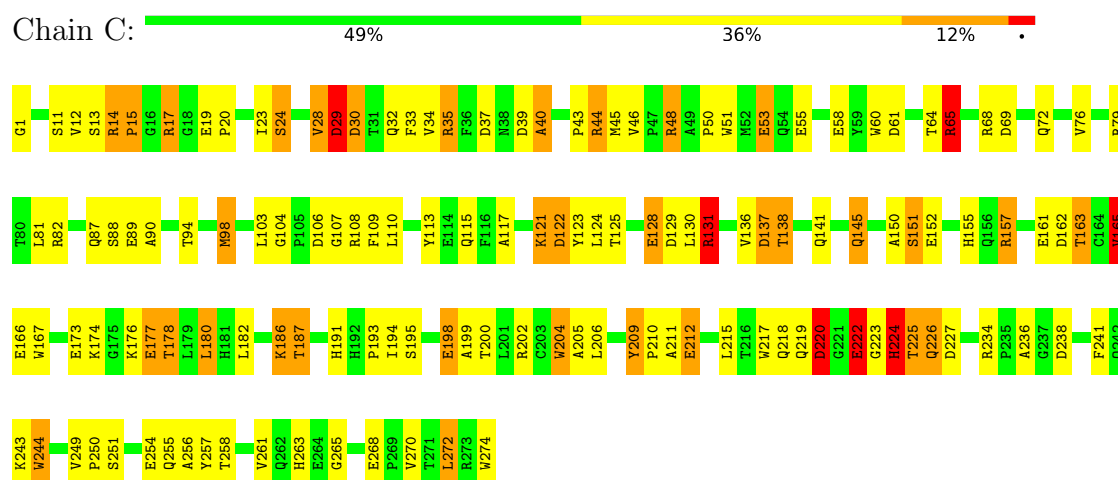
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

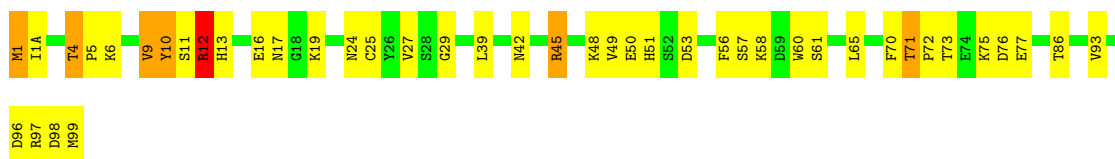


• Molecule 1: HLA CLASS I HISTOCOMPATIBILITY ANTIGEN, ALPHA CHAIN

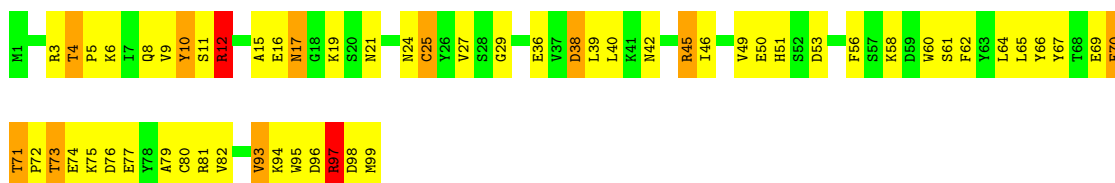


• Molecule 2: BETA-2-MICROGLOBULIN

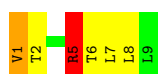




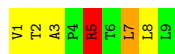
• Molecule 2: BETA-2-MICROGLOBULIN



• Molecule 3: Peptide VTAPRTLTL



• Molecule 3: Peptide VTAPRTLTL



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.40 Å 178.40 Å 87.30 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 3.10	Depositor
% Data completeness (in resolution range)	(Not available) (25.00-3.10)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.238 , 0.280	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6211	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	1.30	3/2265 (0.1%)	1.24	12/3082 (0.4%)
1	C	1.24	11/2278 (0.5%)	1.30	23/3099 (0.7%)
2	B	1.40	4/851 (0.5%)	1.32	9/1152 (0.8%)
2	D	1.28	3/851 (0.4%)	1.26	9/1152 (0.8%)
3	P	2.38	5/69 (7.2%)	3.05	5/93 (5.4%)
3	Q	1.44	1/69 (1.4%)	2.05	3/93 (3.2%)
All	All	1.31	27/6383 (0.4%)	1.32	61/8671 (0.7%)

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	A	0	2
1	C	0	3
3	P	0	1
3	Q	0	1
All	All	0	7

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1	GLY	C-O	9.59	1.39	1.23
3	P	5	ARG	CZ-NH1	-8.86	1.21	1.33
3	P	5	ARG	CG-CD	8.48	1.73	1.51
1	C	131	ARG	CZ-NH2	7.41	1.42	1.33
1	A	21	ARG	CZ-NH1	7.20	1.42	1.33
3	P	2	THR	CB-CG2	-7.20	1.28	1.52
1	A	225	THR	CA-CB	6.72	1.70	1.53
1	C	58	GLU	CG-CD	6.47	1.61	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	157	ARG	CG-CD	-6.26	1.36	1.51
3	P	5	ARG	NE-CZ	-6.22	1.25	1.33
2	B	12	ARG	CZ-NH2	-6.09	1.25	1.33
2	D	10	TYR	C-O	-6.03	1.11	1.23
1	A	154	GLU	CD-OE1	5.92	1.32	1.25
2	B	1	MET	SD-CE	5.85	2.10	1.77
2	B	12	ARG	CZ-NH1	-5.75	1.25	1.33
2	B	9	VAL	CB-CG1	-5.71	1.40	1.52
1	C	204	TRP	CB-CG	-5.70	1.40	1.50
1	C	98	MET	CG-SD	5.64	1.95	1.81
1	C	157	ARG	CZ-NH2	5.63	1.40	1.33
1	C	225	THR	C-O	5.57	1.33	1.23
1	C	1	GLY	N-CA	5.45	1.54	1.46
2	D	70	PHE	CE1-CZ	5.39	1.47	1.37
1	C	165	VAL	CB-CG1	-5.36	1.41	1.52
3	P	1	VAL	C-O	-5.26	1.13	1.23
1	C	212	GLU	CD-OE2	5.23	1.31	1.25
2	D	74	GLU	CD-OE2	5.23	1.31	1.25
3	Q	5	ARG	NE-CZ	-5.13	1.26	1.33

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	5	ARG	NE-CZ-NH1	-17.62	111.49	120.30
3	P	5	ARG	NE-CZ-NH2	12.52	126.56	120.30
1	C	157	ARG	CG-CD-NE	12.02	137.03	111.80
1	C	131	ARG	NE-CZ-NH2	10.97	125.78	120.30
1	A	79	ARG	NE-CZ-NH1	10.73	125.67	120.30
1	A	111	ARG	NE-CZ-NH1	-10.36	115.12	120.30
3	P	5	ARG	CG-CD-NE	-10.17	90.44	111.80
3	Q	5	ARG	NE-CZ-NH2	-10.16	115.22	120.30
2	B	45	ARG	NE-CZ-NH1	9.70	125.15	120.30
2	B	45	ARG	NE-CZ-NH2	-9.69	115.45	120.30
1	C	65	ARG	NE-CZ-NH1	9.39	124.99	120.30
1	A	180	LEU	CA-CB-CG	-8.74	95.21	115.30
2	D	12	ARG	NE-CZ-NH2	8.44	124.52	120.30
1	C	220	ASP	CB-CG-OD1	8.40	125.86	118.30
2	D	45	ARG	NE-CZ-NH1	-8.27	116.17	120.30
3	Q	5	ARG	CG-CD-NE	7.80	128.18	111.80
2	B	12	ARG	NE-CZ-NH1	7.67	124.14	120.30
1	A	35	ARG	NE-CZ-NH1	7.57	124.08	120.30
1	A	28	VAL	N-CA-C	-7.53	90.66	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	97	ARG	NE-CZ-NH1	7.17	123.88	120.30
1	C	180	LEU	CA-CB-CG	-7.11	98.95	115.30
1	C	131	ARG	NH1-CZ-NH2	-7.11	111.58	119.40
1	C	122	ASP	CB-CG-OD2	-7.10	111.91	118.30
1	A	79	ARG	NE-CZ-NH2	-6.94	116.83	120.30
2	B	10	TYR	CB-CG-CD2	-6.84	116.90	121.00
1	C	44	ARG	NE-CZ-NH1	-6.78	116.91	120.30
2	D	25	CYS	CA-CB-SG	-6.76	101.83	114.00
1	C	39	ASP	CB-CG-OD1	-6.75	112.22	118.30
2	B	10	TYR	CB-CG-CD1	6.71	125.02	121.00
3	P	5	ARG	NH1-CZ-NH2	-6.69	112.04	119.40
2	D	45	ARG	NE-CZ-NH2	6.68	123.64	120.30
1	C	79	ARG	NE-CZ-NH1	-6.67	116.96	120.30
1	C	28	VAL	N-CA-C	-6.67	93.00	111.00
1	A	157	ARG	CG-CD-NE	-6.62	97.90	111.80
2	D	38	ASP	CB-CG-OD2	-6.61	112.35	118.30
1	C	137	ASP	CB-CG-OD2	-6.50	112.45	118.30
2	B	97	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	C	131	ARG	CG-CD-NE	-6.40	98.37	111.80
1	C	48	ARG	NE-CZ-NH1	-6.39	117.11	120.30
3	Q	5	ARG	NE-CZ-NH1	6.36	123.48	120.30
1	C	157	ARG	NH1-CZ-NH2	-6.25	112.53	119.40
1	A	21	ARG	NE-CZ-NH2	-6.18	117.21	120.30
1	A	75	ARG	NE-CZ-NH1	-5.81	117.40	120.30
1	C	14	ARG	NE-CZ-NH1	5.79	123.19	120.30
1	C	30	ASP	N-CA-C	5.77	126.59	111.00
1	C	131	ARG	NE-CZ-NH1	-5.67	117.46	120.30
1	C	186	LYS	N-CA-C	-5.65	95.73	111.00
2	D	12	ARG	NH1-CZ-NH2	-5.63	113.20	119.40
1	C	61	ASP	CB-CG-OD1	-5.58	113.27	118.30
2	B	11	SER	N-CA-CB	-5.54	102.19	110.50
2	B	12	ARG	NH1-CZ-NH2	-5.51	113.34	119.40
1	A	24	SER	N-CA-CB	-5.44	102.34	110.50
1	C	17	ARG	N-CA-C	5.43	125.67	111.00
1	A	35	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	A	272	LEU	CA-CB-CG	5.28	127.44	115.30
2	D	97	ARG	NE-CZ-NH2	5.27	122.93	120.30
2	D	64	LEU	CA-CB-CG	5.26	127.39	115.30
1	C	14	ARG	NE-CZ-NH2	-5.18	117.71	120.30
3	P	2	THR	N-CA-C	-5.12	97.18	111.00
1	C	107	GLY	N-CA-C	5.10	125.85	113.10
2	D	97	ARG	CD-NE-CZ	5.05	130.66	123.60

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	209	TYR	Sidechain,Mainchain
1	C	113	TYR	Sidechain
1	C	131	ARG	Sidechain
1	C	209	TYR	Sidechain
3	P	5	ARG	Sidechain
3	Q	5	ARG	Sidechain

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	2200	0	2022	157	1
1	C	2212	0	2039	152	0
2	B	828	0	780	50	0
2	D	828	0	780	58	0
3	P	69	0	83	21	0
3	Q	69	0	83	18	0
4	A	5	0	0	1	0
All	All	6211	0	5787	393	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 33.

All (393) 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:B:1:MET:SD	2:B:1:MET:CE	2.10	1.39
1:A:115:GLN:HG2	1:A:125:THR:HG23	1.34	1.10
1:C:65:ARG:HH11	1:C:65:ARG:CB	1.69	1.06
1:A:65:ARG:HH21	1:C:145:GLN:NE2	1.56	1.02
1:A:263:HIS:HD2	1:A:265:GLY:H	1.10	1.00
1:C:65:ARG:HH11	1:C:65:ARG:HB2	1.26	0.97
1:C:35:ARG:HH11	1:C:35:ARG:HG2	1.30	0.96
3:P:5:ARG:HE	3:Q:5:ARG:HH21	1.15	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:ARG:HH11	1:A:65:ARG:CB	1.81	0.92
2:D:98:ASP:O	2:D:99:MET:HG3	1.69	0.92
1:C:152:GLU:CB	3:Q:7:LEU:HD11	2.00	0.92
1:C:65:ARG:HB2	1:C:65:ARG:NH1	1.85	0.89
1:A:263:HIS:CD2	1:A:265:GLY:H	1.89	0.89
1:C:263:HIS:CD2	1:C:265:GLY:H	1.91	0.88
1:C:115:GLN:HG2	1:C:125:THR:HG23	1.54	0.87
1:C:155:HIS:CE1	3:P:5:ARG:NH2	2.43	0.86
1:A:103:LEU:HD21	1:A:165:VAL:HG23	1.57	0.86
1:C:263:HIS:HD2	1:C:265:GLY:H	1.24	0.85
1:C:204:TRP:HE3	1:C:206:LEU:HD21	1.43	0.83
2:D:9:VAL:CG2	2:D:93:VAL:CG2	2.57	0.83
2:D:42:ASN:HD21	2:D:77:GLU:H	1.27	0.82
2:B:73:THR:HG22	2:B:75:LYS:H	1.45	0.81
1:A:109:PHE:CE2	1:A:161:GLU:HG2	2.17	0.80
2:B:9:VAL:CG2	2:B:93:VAL:HG22	2.10	0.80
2:D:4:THR:HG22	2:D:5:PRO:HD2	1.63	0.80
1:A:65:ARG:HH11	1:A:65:ARG:HB3	1.45	0.80
1:A:152:GLU:CB	3:P:7:LEU:HD11	2.11	0.79
2:D:42:ASN:ND2	2:D:77:GLU:H	1.81	0.79
1:C:193:PRO:HA	1:C:199:ALA:HA	1.63	0.79
1:C:35:ARG:HH11	1:C:35:ARG:CG	1.96	0.79
2:B:73:THR:HG22	2:B:75:LYS:N	1.98	0.79
1:A:50:PRO:HA	1:A:53:GLU:OE2	1.83	0.78
2:B:42:ASN:HD21	2:B:77:GLU:H	1.31	0.78
2:B:42:ASN:ND2	2:B:77:GLU:H	1.81	0.78
1:A:202:ARG:HG3	1:A:202:ARG:HH11	1.48	0.78
2:D:9:VAL:HG23	2:D:93:VAL:CG2	2.16	0.76
1:C:65:ARG:HH11	1:C:65:ARG:HB3	1.51	0.76
1:A:129:ASP:OD2	1:A:131:ARG:HG3	1.86	0.75
1:A:65:ARG:HB2	1:A:65:ARG:NH1	2.00	0.75
1:C:109:PHE:CE2	1:C:161:GLU:HG2	2.22	0.75
1:C:138:THR:HA	1:C:141:GLN:HE21	1.52	0.74
1:C:152:GLU:CG	3:Q:7:LEU:HD13	2.18	0.74
1:C:44:ARG:HA	1:C:64:THR:HG23	1.69	0.74
1:A:79:ARG:HH11	1:C:72:GLN:NE2	1.84	0.74
1:C:117:ALA:HB2	2:D:60:TRP:CE2	2.23	0.74
1:A:204:TRP:HE3	1:A:206:LEU:HD21	1.51	0.73
2:D:29:GLY:HA2	2:D:61:SER:OG	1.89	0.73
1:C:152:GLU:HB2	3:Q:7:LEU:HD11	1.70	0.73
2:D:73:THR:HG22	2:D:75:LYS:H	1.54	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:49:VAL:HG12	2:B:50:GLU:N	2.04	0.72
1:A:145:GLN:HG3	1:A:146:LYS:N	2.05	0.72
1:C:152:GLU:CB	3:Q:7:LEU:CD1	2.68	0.71
1:A:65:ARG:HH21	1:C:145:GLN:HE22	1.38	0.71
2:D:42:ASN:HD22	2:D:77:GLU:HB2	1.56	0.71
2:D:73:THR:HG22	2:D:75:LYS:N	2.06	0.71
3:P:5:ARG:HE	3:Q:5:ARG:NH2	1.87	0.70
1:A:65:ARG:CB	1:A:65:ARG:NH1	2.54	0.70
1:A:65:ARG:HH21	1:C:145:GLN:HE21	1.38	0.70
2:B:9:VAL:CG2	2:B:93:VAL:CG2	2.70	0.70
1:A:65:ARG:NH2	1:C:145:GLN:NE2	2.36	0.69
1:C:33:PHE:CD2	1:C:34:VAL:HG13	2.27	0.69
1:A:65:ARG:HH11	1:A:65:ARG:HB2	1.53	0.69
1:A:79:ARG:NH1	1:C:72:GLN:NE2	2.41	0.69
1:A:121:LYS:HG3	1:A:122:ASP:N	2.08	0.69
1:C:98:MET:HE2	1:C:115:GLN:OE1	1.92	0.69
2:B:4:THR:HG23	2:B:86:THR:OG1	1.92	0.68
1:C:152:GLU:CG	3:Q:7:LEU:CD1	2.71	0.68
1:C:28:VAL:O	1:C:29:ASP:O	2.10	0.68
1:C:152:GLU:CD	3:Q:7:LEU:HD13	2.14	0.68
1:A:55:GLU:HA	1:A:55:GLU:OE1	1.92	0.68
1:A:152:GLU:HB2	3:P:7:LEU:HD11	1.75	0.68
1:A:79:ARG:NH2	4:A:275:SO4:O2	2.25	0.68
1:A:195:SER:OG	1:A:198:GLU:HG3	1.94	0.68
1:A:152:GLU:CB	3:P:7:LEU:CD1	2.71	0.67
1:C:193:PRO:HA	1:C:199:ALA:CB	2.25	0.67
1:C:204:TRP:CE3	1:C:206:LEU:HD21	2.29	0.67
1:C:50:PRO:HA	1:C:53:GLU:OE2	1.94	0.66
3:P:5:ARG:NE	3:Q:5:ARG:HH21	1.90	0.66
1:A:255:GLN:CD	1:A:255:GLN:N	2.48	0.66
1:C:193:PRO:HA	1:C:199:ALA:CA	2.25	0.66
2:B:9:VAL:HG21	2:B:93:VAL:HG22	1.77	0.66
2:D:4:THR:CG2	2:D:5:PRO:HD2	2.26	0.66
2:D:96:ASP:O	2:D:98:ASP:N	2.29	0.65
1:C:35:ARG:NH1	1:C:46:VAL:HG21	2.11	0.65
1:A:255:GLN:H	1:A:255:GLN:NE2	1.94	0.65
2:D:39:LEU:HB3	2:D:46:ILE:HD12	1.78	0.65
1:A:187:THR:HB	1:A:272:LEU:HD11	1.78	0.65
1:A:48:ARG:NH1	2:B:53:ASP:OD2	2.30	0.65
1:A:79:ARG:HD2	1:C:72:GLN:NE2	2.12	0.65
1:C:249:VAL:HG22	1:C:257:TYR:CE1	2.32	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:TRP:HB2	1:A:174:LYS:O	1.97	0.64
1:A:193:PRO:HA	1:A:199:ALA:HA	1.79	0.64
1:A:8:PHE:HB3	2:B:56:PHE:CE2	2.33	0.64
1:A:109:PHE:CD1	1:A:110:LEU:N	2.66	0.64
1:A:194:ILE:HD12	1:A:194:ILE:N	2.13	0.64
2:D:42:ASN:ND2	2:D:77:GLU:HB2	2.14	0.63
2:B:73:THR:CG2	2:B:75:LYS:H	2.11	0.63
1:A:81:LEU:HD21	1:A:123:TYR:CZ	2.34	0.63
1:A:145:GLN:HG3	1:A:146:LYS:H	1.60	0.63
1:A:182:LEU:HD23	1:A:210:PRO:HD3	1.80	0.63
1:A:249:VAL:HG22	1:A:257:TYR:CE1	2.34	0.63
2:B:42:ASN:HD22	2:B:77:GLU:HB2	1.64	0.63
1:A:172:LEU:HD23	1:A:179:LEU:HD23	1.81	0.62
1:A:138:THR:HA	1:A:141:GLN:HG3	1.80	0.62
1:A:225:THR:O	1:A:227:ASP:N	2.31	0.62
2:D:9:VAL:CG2	2:D:93:VAL:HG23	2.27	0.62
2:B:49:VAL:CG1	2:B:50:GLU:N	2.63	0.62
1:A:193:PRO:HA	1:A:199:ALA:CB	2.29	0.62
2:B:4:THR:HG22	2:B:5:PRO:HD2	1.80	0.62
1:C:152:GLU:HB3	3:Q:7:LEU:HD11	1.80	0.62
1:C:215:LEU:CD2	1:C:261:VAL:HG22	2.30	0.62
2:D:9:VAL:CG2	2:D:93:VAL:HG22	2.28	0.62
1:C:194:ILE:HD12	1:C:194:ILE:N	2.14	0.61
2:D:73:THR:CG2	2:D:75:LYS:H	2.11	0.61
1:A:72:GLN:HE21	1:C:76:VAL:HG13	1.64	0.61
1:A:187:THR:O	1:A:188:HIS:HB3	2.00	0.61
1:A:152:GLU:CG	3:P:7:LEU:HD13	2.30	0.61
1:C:137:ASP:O	1:C:141:GLN:HG3	2.01	0.61
2:D:16:GLU:HB3	2:D:19:LYS:CB	2.31	0.61
1:C:155:HIS:CE1	3:P:5:ARG:HH21	2.18	0.61
1:C:212:GLU:HA	1:C:212:GLU:OE1	2.01	0.60
2:D:16:GLU:O	2:D:19:LYS:N	2.30	0.60
1:C:150:ALA:O	1:C:151:SER:CB	2.49	0.60
2:D:51:HIS:HA	2:D:65:LEU:O	2.01	0.60
1:A:13:SER:HA	1:A:20:PRO:HB3	1.84	0.60
1:A:270:VAL:HG12	1:A:272:LEU:CD2	2.31	0.60
1:C:82:ARG:HE	1:C:89:GLU:HB2	1.66	0.60
1:C:152:GLU:HG3	3:Q:7:LEU:CD1	2.32	0.60
1:C:155:HIS:CE1	3:P:5:ARG:HH22	2.19	0.60
1:C:81:LEU:HD21	1:C:123:TYR:CZ	2.36	0.59
1:A:225:THR:C	1:A:227:ASP:H	2.05	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:32:GLN:NE2	2:D:53:ASP:OD2	2.35	0.59
1:C:234:ARG:NH1	2:D:8:GLN:OE1	2.34	0.59
1:C:209:TYR:HA	1:C:210:PRO:O	2.03	0.59
1:A:73:ILE:HD11	3:Q:8:LEU:HD22	1.84	0.59
1:C:255:GLN:H	1:C:255:GLN:NE2	2.01	0.59
1:A:152:GLU:HB3	3:P:7:LEU:HD11	1.85	0.58
1:C:51:TRP:HB2	1:C:174:LYS:O	2.03	0.58
1:A:250:PRO:HG2	1:A:253:GLU:OE2	2.03	0.58
1:C:65:ARG:CB	1:C:65:ARG:NH1	2.47	0.58
1:A:263:HIS:HD2	1:A:265:GLY:N	1.92	0.58
2:B:9:VAL:HG23	2:B:93:VAL:CG2	2.33	0.58
2:B:42:ASN:ND2	2:B:77:GLU:HB2	2.18	0.58
1:A:35:ARG:HG2	1:A:35:ARG:HH11	1.68	0.58
1:A:202:ARG:HG3	1:A:202:ARG:NH1	2.15	0.58
1:C:89:GLU:O	1:C:89:GLU:HG2	2.04	0.58
1:C:109:PHE:CD1	1:C:110:LEU:N	2.71	0.58
3:P:5:ARG:HH11	3:Q:5:ARG:HH21	1.52	0.58
1:A:79:ARG:HH11	1:C:72:GLN:HE21	1.52	0.57
1:A:194:ILE:HD11	1:A:200:THR:OG1	2.04	0.57
1:A:155:HIS:CE1	3:P:5:ARG:NH2	2.72	0.57
1:C:98:MET:CE	1:C:115:GLN:OE1	2.51	0.57
1:C:234:ARG:HG3	2:D:10:TYR:CZ	2.39	0.57
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.39	0.57
1:C:211:ALA:HB2	1:C:241:PHE:CE1	2.40	0.57
2:D:25:CYS:HB2	2:D:39:LEU:HD21	1.87	0.57
2:D:73:THR:HG22	2:D:76:ASP:H	1.70	0.57
1:A:6:LYS:HD2	1:A:113:TYR:OH	2.05	0.56
1:A:152:GLU:HB3	3:P:7:LEU:CD1	2.35	0.56
1:C:35:ARG:HG2	1:C:35:ARG:NH1	2.10	0.56
1:C:43:PRO:O	1:C:68:ARG:NH2	2.39	0.56
1:C:236:ALA:O	2:D:12:ARG:HD3	2.05	0.56
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.87	0.56
2:B:98:ASP:C	2:B:99:MET:HG3	2.25	0.56
1:A:109:PHE:CE2	1:A:161:GLU:HA	2.41	0.55
1:A:193:PRO:HA	1:A:199:ALA:CA	2.36	0.55
1:A:227:ASP:O	1:A:247:VAL:HG23	2.06	0.55
1:C:162:ASP:O	1:C:163:THR:C	2.44	0.55
2:B:51:HIS:HA	2:B:65:LEU:O	2.07	0.55
1:A:109:PHE:HD1	1:A:110:LEU:N	2.03	0.55
1:C:45:MET:O	1:C:60:TRP:CE3	2.60	0.55
1:C:12:VAL:HG22	1:C:94:THR:HG23	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:194:ILE:HD11	1:C:200:THR:OG1	2.06	0.55
2:D:49:VAL:HG12	2:D:50:GLU:N	2.21	0.55
1:A:150:ALA:O	1:A:151:SER:CB	2.56	0.54
1:C:35:ARG:HD3	2:D:53:ASP:CG	2.28	0.54
1:C:152:GLU:HB3	3:Q:7:LEU:CD1	2.38	0.54
1:A:55:GLU:CD	1:A:174:LYS:NZ	2.60	0.54
1:A:35:ARG:HG3	1:A:36:PHE:N	2.20	0.54
1:C:249:VAL:HG13	1:C:250:PRO:HD2	1.90	0.54
1:C:117:ALA:HB2	2:D:60:TRP:CD2	2.44	0.53
1:A:204:TRP:CE3	1:A:206:LEU:HD21	2.37	0.53
2:D:9:VAL:HG23	2:D:93:VAL:HG23	1.85	0.53
1:A:136:VAL:O	1:A:136:VAL:CG1	2.56	0.53
1:A:225:THR:C	1:A:227:ASP:N	2.62	0.53
1:C:254:GLU:HG2	1:C:274:TRP:CE3	2.44	0.53
1:C:225:THR:CG2	1:C:226:GLN:N	2.72	0.53
2:B:4:THR:CG2	2:B:5:PRO:HD2	2.38	0.53
1:C:223:GLY:O	1:C:224:HIS:O	2.27	0.53
2:D:39:LEU:CB	2:D:46:ILE:HD12	2.37	0.53
1:C:35:ARG:CG	1:C:35:ARG:NH1	2.67	0.52
1:C:55:GLU:HA	1:C:55:GLU:OE1	2.09	0.52
3:Q:2:THR:HG22	3:Q:3:ALA:N	2.23	0.52
3:Q:5:ARG:HH11	3:Q:5:ARG:CG	2.23	0.52
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.44	0.52
1:A:189:VAL:HG12	1:A:190:THR:N	2.24	0.52
1:C:55:GLU:CD	1:C:174:LYS:NZ	2.63	0.52
1:C:128:GLU:O	1:C:130:LEU:HG	2.09	0.52
1:C:13:SER:HA	1:C:20:PRO:HB3	1.92	0.52
2:D:49:VAL:CG1	2:D:50:GLU:N	2.72	0.52
1:C:88:SER:OG	1:C:90:ALA:HB3	2.10	0.52
2:B:29:GLY:HA2	2:B:61:SER:OG	2.11	0.52
1:C:35:ARG:HH12	1:C:46:VAL:HG21	1.74	0.52
1:A:1:GLY:O	1:A:105:PRO:HA	2.10	0.51
2:B:10:TYR:O	2:B:24:ASN:HB2	2.10	0.51
2:B:45:ARG:O	2:B:45:ARG:HG3	2.10	0.51
1:C:103:LEU:HA	1:C:108:ARG:O	2.10	0.51
2:D:9:VAL:HG21	2:D:93:VAL:CG2	2.39	0.51
1:C:69:ASP:O	1:C:72:GLN:HB2	2.10	0.51
2:D:29:GLY:HA2	2:D:61:SER:CB	2.40	0.51
2:B:9:VAL:HG21	2:B:93:VAL:CG2	2.39	0.51
1:C:37:ASP:HB3	1:C:40:ALA:HB2	1.92	0.51
1:C:219:GLN:HG3	1:C:219:GLN:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:LYS:CG	1:A:122:ASP:N	2.74	0.51
1:A:152:GLU:CG	3:P:7:LEU:CD1	2.89	0.51
1:A:35:ARG:HD3	2:B:53:ASP:OD2	2.11	0.51
1:C:103:LEU:HD21	1:C:165:VAL:HG23	1.93	0.50
1:C:152:GLU:OE1	3:P:5:ARG:NH1	2.44	0.50
3:Q:5:ARG:HH11	3:Q:5:ARG:HG2	1.76	0.50
2:B:4:THR:HG22	2:B:5:PRO:CD	2.42	0.50
1:C:82:ARG:NE	1:C:89:GLU:HB2	2.26	0.50
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.46	0.50
2:D:9:VAL:HA	2:D:24:ASN:O	2.12	0.50
1:A:109:PHE:HE1	1:A:111:ARG:CA	2.25	0.50
1:C:176:LYS:O	1:C:178:THR:N	2.41	0.50
1:A:135:ALA:HB1	1:A:140:ALA:CB	2.42	0.49
1:A:218:GLN:HB2	1:A:222:GLU:O	2.12	0.49
1:C:122:ASP:O	1:C:136:VAL:HG11	2.12	0.49
1:C:129:ASP:OD2	1:C:131:ARG:HB2	2.12	0.49
2:D:9:VAL:HG21	2:D:93:VAL:HG23	1.95	0.49
1:A:210:PRO:O	1:A:263:HIS:HE1	1.95	0.49
1:C:263:HIS:HD2	1:C:265:GLY:N	2.02	0.49
1:C:82:ARG:CG	1:C:87:GLN:O	2.61	0.49
2:B:39:LEU:HD23	2:B:39:LEU:HA	1.60	0.48
1:C:194:ILE:H	1:C:199:ALA:HA	1.77	0.48
2:D:98:ASP:C	2:D:99:MET:HG3	2.32	0.48
1:A:184:PRO:HB3	1:A:265:GLY:O	2.13	0.48
3:P:5:ARG:HH11	3:Q:5:ARG:NH2	2.10	0.48
1:C:217:TRP:O	1:C:224:HIS:HA	2.13	0.48
1:C:219:GLN:HB3	1:C:257:TYR:CE2	2.49	0.48
1:A:175:GLY:O	1:A:179:LEU:HB2	2.14	0.48
1:A:255:GLN:CD	1:A:255:GLN:H	2.15	0.48
1:A:109:PHE:CD2	1:A:161:GLU:HA	2.49	0.47
1:A:215:LEU:CD2	1:A:261:VAL:HG22	2.44	0.47
2:B:73:THR:HB	2:B:76:ASP:HB2	1.96	0.47
1:C:234:ARG:HG3	2:D:10:TYR:OH	2.14	0.47
1:A:35:ARG:HG2	1:A:35:ARG:NH1	2.27	0.47
1:A:186:LYS:HD3	1:A:186:LYS:N	2.28	0.47
1:A:129:ASP:OD2	1:A:131:ARG:CG	2.59	0.47
1:C:45:MET:O	1:C:60:TRP:HE3	1.97	0.47
1:C:157:ARG:HH21	1:C:157:ARG:HB3	1.79	0.47
1:C:236:ALA:HB3	1:C:238:ASP:OD1	2.15	0.47
2:D:79:ALA:HB2	2:D:94:LYS:HA	1.97	0.47
1:A:152:GLU:CD	3:P:7:LEU:HD13	2.35	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:270:VAL:HG12	1:C:272:LEU:CD2	2.45	0.47
2:D:9:VAL:HG21	2:D:93:VAL:O	2.15	0.47
2:D:51:HIS:HB3	2:D:66:TYR:CD2	2.49	0.47
1:C:191:HIS:ND1	1:C:274:TRP:CH2	2.72	0.47
1:A:155:HIS:CE1	1:C:155:HIS:CE1	3.03	0.47
1:A:235:PRO:O	2:B:10:TYR:OH	2.19	0.47
2:B:16:GLU:HG2	2:B:19:LYS:CB	2.45	0.47
1:C:202:ARG:HG3	1:C:202:ARG:HH11	1.80	0.47
1:C:44:ARG:HA	1:C:64:THR:CG2	2.41	0.47
1:C:194:ILE:N	1:C:194:ILE:CD1	2.77	0.47
1:A:82:ARG:HG3	1:A:87:GLN:HB2	1.98	0.46
1:A:127:ASN:HB2	1:A:132:SER:OG	2.14	0.46
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.97	0.46
1:C:109:PHE:CE2	1:C:161:GLU:HA	2.50	0.46
1:C:219:GLN:O	1:C:220:ASP:C	2.52	0.46
1:C:249:VAL:CG2	1:C:257:TYR:CE1	2.99	0.46
1:C:35:ARG:NH1	1:C:46:VAL:CG2	2.78	0.46
1:A:83:GLY:O	1:A:84:TYR:C	2.53	0.46
1:A:227:ASP:O	1:A:227:ASP:OD1	2.33	0.46
1:C:121:LYS:HG3	1:C:122:ASP:N	2.31	0.46
2:D:4:THR:HG22	2:D:5:PRO:CD	2.39	0.46
2:D:36:GLU:O	2:D:82:VAL:HA	2.16	0.46
1:A:103:LEU:CD2	1:A:165:VAL:HG23	2.38	0.46
1:C:180:LEU:HD23	1:C:180:LEU:HA	1.67	0.46
1:C:195:SER:OG	1:C:198:GLU:HG3	2.15	0.46
2:B:96:ASP:OD1	2:B:98:ASP:HB2	2.16	0.46
1:C:124:LEU:HD12	1:C:124:LEU:HA	1.73	0.46
1:C:28:VAL:O	1:C:29:ASP:C	2.55	0.45
1:C:205:ALA:O	1:C:206:LEU:HD23	2.17	0.45
2:D:16:GLU:O	2:D:17:ASN:C	2.55	0.45
2:D:40:LEU:HD11	2:D:81:ARG:HB2	1.98	0.45
1:A:211:ALA:HB2	1:A:241:PHE:CE1	2.52	0.45
1:C:176:LYS:C	1:C:178:THR:H	2.19	0.45
2:D:16:GLU:HG2	2:D:19:LYS:CB	2.47	0.45
1:A:234:ARG:HG3	2:B:10:TYR:CZ	2.51	0.45
1:A:28:VAL:O	1:A:29:ASP:C	2.54	0.45
1:A:42:SER:O	1:A:44:ARG:N	2.44	0.45
1:A:109:PHE:HE1	1:A:111:ARG:HA	1.82	0.45
1:A:211:ALA:HB2	1:A:241:PHE:CD1	2.52	0.45
2:B:73:THR:CG2	2:B:76:ASP:H	2.30	0.45
1:A:135:ALA:HB1	1:A:140:ALA:HB3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:THR:O	1:A:188:HIS:CB	2.64	0.45
1:A:219:GLN:HG3	1:A:219:GLN:O	2.17	0.45
2:B:12:ARG:HG2	2:B:13:HIS:NE2	2.32	0.45
1:A:44:ARG:HA	1:A:64:THR:HG23	1.99	0.45
1:A:104:GLY:HA3	1:A:106:ASP:OD1	2.17	0.45
1:A:33:PHE:CD1	1:A:33:PHE:N	2.84	0.45
2:D:3:ARG:HG2	2:D:3:ARG:HH11	1.82	0.45
1:A:142:ILE:HG23	1:C:65:ARG:HE	1.82	0.44
1:A:191:HIS:NE2	1:A:199:ALA:CB	2.80	0.44
1:C:243:LYS:O	1:C:244:TRP:HB3	2.17	0.44
1:C:182:LEU:HD23	1:C:182:LEU:HA	1.71	0.44
2:D:71:THR:HA	2:D:72:PRO:HD2	1.69	0.44
2:B:73:THR:HG22	2:B:76:ASP:H	1.82	0.44
1:C:219:GLN:HA	1:C:256:ALA:O	2.17	0.44
1:A:150:ALA:O	1:A:151:SER:HB2	2.17	0.44
2:B:6:LYS:HB2	2:B:6:LYS:HE3	1.89	0.44
2:B:57:SER:O	2:B:60:TRP:N	2.49	0.44
1:C:109:PHE:CD1	1:C:109:PHE:C	2.90	0.44
1:C:209:TYR:HA	1:C:210:PRO:C	2.36	0.44
2:D:15:ALA:HB2	2:D:95:TRP:CZ2	2.52	0.44
1:A:225:THR:CG2	1:A:226:GLN:N	2.80	0.44
1:A:255:GLN:N	1:A:255:GLN:NE2	2.63	0.44
1:C:131:ARG:HH11	1:C:157:ARG:NH1	2.16	0.44
1:A:79:ARG:NH1	1:C:72:GLN:HE22	2.15	0.44
1:C:243:LYS:HG2	1:C:244:TRP:N	2.33	0.43
1:A:152:GLU:CB	3:P:7:LEU:HD13	2.48	0.43
2:B:16:GLU:HB3	2:B:19:LYS:CB	2.48	0.43
1:C:162:ASP:HB2	1:C:163:THR:H	1.45	0.43
1:C:194:ILE:HD13	1:C:199:ALA:C	2.38	0.43
1:A:47:PRO:HB3	1:A:60:TRP:CZ2	2.54	0.43
1:A:85:TYR:O	1:A:86:ASN:C	2.56	0.43
1:A:249:VAL:CG2	1:A:257:TYR:CE1	3.00	0.43
1:A:37:ASP:HB3	1:A:40:ALA:HB2	2.00	0.43
1:A:162:ASP:HB2	1:A:163:THR:H	1.42	0.43
1:A:121:LYS:HG3	1:A:122:ASP:H	1.80	0.43
1:A:65:ARG:NH2	1:C:145:GLN:HE21	2.08	0.43
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.54	0.43
1:A:142:ILE:HG23	1:C:65:ARG:NE	2.34	0.43
1:A:129:ASP:O	1:A:130:LEU:HB2	2.18	0.43
1:A:249:VAL:HA	1:A:250:PRO:HD3	1.80	0.43
1:A:137:ASP:OD2	1:A:138:THR:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:VAL:CG1	1:A:190:THR:N	2.82	0.43
1:A:194:ILE:N	1:A:194:ILE:CD1	2.80	0.43
2:B:29:GLY:HA2	2:B:61:SER:CB	2.49	0.43
1:A:24:SER:HB3	1:A:36:PHE:HB3	2.00	0.42
1:A:270:VAL:HG12	1:A:272:LEU:HD23	2.01	0.42
1:C:44:ARG:HH11	1:C:44:ARG:HD3	1.69	0.42
2:D:39:LEU:C	2:D:46:ILE:HD12	2.39	0.42
2:D:39:LEU:HD23	2:D:39:LEU:HA	1.70	0.42
1:A:104:GLY:C	1:A:106:ASP:N	2.71	0.42
1:A:109:PHE:CD1	1:A:109:PHE:C	2.93	0.42
1:A:188:HIS:HA	1:A:272:LEU:HD12	2.01	0.42
1:C:187:THR:HB	1:C:272:LEU:HD11	2.01	0.42
1:A:216:THR:O	1:A:259:CYS:HA	2.18	0.42
2:B:27:VAL:O	2:B:27:VAL:HG23	2.20	0.42
1:A:76:VAL:HG11	3:P:8:LEU:HD11	2.01	0.42
1:C:163:THR:O	1:C:167:TRP:HD1	2.03	0.42
1:A:73:ILE:HG21	3:P:6:THR:HG23	2.01	0.42
1:C:104:GLY:C	1:C:106:ASP:N	2.73	0.42
1:C:244:TRP:C	1:C:244:TRP:HE3	2.23	0.42
2:B:16:GLU:O	2:B:17:ASN:C	2.58	0.42
2:B:49:VAL:CG1	2:B:50:GLU:H	2.33	0.42
1:C:109:PHE:HD1	1:C:110:LEU:N	2.18	0.42
1:C:255:GLN:NE2	1:C:255:GLN:N	2.66	0.42
1:A:191:HIS:ND1	1:A:274:TRP:CH2	2.75	0.42
1:C:109:PHE:CD2	1:C:161:GLU:HA	2.55	0.42
1:C:136:VAL:O	1:C:136:VAL:CG1	2.66	0.42
2:D:6:LYS:O	2:D:27:VAL:HA	2.19	0.42
1:A:198:GLU:CB	1:A:249:VAL:O	2.68	0.41
2:D:72:PRO:HB2	2:D:97:ARG:NH1	2.34	0.41
2:B:45:ARG:O	2:B:45:ARG:CG	2.67	0.41
2:D:38:ASP:O	2:D:80:CYS:HA	2.20	0.41
1:A:137:ASP:H	1:A:140:ALA:HB3	1.84	0.41
1:C:211:ALA:HB2	1:C:241:PHE:CD1	2.55	0.41
1:C:218:GLN:HB2	1:C:222:GLU:O	2.20	0.41
1:A:181:HIS:CE1	1:A:182:LEU:O	2.73	0.41
1:A:255:GLN:C	1:A:257:TYR:N	2.72	0.41
2:B:71:THR:HA	2:B:72:PRO:HD2	1.77	0.41
1:C:23:ILE:HG22	1:C:24:SER:N	2.36	0.41
1:C:150:ALA:O	1:C:151:SER:HB2	2.21	0.41
1:A:8:PHE:HD2	2:B:56:PHE:CE1	2.39	0.41
1:A:121:LYS:HB2	2:B:1(A):ILE:HG13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:MET:O	1:A:60:TRP:CE3	2.73	0.41
1:A:109:PHE:HE2	1:A:161:GLU:HA	1.84	0.41
1:C:122:ASP:HB3	1:C:136:VAL:HG21	2.03	0.41
2:D:56:PHE:HA	2:D:62:PHE:HA	2.03	0.41
2:D:21:ASN:O	2:D:69:GLU:HG3	2.20	0.41
1:A:145:GLN:CD	1:C:65:ARG:HH21	2.23	0.41
1:C:166:GLU:O	1:C:166:GLU:HG2	2.21	0.41
1:C:270:VAL:HG12	1:C:272:LEU:HD23	2.02	0.41
1:C:225:THR:O	1:C:227:ASP:N	2.54	0.40
1:A:70:THR:O	1:A:71:ALA:C	2.58	0.40
1:A:222:GLU:H	1:A:222:GLU:HG3	1.57	0.40
1:C:212:GLU:OE1	1:C:212:GLU:CA	2.66	0.40
2:D:67:TYR:CD1	2:D:67:TYR:N	2.88	0.40
1:A:155:HIS:NE2	1:C:151:SER:HB3	2.37	0.40
1:A:162:ASP:O	1:A:163:THR:C	2.59	0.40
1:A:194:ILE:H	1:A:199:ALA:HA	1.86	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:A:111:ARG:NH1	1:A:268:GLU:OE2[5_555]	2.08	0.12

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	272/274 (99%)	232 (85%)	25 (9%)	15 (6%)	2	11
1	C	272/274 (99%)	231 (85%)	30 (11%)	11 (4%)	3	17
2	B	98/100 (98%)	92 (94%)	6 (6%)	0	100	100
2	D	98/100 (98%)	85 (87%)	11 (11%)	2 (2%)	7	31

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	P	7/9 (78%)	7 (100%)	0	0	100	100
3	Q	7/9 (78%)	7 (100%)	0	0	100	100
All	All	754/766 (98%)	654 (87%)	72 (10%)	28 (4%)	3	19

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	177	GLU
1	A	222	GLU
1	A	224	HIS
1	A	226	GLN
1	C	29	ASP
1	C	177	GLU
1	C	222	GLU
1	C	224	HIS
1	C	226	GLN
2	D	97	ARG
1	A	15	PRO
1	A	29	ASP
1	A	254	GLU
1	C	15	PRO
1	A	2	SER
1	A	178	THR
1	C	30	ASP
1	C	40	ALA
1	C	163	THR
1	C	178	THR
1	C	220	ASP
1	A	40	ALA
1	A	57	SER
1	A	220	ASP
2	D	17	ASN
1	A	18	GLY
1	A	19	GLU
1	A	210	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	229/234 (98%)	204 (89%)	25 (11%)	6	25
1	C	231/234 (99%)	202 (87%)	29 (13%)	4	18
2	B	93/95 (98%)	87 (94%)	6 (6%)	17	47
2	D	93/95 (98%)	83 (89%)	10 (11%)	6	25
3	P	8/8 (100%)	6 (75%)	2 (25%)	0	2
3	Q	8/8 (100%)	5 (62%)	3 (38%)	0	0
All	All	662/674 (98%)	587 (89%)	75 (11%)	6	23

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	15	PRO
1	A	29	ASP
1	A	35	ARG
1	A	48	ARG
1	A	53	GLU
1	A	109	PHE
1	A	138	THR
1	A	145	GLN
1	A	151	SER
1	A	155	HIS
1	A	173	GLU
1	A	178	THR
1	A	186	LYS
1	A	187	THR
1	A	198	GLU
1	A	210	PRO
1	A	222	GLU
1	A	244	TRP
1	A	251	SER
1	A	255	GLN
1	A	258	THR
1	A	268	GLU
1	A	272	LEU
1	A	273	ARG
2	B	4	THR
2	B	12	ARG

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Mol	Chain	Res	Type
2	B	48	LYS
2	B	58	LYS
2	B	70	PHE
2	B	71	THR
1	C	11	SER
1	C	14	ARG
1	C	15	PRO
1	C	17	ARG
1	C	19	GLU
1	C	24	SER
1	C	29	ASP
1	C	35	ARG
1	C	48	ARG
1	C	53	GLU
1	C	65	ARG
1	C	121	LYS
1	C	128	GLU
1	C	138	THR
1	C	145	GLN
1	C	151	SER
1	C	165	VAL
1	C	173	GLU
1	C	177	GLU
1	C	186	LYS
1	C	187	THR
1	C	198	GLU
1	C	222	GLU
1	C	224	HIS
1	C	244	TRP
1	C	251	SER
1	C	258	THR
1	C	268	GLU
1	C	272	LEU
2	D	4	THR
2	D	11	SER
2	D	12	ARG
2	D	45	ARG
2	D	58	LYS
2	D	70	PHE
2	D	71	THR
2	D	73	THR
2	D	93	VAL

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Mol	Chain	Res	Type
2	D	97	ARG
3	P	1	VAL
3	P	5	ARG
3	Q	1	VAL
3	Q	5	ARG
3	Q	7	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	87	GLN
1	A	127	ASN
1	A	141	GLN
1	A	145	GLN
1	A	148	ASN
1	A	181	HIS
1	A	255	GLN
1	A	263	HIS
2	B	2	GLN
2	B	42	ASN
2	B	89	GLN
1	C	72	GLN
1	C	87	GLN
1	C	141	GLN
1	C	145	GLN
1	C	255	GLN
1	C	263	HIS
2	D	2	GLN
2	D	42	ASN
2	D	89	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	A	275	-	4,4,4	1.02	0	6,6,6	0.30	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	275	SO4	1	0

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.