



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:32 PM GMT

PDB ID : 1KPR
Title : The human non-classical major histocompatibility complex molecule HLA-E
Authors : Holmes, M.A.; Strong, R.K.
Deposited on : 2002-01-02
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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

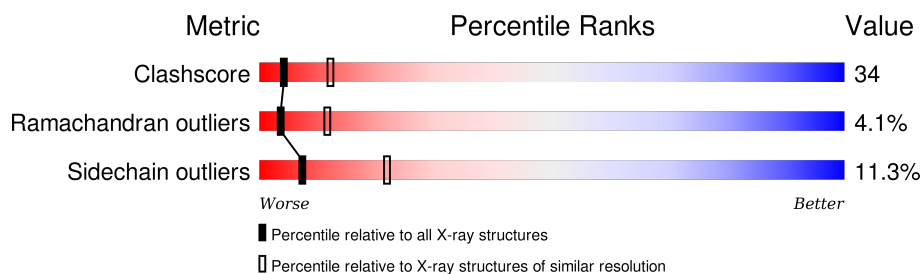
1 Overall quality at a glance

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	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (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	A	274	
1	C	274	
2	B	100	
2	D	100	
3	P	9	
3	Q	9	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6210 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
			2206	1380	391	428	7			
1	C	274	Total	C	N	O	S	0	0	0
			2205	1379	393	426	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	GLY	ARG	ENGINEERED	UNP P13747
A	256	ALA	ARG	ENGINEERED	UNP P13747
C	107	GLY	ARG	ENGINEERED	UNP P13747
C	256	ALA	ARG	ENGINEERED	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 P61769
D	1	MET	-	CLONING ARTIFACT	UNP P61769

- Molecule 3 is a protein called Peptide VMAPRTVLL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	9	Total	C	N	O	S	0	0	0
			69	45	12	11	1			
3	Q	9	Total	C	N	O	S	0	0	0
			69	45	12	11	1			

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



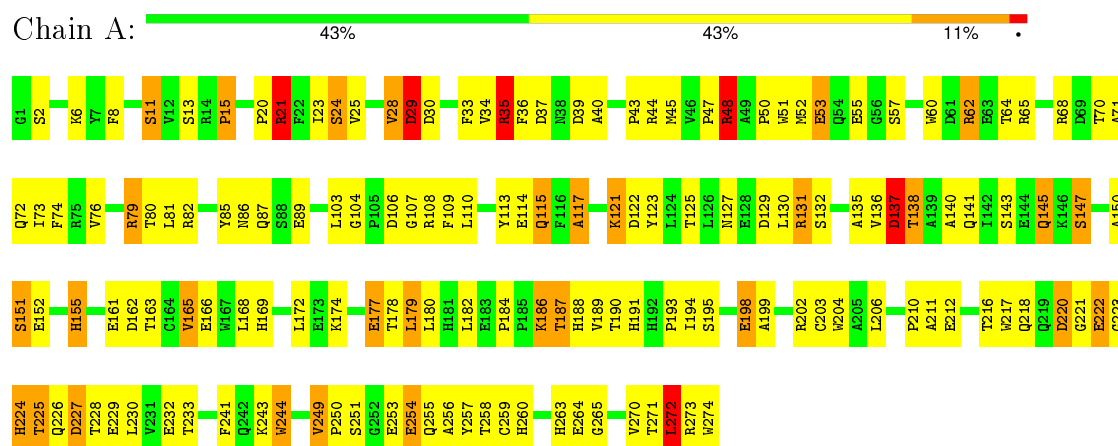
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

3 Residue-property plots

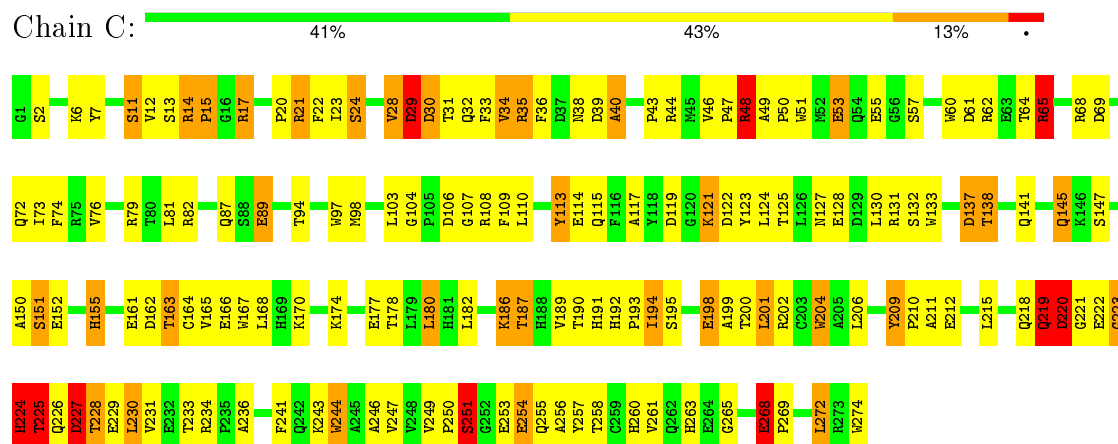
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

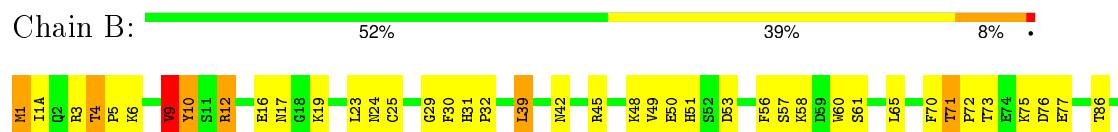
• Molecule 1: HLA CLASS I HISTOCOMPATIBILITY ANTIGEN, ALPHA CHAIN



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



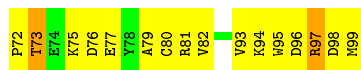
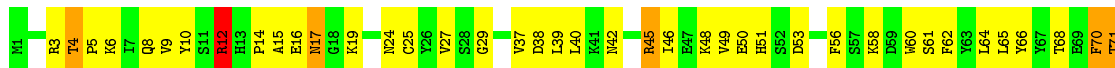
• Molecule 2: BETA-2-MICROGLOBULIN





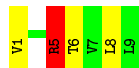
- Molecule 2: BETA-2-MICROGLOBULIN

Chain D: 44% 48% 7%



- Molecule 3: Peptide VMAPRTVLL

Chain P: 56% 33% 11%



- Molecule 3: Peptide VMAPRTVLL

Chain Q: 67% 33%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.70 Å 178.70 Å 88.20 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.80	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.80)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.235 , 0.289	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6210	wwPDB-VP
Average B, all atoms (Å ²)	64.0	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	1.25	5/2272 (0.2%)	1.25	16/3092 (0.5%)
1	C	1.32	9/2271 (0.4%)	1.34	26/3089 (0.8%)
2	B	1.52	6/851 (0.7%)	1.44	9/1152 (0.8%)
2	D	1.15	0/851	1.30	8/1152 (0.7%)
3	P	1.53	0/69	1.45	1/92 (1.1%)
3	Q	1.22	0/69	1.35	0/92
All	All	1.30	20/6383 (0.3%)	1.32	60/8669 (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	C	0	2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	34	VAL	CB-CG1	9.61	1.73	1.52
1	C	256	ALA	CA-CB	9.27	1.72	1.52
2	B	10	TYR	CE2-CZ	8.95	1.50	1.38
2	B	10	TYR	CG-CD1	7.30	1.48	1.39
1	C	224	HIS	CA-CB	7.00	1.69	1.53
1	C	204	TRP	CB-CG	-6.40	1.38	1.50
1	C	268	GLU	CD-OE1	6.06	1.32	1.25
1	A	115	GLN	CB-CG	-6.05	1.36	1.52
1	C	98	MET	CG-SD	5.87	1.96	1.81
1	C	7	TYR	CG-CD1	5.71	1.46	1.39
2	B	10	TYR	C-O	-5.68	1.12	1.23
1	C	133	TRP	CB-CG	-5.60	1.40	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	114	GLU	CB-CG	-5.59	1.41	1.52
2	B	9	VAL	CB-CG2	-5.46	1.41	1.52
2	B	1	MET	SD-CE	5.38	2.08	1.77
1	A	114	GLU	CB-CG	-5.27	1.42	1.52
1	A	48	ARG	CG-CD	-5.25	1.38	1.51
1	A	117	ALA	CA-CB	-5.16	1.41	1.52
1	A	121	LYS	CB-CG	-5.03	1.39	1.52
2	B	30	PHE	CE2-CZ	5.03	1.46	1.37

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	12	ARG	NE-CZ-NH1	10.75	125.68	120.30
1	C	79	ARG	NE-CZ-NH2	-9.78	115.41	120.30
2	B	45	ARG	NE-CZ-NH1	9.52	125.06	120.30
1	C	224	HIS	N-CA-C	8.73	134.57	111.00
1	A	180	LEU	CA-CB-CG	-8.67	95.35	115.30
2	B	97	ARG	NE-CZ-NH1	8.55	124.57	120.30
1	C	131	ARG	NE-CZ-NH1	-8.45	116.07	120.30
2	B	45	ARG	NE-CZ-NH2	-8.30	116.15	120.30
2	B	97	ARG	NE-CZ-NH2	-8.21	116.20	120.30
1	A	28	VAL	N-CA-C	-8.17	88.94	111.00
1	A	62	ARG	NE-CZ-NH1	8.11	124.36	120.30
2	D	45	ARG	NE-CZ-NH1	-8.07	116.26	120.30
1	A	131	ARG	NE-CZ-NH2	-7.89	116.35	120.30
2	D	12	ARG	NE-CZ-NH2	7.70	124.15	120.30
1	A	131	ARG	NE-CZ-NH1	7.57	124.09	120.30
2	D	45	ARG	NE-CZ-NH2	7.46	124.03	120.30
1	A	62	ARG	NE-CZ-NH2	-7.39	116.61	120.30
1	A	224	HIS	N-CA-C	-7.27	91.38	111.00
2	B	12	ARG	NH1-CZ-NH2	-7.18	111.50	119.40
2	B	3	ARG	NE-CZ-NH2	7.09	123.85	120.30
1	C	131	ARG	NE-CZ-NH2	7.04	123.82	120.30
1	C	224	HIS	CA-C-N	-7.04	101.72	117.20
1	C	28	VAL	N-CA-C	-6.84	92.52	111.00
1	C	225	THR	N-CA-C	6.76	129.26	111.00
1	A	24	SER	N-CA-CB	-6.63	100.55	110.50
1	C	62	ARG	NE-CZ-NH1	-6.62	116.99	120.30
2	B	98	ASP	CB-CG-OD1	6.55	124.20	118.30
1	A	137	ASP	CB-CG-OD2	-6.53	112.43	118.30
1	C	65	ARG	NE-CZ-NH1	6.52	123.56	120.30
1	C	224	HIS	C-N-CA	6.52	137.99	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	12	ARG	NH1-CZ-NH2	-6.48	112.27	119.40
1	C	223	GLY	C-N-CA	6.48	137.89	121.70
1	C	180	LEU	CA-CB-CG	-6.30	100.81	115.30
1	C	62	ARG	NE-CZ-NH2	6.24	123.42	120.30
2	D	12	ARG	NE-CZ-NH1	6.21	123.40	120.30
2	D	97	ARG	NE-CZ-NH2	6.17	123.39	120.30
1	A	21	ARG	NE-CZ-NH2	-6.13	117.23	120.30
1	C	119	ASP	CB-CG-OD2	-6.04	112.86	118.30
1	C	227	ASP	CA-C-N	-5.95	104.11	117.20
1	C	48	ARG	NE-CZ-NH1	-5.91	117.35	120.30
1	C	69	ASP	CB-CG-OD1	5.84	123.56	118.30
1	C	107	GLY	N-CA-C	5.78	127.55	113.10
1	C	201	LEU	CA-CB-CG	5.66	128.32	115.30
1	A	225	THR	CA-C-N	-5.60	104.89	117.20
1	C	24	SER	N-CA-CB	-5.58	102.12	110.50
1	C	186	LYS	N-CA-C	-5.44	96.31	111.00
1	A	107	GLY	N-CA-C	5.44	126.69	113.10
1	C	268	GLU	OE1-CD-OE2	5.41	129.79	123.30
1	A	249	VAL	N-CA-CB	-5.31	99.82	111.50
1	A	272	LEU	CA-CB-CG	5.27	127.42	115.30
3	P	5	ARG	NE-CZ-NH1	5.26	122.93	120.30
2	D	64	LEU	CA-CB-CG	5.22	127.31	115.30
2	D	64	LEU	CB-CG-CD1	-5.22	102.13	111.00
2	B	39	LEU	CA-CB-CG	-5.17	103.40	115.30
1	C	113	TYR	CD1-CE1-CZ	-5.15	115.17	119.80
1	A	35	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	A	79	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	C	17	ARG	N-CA-C	5.07	124.69	111.00
1	C	220	ASP	CB-CG-OD1	5.05	122.85	118.30
1	C	30	ASP	N-CA-C	5.00	124.51	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	209	TYR	Sidechain
1	C	22	PHE	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	A	2206	0	2028	150	0
1	C	2205	0	2027	167	0
2	B	828	0	780	49	0
2	D	828	0	780	56	0
3	P	69	0	83	12	0
3	Q	69	0	83	10	0
4	A	5	0	0	1	0
All	All	6210	0	5781	403	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 34.

All (403) 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:CE	2:B:1:MET:SD	2.08	1.41
1:A:218:GLN:HB2	1:A:223:GLY:HA2	1.25	1.16
1:C:65:ARG:HH11	1:C:65:ARG:CB	1.67	1.08
1:C:219:GLN:HG3	1:C:220:ASP:N	1.68	1.04
1:C:65:ARG:HB2	1:C:65:ARG:HH11	1.20	1.03
1:C:219:GLN:HG3	1:C:220:ASP:H	0.89	1.03
2:D:98:ASP:O	2:D:99:MET:HG3	1.59	1.01
1:A:65:ARG:CB	1:A:65:ARG:HH11	1.75	0.99
1:C:263:HIS:HD2	1:C:265:GLY:H	1.07	0.97
1:C:65:ARG:HB2	1:C:65:ARG:NH1	1.80	0.94
1:C:219:GLN:CG	1:C:220:ASP:H	1.80	0.93
1:A:263:HIS:HD2	1:A:265:GLY:H	1.07	0.92
1:A:65:ARG:HH21	1:C:145:GLN:HE22	1.13	0.92
1:C:115:GLN:HG2	1:C:125:THR:HG23	1.51	0.91
1:A:65:ARG:HH21	1:C:145:GLN:NE2	1.67	0.91
1:C:263:HIS:CD2	1:C:265:GLY:H	1.89	0.89
1:A:218:GLN:CB	1:A:223:GLY:HA2	2.03	0.87
1:A:138:THR:HA	1:A:141:GLN:HG3	1.55	0.87
1:A:263:HIS:CD2	1:A:265:GLY:H	1.94	0.85
2:B:9:VAL:CG2	2:B:93:VAL:HG22	2.06	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:ARG:HB2	1:A:65:ARG:NH1	1.91	0.85
1:C:21:ARG:HH11	1:C:21:ARG:HG3	1.42	0.85
1:C:187:THR:HB	1:C:272:LEU:HD11	1.59	0.84
1:A:187:THR:HB	1:A:272:LEU:HD11	1.59	0.83
1:A:65:ARG:HB3	1:A:65:ARG:HH11	1.42	0.82
2:B:73:THR:HG22	2:B:75:LYS:H	1.45	0.82
2:B:73:THR:HG22	2:B:75:LYS:N	1.95	0.81
1:A:115:GLN:HG2	1:A:125:THR:HG23	1.61	0.81
1:C:219:GLN:O	1:C:221:GLY:N	2.14	0.81
1:C:33:PHE:CD2	1:C:34:VAL:HG13	2.16	0.80
1:C:35:ARG:NH1	1:C:46:VAL:HG21	1.97	0.80
2:D:9:VAL:CG2	2:D:93:VAL:CG2	2.60	0.80
2:B:42:ASN:ND2	2:B:77:GLU:H	1.79	0.79
1:A:218:GLN:HB2	1:A:223:GLY:CA	2.10	0.79
3:P:5:ARG:NH2	3:Q:5:ARG:NE	2.30	0.79
1:C:204:TRP:HE3	1:C:206:LEU:HD21	1.49	0.78
1:C:117:ALA:HB2	2:D:60:TRP:CE2	2.18	0.78
1:A:65:ARG:HB2	1:A:65:ARG:HH11	1.43	0.78
2:D:42:ASN:ND2	2:D:77:GLU:H	1.84	0.76
2:D:42:ASN:HD21	2:D:77:GLU:H	1.32	0.76
1:A:35:ARG:HG2	1:A:35:ARG:HH11	1.49	0.76
2:B:42:ASN:HD21	2:B:77:GLU:H	1.34	0.75
1:C:211:ALA:HB2	1:C:241:PHE:CE1	2.21	0.74
2:B:4:THR:HG23	2:B:86:THR:OG1	1.88	0.74
1:A:21:ARG:HH11	1:A:21:ARG:HG3	1.52	0.74
1:A:79:ARG:NH2	4:A:275:SO4:O2	2.21	0.73
1:C:230:LEU:HD12	1:C:243:LYS:HE3	1.69	0.73
2:D:73:THR:HG22	2:D:75:LYS:H	1.53	0.73
1:C:65:ARG:HB3	1:C:65:ARG:HH11	1.53	0.73
1:A:121:LYS:HG3	1:A:122:ASP:N	2.03	0.73
2:D:73:THR:HG22	2:D:75:LYS:N	2.04	0.73
2:D:4:THR:HG22	2:D:5:PRO:HD2	1.71	0.73
2:D:9:VAL:HG23	2:D:93:VAL:CG2	2.19	0.72
1:A:194:ILE:HD12	1:A:194:ILE:N	2.05	0.72
1:C:193:PRO:HA	1:C:199:ALA:CB	2.20	0.71
1:A:189:VAL:HG12	1:A:190:THR:N	2.05	0.71
2:D:9:VAL:CG2	2:D:93:VAL:HG22	2.20	0.70
1:C:13:SER:HA	1:C:20:PRO:HB3	1.73	0.70
1:A:109:PHE:CD1	1:A:110:LEU:N	2.60	0.70
1:A:204:TRP:HE3	1:A:206:LEU:HD21	1.56	0.70
2:B:73:THR:CG2	2:B:75:LYS:H	2.05	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:230:LEU:CD1	1:C:243:LYS:HE3	2.22	0.69
1:A:211:ALA:HB2	1:A:241:PHE:CE1	2.27	0.69
2:D:51:HIS:HA	2:D:65:LEU:O	1.92	0.69
1:A:202:ARG:HH11	1:A:202:ARG:HG3	1.57	0.69
1:C:50:PRO:HA	1:C:53:GLU:OE2	1.93	0.69
1:A:137:ASP:H	1:A:140:ALA:HB3	1.58	0.69
1:A:220:ASP:N	1:A:256:ALA:O	2.25	0.69
1:A:55:GLU:HA	1:A:55:GLU:OE1	1.93	0.69
1:A:218:GLN:OE1	1:A:223:GLY:HA3	1.93	0.68
3:P:5:ARG:HH21	3:Q:5:ARG:CD	2.06	0.68
1:A:177:GLU:CD	1:A:177:GLU:H	1.96	0.68
1:A:50:PRO:HA	1:A:53:GLU:OE2	1.94	0.68
1:A:65:ARG:NH1	1:A:65:ARG:CB	2.47	0.68
1:C:44:ARG:HA	1:C:64:THR:HG23	1.75	0.68
1:C:89:GLU:O	1:C:89:GLU:HG2	1.93	0.68
1:C:194:ILE:HD12	1:C:194:ILE:N	2.08	0.68
1:A:193:PRO:HA	1:A:199:ALA:CB	2.23	0.67
1:C:152:GLU:OE1	3:P:5:ARG:NH1	2.23	0.67
1:C:109:PHE:CE2	1:C:161:GLU:HG2	2.30	0.67
2:D:39:LEU:HB3	2:D:46:ILE:HD12	1.77	0.67
1:C:202:ARG:HH11	1:C:202:ARG:HG3	1.61	0.66
1:C:211:ALA:HB2	1:C:241:PHE:CD1	2.30	0.66
3:P:5:ARG:NH2	3:Q:5:ARG:CD	2.59	0.66
2:B:9:VAL:CG2	2:B:93:VAL:CG2	2.74	0.66
1:C:236:ALA:O	2:D:12:ARG:HD3	1.95	0.66
1:C:12:VAL:HG22	1:C:94:THR:HG23	1.77	0.65
2:B:49:VAL:HG12	2:B:50:GLU:N	2.11	0.65
2:B:9:VAL:HG21	2:B:93:VAL:HG22	1.78	0.65
1:A:255:GLN:N	1:A:255:GLN:NE2	2.43	0.65
1:C:234:ARG:NH1	2:D:8:GLN:OE1	2.30	0.65
1:A:79:ARG:HH11	1:C:72:GLN:NE2	1.94	0.65
3:P:5:ARG:NH2	3:Q:5:ARG:HD3	2.12	0.65
2:D:29:GLY:HA2	2:D:61:SER:OG	1.97	0.65
1:C:117:ALA:HB2	2:D:60:TRP:CD2	2.33	0.64
2:D:4:THR:CG2	2:D:5:PRO:HD2	2.27	0.64
1:A:211:ALA:HB2	1:A:241:PHE:CD1	2.33	0.64
1:C:28:VAL:O	1:C:29:ASP:O	2.15	0.63
1:A:271:THR:C	1:A:272:LEU:HD23	2.19	0.63
1:C:82:ARG:HE	1:C:89:GLU:HB2	1.63	0.63
1:C:103:LEU:HA	1:C:108:ARG:O	1.98	0.63
1:C:249:VAL:HG22	1:C:257:TYR:CE1	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:PHE:HB3	2:B:56:PHE:CE2	2.34	0.63
1:A:172:LEU:HD23	1:A:179:LEU:HD23	1.81	0.63
2:D:16:GLU:HB3	2:D:19:LYS:CB	2.29	0.63
1:A:249:VAL:HG22	1:A:257:TYR:CE1	2.34	0.62
2:D:73:THR:CG2	2:D:75:LYS:H	2.11	0.62
2:D:42:ASN:HD22	2:D:77:GLU:HB2	1.65	0.62
1:A:189:VAL:CG1	1:A:190:THR:N	2.62	0.62
1:C:204:TRP:CE3	1:C:206:LEU:HD21	2.34	0.62
1:A:109:PHE:CE2	1:A:161:GLU:HG2	2.35	0.62
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.82	0.62
1:C:263:HIS:HD2	1:C:265:GLY:N	1.90	0.61
1:C:21:ARG:CG	1:C:21:ARG:HH11	2.11	0.61
1:A:220:ASP:OD2	1:A:220:ASP:O	2.17	0.61
1:C:35:ARG:HH11	1:C:46:VAL:CG2	2.14	0.61
1:C:194:ILE:HD13	1:C:199:ALA:HA	1.82	0.61
1:A:255:GLN:CD	1:A:255:GLN:N	2.53	0.61
1:A:6:LYS:HD2	1:A:113:TYR:OH	2.01	0.60
1:C:32:GLN:NE2	2:D:53:ASP:OD2	2.32	0.60
1:A:243:LYS:HG2	1:A:244:TRP:N	2.16	0.60
1:C:65:ARG:CB	1:C:65:ARG:NH1	2.46	0.60
1:A:184:PRO:HB3	1:A:265:GLY:O	2.01	0.60
1:C:127:ASN:ND2	1:C:132:SER:OG	2.33	0.60
2:B:42:ASN:HD22	2:B:77:GLU:HB2	1.67	0.59
2:B:51:HIS:HA	2:B:65:LEU:O	2.03	0.59
1:A:150:ALA:O	1:A:151:SER:CB	2.50	0.59
1:C:150:ALA:O	1:C:151:SER:CB	2.48	0.59
1:C:190:THR:O	1:C:201:LEU:HA	2.02	0.59
1:C:109:PHE:CD1	1:C:110:LEU:N	2.71	0.59
1:C:229:GLU:OE2	1:C:244:TRP:HH2	1.86	0.58
1:C:209:TYR:HA	1:C:210:PRO:O	2.03	0.58
1:C:138:THR:HA	1:C:141:GLN:HE21	1.68	0.58
2:D:16:GLU:O	2:D:17:ASN:C	2.42	0.58
3:P:5:ARG:HH21	3:Q:5:ARG:HD3	1.68	0.58
1:C:24:SER:HB3	1:C:36:PHE:HB3	1.86	0.57
2:D:42:ASN:ND2	2:D:77:GLU:HB2	2.20	0.57
1:C:121:LYS:HG3	1:C:122:ASP:N	2.19	0.57
1:C:35:ARG:HH11	1:C:46:VAL:HG21	1.67	0.57
1:A:65:ARG:NH2	1:C:145:GLN:NE2	2.45	0.57
1:C:128:GLU:O	1:C:130:LEU:HG	2.03	0.57
1:A:194:ILE:HD13	1:A:199:ALA:HA	1.86	0.57
1:A:47:PRO:HB3	1:A:60:TRP:CZ2	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:49:VAL:CG1	2:B:50:GLU:N	2.68	0.56
1:C:47:PRO:HB3	1:C:60:TRP:CZ2	2.41	0.56
2:D:98:ASP:C	2:D:99:MET:HG3	2.24	0.56
1:C:82:ARG:NE	1:C:89:GLU:HB2	2.19	0.56
1:A:194:ILE:CD1	1:A:194:ILE:N	2.69	0.56
1:C:44:ARG:NH2	1:C:61:ASP:OD1	2.39	0.55
1:A:13:SER:HA	1:A:20:PRO:HB3	1.89	0.55
1:C:221:GLY:C	1:C:222:GLU:HG3	2.27	0.55
1:C:137:ASP:O	1:C:141:GLN:HG3	2.06	0.55
1:A:204:TRP:CE3	1:A:206:LEU:HD21	2.41	0.55
2:D:96:ASP:O	2:D:98:ASP:N	2.39	0.55
1:A:79:ARG:NH1	1:C:72:GLN:NE2	2.53	0.55
1:C:194:ILE:CD1	1:C:194:ILE:N	2.70	0.54
1:C:166:GLU:HG2	1:C:166:GLU:O	2.06	0.54
2:B:9:VAL:HG23	2:B:93:VAL:CG2	2.36	0.54
1:A:270:VAL:HG12	1:A:272:LEU:CD2	2.37	0.54
1:C:254:GLU:HG2	1:C:274:TRP:CE3	2.43	0.54
1:A:135:ALA:HB1	1:A:140:ALA:CB	2.38	0.54
1:A:155:HIS:CE1	1:C:155:HIS:CE1	2.95	0.54
2:D:16:GLU:O	2:D:19:LYS:N	2.27	0.53
1:A:193:PRO:HA	1:A:199:ALA:HA	1.88	0.53
1:C:209:TYR:HA	1:C:210:PRO:C	2.27	0.53
1:A:103:LEU:HA	1:A:108:ARG:O	2.07	0.53
1:A:182:LEU:HD23	1:A:210:PRO:HD3	1.90	0.53
2:B:4:THR:HG22	2:B:5:PRO:HD2	1.89	0.53
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.43	0.53
1:A:212:GLU:OE1	1:A:212:GLU:HA	2.08	0.53
1:C:219:GLN:CG	1:C:220:ASP:N	2.46	0.53
1:A:193:PRO:HA	1:A:199:ALA:CA	2.39	0.53
1:C:202:ARG:NH1	1:C:202:ARG:HG3	2.22	0.53
2:D:16:GLU:HG2	2:D:19:LYS:CB	2.39	0.53
1:A:250:PRO:HG2	1:A:253:GLU:OE2	2.09	0.52
1:C:234:ARG:HG3	2:D:10:TYR:CZ	2.44	0.52
1:C:51:TRP:HB2	1:C:174:LYS:O	2.09	0.52
1:A:145:GLN:CD	1:C:65:ARG:HH21	2.11	0.52
1:A:194:ILE:CD1	1:A:199:ALA:HA	2.40	0.52
1:A:204:TRP:HZ2	2:B:98:ASP:O	1.92	0.52
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.92	0.52
2:B:4:THR:CG2	2:B:5:PRO:HD2	2.40	0.52
1:A:48:ARG:NH1	2:B:53:ASP:OD2	2.42	0.52
2:D:40:LEU:HD11	2:D:81:ARG:HB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:ILE:HG22	1:A:24:SER:N	2.25	0.52
1:A:182:LEU:HD22	1:A:264:GLU:HB3	1.92	0.51
1:C:14:ARG:HD2	1:C:17:ARG:NH1	2.25	0.51
1:A:222:GLU:O	1:A:223:GLY:C	2.47	0.51
2:D:9:VAL:HG21	2:D:93:VAL:CG2	2.37	0.51
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.45	0.51
2:B:29:GLY:HA2	2:B:61:SER:OG	2.10	0.51
2:B:17:ASN:OD1	2:B:73:THR:C	2.48	0.51
2:B:73:THR:CG2	2:B:76:ASP:H	2.23	0.51
1:C:44:ARG:HA	1:C:64:THR:CG2	2.41	0.51
2:D:9:VAL:CG2	2:D:93:VAL:HG23	2.41	0.51
3:P:5:ARG:NH2	3:Q:5:ARG:HE	2.07	0.50
1:A:152:GLU:OE2	3:Q:5:ARG:NH1	2.44	0.50
2:D:25:CYS:HB2	2:D:39:LEU:HD21	1.93	0.50
1:C:193:PRO:HA	1:C:199:ALA:HA	1.93	0.50
1:A:243:LYS:HG2	1:A:244:TRP:H	1.74	0.50
1:C:215:LEU:CD2	1:C:261:VAL:HG22	2.42	0.50
1:C:55:GLU:HA	1:C:55:GLU:OE1	2.12	0.50
1:C:23:ILE:CG2	1:C:24:SER:N	2.75	0.50
1:C:191:HIS:ND1	1:C:274:TRP:CZ3	2.76	0.50
1:C:35:ARG:HH12	1:C:46:VAL:HG21	1.76	0.49
1:A:51:TRP:HB2	1:A:174:LYS:O	2.12	0.49
2:D:29:GLY:HA2	2:D:61:SER:CB	2.42	0.49
1:C:138:THR:HA	1:C:141:GLN:HG3	1.93	0.49
1:A:202:ARG:HG3	1:A:202:ARG:NH1	2.25	0.49
2:B:73:THR:HG22	2:B:76:ASP:H	1.76	0.49
2:D:9:VAL:HG21	2:D:93:VAL:O	2.13	0.49
1:A:79:ARG:HD2	1:C:72:GLN:NE2	2.27	0.49
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.47	0.49
1:A:82:ARG:HG3	1:A:87:GLN:HB2	1.94	0.49
1:A:263:HIS:HD2	1:A:265:GLY:N	1.92	0.49
2:D:39:LEU:CB	2:D:46:ILE:HD12	2.42	0.49
1:A:23:ILE:CG2	1:A:24:SER:N	2.75	0.49
2:D:9:VAL:HA	2:D:24:ASN:O	2.13	0.49
1:C:73:ILE:N	1:C:73:ILE:HD13	2.26	0.49
1:C:226:GLN:O	1:C:227:ASP:O	2.31	0.49
1:A:187:THR:O	1:A:188:HIS:HB3	2.13	0.49
1:C:268:GLU:CB	1:C:269:PRO:CD	2.89	0.49
1:A:25:VAL:HG23	1:A:25:VAL:O	2.11	0.49
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.47	0.49
1:A:37:ASP:C	1:A:39:ASP:H	2.16	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:244:TRP:C	1:C:244:TRP:HE3	2.15	0.49
1:C:218:GLN:HE21	1:C:260:HIS:CE1	2.31	0.49
1:C:219:GLN:NE2	1:C:220:ASP:OD2	2.44	0.48
1:A:254:GLU:N	1:A:254:GLU:OE1	2.42	0.48
1:A:202:ARG:HD3	1:A:244:TRP:CD2	2.49	0.48
1:A:127:ASN:ND2	1:A:132:SER:OG	2.44	0.48
1:C:81:LEU:HD21	1:C:123:TYR:CZ	2.49	0.48
1:A:138:THR:CA	1:A:141:GLN:HG3	2.36	0.48
1:C:193:PRO:HA	1:C:199:ALA:CA	2.43	0.48
1:A:221:GLY:O	1:A:223:GLY:N	2.46	0.48
2:D:73:THR:HG22	2:D:76:ASP:H	1.77	0.48
1:A:85:TYR:O	1:A:86:ASN:C	2.50	0.48
1:A:76:VAL:HG11	3:P:8:LEU:HD11	1.96	0.48
1:C:35:ARG:HH11	1:C:35:ARG:CG	2.27	0.48
1:C:163:THR:O	1:C:167:TRP:HD1	1.96	0.48
1:A:195:SER:OG	1:A:198:GLU:HG3	2.14	0.48
1:A:11:SER:HB2	1:A:74:PHE:HD2	1.79	0.48
1:A:72:GLN:HE21	1:C:76:VAL:HG13	1.78	0.48
1:A:51:TRP:CZ3	1:A:52:MET:SD	3.07	0.47
1:C:152:GLU:CD	3:P:5:ARG:HH12	2.11	0.47
1:A:109:PHE:CE2	1:A:161:GLU:HA	2.50	0.47
1:A:79:ARG:HH11	1:C:72:GLN:HE21	1.61	0.47
1:C:224:HIS:HB2	1:C:225:THR:H	1.25	0.47
1:C:11:SER:HB2	1:C:74:PHE:HD2	1.79	0.47
2:D:37:VAL:HG13	2:D:82:VAL:HG22	1.95	0.47
2:B:57:SER:O	2:B:60:TRP:N	2.47	0.47
1:C:218:GLN:CG	1:C:260:HIS:CE1	2.97	0.47
1:C:228:THR:HA	1:C:246:ALA:O	2.14	0.47
1:A:210:PRO:O	1:A:263:HIS:HE1	1.97	0.47
1:C:162:ASP:O	1:C:163:THR:C	2.52	0.47
1:A:47:PRO:HB3	1:A:60:TRP:CH2	2.50	0.47
1:A:73:ILE:HG21	3:P:6:THR:HG23	1.96	0.47
1:A:70:THR:O	1:A:71:ALA:C	2.52	0.47
2:B:42:ASN:ND2	2:B:77:GLU:HB2	2.29	0.47
1:C:82:ARG:CG	1:C:87:GLN:O	2.62	0.47
1:C:162:ASP:O	1:C:165:VAL:N	2.48	0.47
1:A:216:THR:O	1:A:259:CYS:HA	2.15	0.47
2:D:38:ASP:O	2:D:80:CYS:HA	2.15	0.46
2:D:49:VAL:CG1	2:D:50:GLU:N	2.77	0.46
3:P:5:ARG:HH21	3:Q:5:ARG:NE	2.08	0.46
1:C:127:ASN:O	1:C:130:LEU:N	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:212:GLU:OE1	1:C:212:GLU:HA	2.15	0.46
1:A:28:VAL:O	1:A:29:ASP:O	2.34	0.46
1:A:255:GLN:H	1:A:255:GLN:NE2	2.10	0.46
1:C:109:PHE:CE2	1:C:161:GLU:HA	2.50	0.46
2:B:31:HIS:ND1	2:B:32:PRO:HA	2.31	0.46
2:B:16:GLU:O	2:B:17:ASN:C	2.53	0.46
1:A:177:GLU:N	1:A:177:GLU:CD	2.64	0.46
1:C:23:ILE:HG22	1:C:24:SER:N	2.31	0.46
1:A:218:GLN:CG	1:A:260:HIS:CE1	2.98	0.46
1:C:55:GLU:CD	1:C:174:LYS:NZ	2.69	0.46
1:C:189:VAL:HG12	1:C:190:THR:N	2.30	0.46
2:D:49:VAL:HG12	2:D:50:GLU:N	2.30	0.46
1:C:104:GLY:HA3	1:C:106:ASP:OD1	2.16	0.46
1:A:155:HIS:CE1	1:C:155:HIS:ND1	2.83	0.46
1:A:129:ASP:OD2	1:A:131:ARG:HG3	2.16	0.46
1:A:104:GLY:C	1:A:106:ASP:N	2.69	0.45
2:B:73:THR:HB	2:B:76:ASP:HB2	1.98	0.45
1:A:198:GLU:HB3	1:A:250:PRO:HA	1.98	0.45
1:A:103:LEU:HG	1:A:168:LEU:HD23	1.96	0.45
1:A:37:ASP:C	1:A:39:ASP:N	2.69	0.45
2:B:16:GLU:HG2	2:B:19:LYS:CB	2.47	0.45
1:A:202:ARG:HD3	1:A:244:TRP:CE2	2.51	0.45
1:C:250:PRO:O	1:C:251:SER:C	2.55	0.45
1:A:135:ALA:HB1	1:A:140:ALA:HB3	1.97	0.45
1:A:198:GLU:CB	1:A:249:VAL:O	2.64	0.45
1:A:232:GLU:O	1:A:233:THR:C	2.54	0.45
1:A:121:LYS:HE2	2:B:1(A):ILE:HD11	1.99	0.45
2:B:98:ASP:C	2:B:99:MET:HG3	2.36	0.45
1:C:115:GLN:CG	1:C:125:THR:HG23	2.34	0.45
2:D:51:HIS:HB3	2:D:66:TYR:CD2	2.52	0.45
1:C:180:LEU:HA	1:C:180:LEU:HD23	1.66	0.45
1:C:244:TRP:C	1:C:244:TRP:CE3	2.90	0.45
1:A:43:PRO:O	1:A:68:ARG:NH2	2.50	0.45
1:C:14:ARG:CD	1:C:17:ARG:NH1	2.80	0.44
2:D:56:PHE:HA	2:D:62:PHE:HA	1.99	0.44
1:C:21:ARG:HD3	1:C:38:ASN:OD1	2.17	0.44
1:A:79:ARG:NH1	1:C:72:GLN:HE22	2.15	0.44
1:C:198:GLU:HB3	1:C:250:PRO:HA	2.00	0.44
2:D:79:ALA:HB2	2:D:94:LYS:HA	1.98	0.44
2:B:9:VAL:HG21	2:B:93:VAL:CG2	2.44	0.44
1:C:230:LEU:HD13	1:C:230:LEU:HA	1.73	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:GLN:CB	1:A:223:GLY:CA	2.83	0.44
1:A:270:VAL:HG12	1:A:272:LEU:HD21	1.98	0.44
1:C:253:GLU:O	1:C:255:GLN:N	2.51	0.44
1:A:162:ASP:HB2	1:A:163:THR:H	1.40	0.44
2:B:96:ASP:OD1	2:B:98:ASP:HB2	2.18	0.44
1:A:228:THR:HG22	1:A:230:LEU:HD22	1.99	0.44
1:A:73:ILE:HD11	3:Q:8:LEU:HD22	1.99	0.44
1:A:166:GLU:O	1:A:169:HIS:HB2	2.18	0.44
1:A:191:HIS:ND1	1:A:274:TRP:CH2	2.83	0.43
1:A:81:LEU:HD21	1:A:123:TYR:CZ	2.53	0.43
2:B:89:GLN:O	2:B:90:PRO:C	2.55	0.43
1:C:53:GLU:C	1:C:55:GLU:H	2.21	0.43
1:C:82:ARG:HG3	1:C:87:GLN:O	2.18	0.43
1:C:29:ASP:O	1:C:31:THR:N	2.50	0.43
2:D:9:VAL:HG21	2:D:93:VAL:HG22	1.94	0.43
2:B:17:ASN:OD1	2:B:73:THR:CA	2.66	0.43
1:C:194:ILE:H	1:C:199:ALA:HA	1.82	0.43
1:C:233:THR:HG22	1:C:234:ARG:N	2.34	0.43
1:A:155:HIS:HE1	1:C:155:HIS:ND1	2.17	0.43
2:B:4:THR:HG22	2:B:5:PRO:CD	2.49	0.43
2:B:39:LEU:HA	2:B:39:LEU:HD23	1.62	0.43
2:D:15:ALA:HB2	2:D:95:TRP:CZ2	2.54	0.43
1:C:33:PHE:CD1	1:C:33:PHE:N	2.86	0.43
1:A:145:GLN:OE1	1:C:65:ARG:NH2	2.46	0.43
1:A:272:LEU:N	1:A:272:LEU:HD23	2.34	0.43
1:C:166:GLU:OE2	1:C:170:LYS:HE3	2.19	0.43
1:A:89:GLU:HG2	1:A:89:GLU:O	2.19	0.43
1:C:243:LYS:HG2	1:C:244:TRP:N	2.34	0.43
1:C:128:GLU:HG2	1:C:128:GLU:H	1.58	0.43
1:A:44:ARG:HA	1:A:64:THR:HG23	2.01	0.43
3:P:5:ARG:CZ	3:Q:5:ARG:NE	2.82	0.42
1:C:103:LEU:HG	1:C:168:LEU:HD23	2.01	0.42
1:A:129:ASP:O	1:A:130:LEU:HB2	2.19	0.42
1:A:109:PHE:CD2	1:A:161:GLU:HA	2.54	0.42
2:B:6:LYS:HE3	2:B:6:LYS:HB2	1.89	0.42
1:C:33:PHE:C	1:C:48:ARG:HB2	2.40	0.42
1:C:194:ILE:HD13	1:C:199:ALA:CA	2.46	0.42
1:A:255:GLN:C	1:A:257:TYR:N	2.72	0.42
1:C:117:ALA:HB2	2:D:60:TRP:CZ2	2.54	0.42
1:C:249:VAL:HA	1:C:250:PRO:HD3	1.69	0.42
1:C:162:ASP:O	1:C:164:CYS:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:LYS:N	1:A:186:LYS:HD3	2.33	0.42
1:C:6:LYS:HD2	1:C:113:TYR:OH	2.20	0.42
1:A:249:VAL:HA	1:A:250:PRO:HD3	1.77	0.42
1:A:224:HIS:C	1:A:225:THR:HG23	2.40	0.42
2:D:6:LYS:O	2:D:27:VAL:HA	2.19	0.42
1:C:229:GLU:OE2	1:C:244:TRP:CH2	2.70	0.42
1:C:194:ILE:HG22	1:C:195:SER:N	2.35	0.42
1:C:218:GLN:NE2	1:C:260:HIS:CE1	2.88	0.42
1:C:209:TYR:CA	1:C:210:PRO:O	2.67	0.42
1:C:191:HIS:ND1	1:C:274:TRP:CH2	2.77	0.42
1:A:21:ARG:HH11	1:A:21:ARG:CG	2.24	0.42
1:C:39:ASP:O	1:C:40:ALA:O	2.38	0.42
1:A:226:GLN:O	1:A:227:ASP:C	2.56	0.42
1:C:43:PRO:O	1:C:68:ARG:NH2	2.53	0.41
1:C:97:TRP:HD1	1:C:115:GLN:O	2.02	0.41
1:C:225:THR:CB	1:C:247:VAL:HG21	2.50	0.41
2:B:16:GLU:HB3	2:B:19:LYS:CB	2.50	0.41
1:A:24:SER:HB3	1:A:36:PHE:HB3	2.01	0.41
1:C:182:LEU:HD23	1:C:182:LEU:HA	1.81	0.41
1:C:230:LEU:HD11	1:C:243:LYS:HE3	1.99	0.41
1:A:122:ASP:O	1:A:136:VAL:HG11	2.21	0.41
1:C:193:PRO:HA	1:C:199:ALA:HB1	1.99	0.41
1:A:45:MET:O	1:A:60:TRP:CE3	2.73	0.41
1:A:33:PHE:CD1	1:A:33:PHE:N	2.89	0.41
2:B:71:THR:HA	2:B:72:PRO:HD2	1.69	0.41
1:A:79:ARG:O	1:A:80:THR:C	2.57	0.41
1:C:48:ARG:HD2	1:C:48:ARG:HA	1.47	0.41
1:A:191:HIS:CE1	1:A:254:GLU:OE2	2.74	0.41
1:A:143:SER:O	1:A:147:SER:HB2	2.20	0.41
2:D:71:THR:HA	2:D:72:PRO:HD2	1.67	0.41
1:C:35:ARG:NH1	1:C:46:VAL:CG2	2.70	0.41
2:D:9:VAL:HG23	2:D:93:VAL:HG23	1.99	0.41
1:C:231:VAL:HG13	1:C:244:TRP:CZ3	2.56	0.41
1:C:194:ILE:CD1	1:C:199:ALA:HA	2.49	0.41
1:C:14:ARG:HD3	1:C:17:ARG:HH12	1.86	0.41
1:C:124:LEU:HA	1:C:124:LEU:HD12	1.83	0.41
2:B:29:GLY:HA2	2:B:61:SER:CB	2.51	0.41
2:D:39:LEU:HD13	2:D:68:THR:HG22	2.02	0.40
2:B:23:LEU:HD12	2:B:23:LEU:HA	1.79	0.40
2:D:70:PHE:C	2:D:70:PHE:CD1	2.93	0.40
2:D:98:ASP:O	2:D:99:MET:CG	2.49	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:GLN:HB3	2:D:60:TRP:CH2	2.55	0.40
2:B:49:VAL:CG1	2:B:50:GLU:H	2.34	0.40
2:D:3:ARG:HG2	2:D:3:ARG:HH11	1.86	0.40
1:C:117:ALA:HA	1:C:121:LYS:O	2.21	0.40
1:C:231:VAL:HG11	1:C:244:TRP:CE2	2.56	0.40
2:D:4:THR:HG22	2:D:5:PRO:CD	2.45	0.40
1:A:55:GLU:CD	1:A:174:LYS:NZ	2.75	0.40
1:C:109:PHE:CD2	1:C:161:GLU:HA	2.56	0.40
1:C:192:HIS:O	1:C:200:THR:N	2.50	0.40
2:B:9:VAL:O	2:B:10:TYR:HB3	2.21	0.40
2:B:16:GLU:O	2:B:19:LYS:N	2.46	0.40
1:C:33:PHE:HA	1:C:49:ALA:HB2	2.03	0.40
1:C:231:VAL:HG11	1:C:244:TRP:CZ2	2.57	0.40
1:A:161:GLU:O	1:A:165:VAL:CG1	2.70	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	A	272/274 (99%)	231 (85%)	31 (11%)	10 (4%)	4	14
1	C	272/274 (99%)	230 (85%)	24 (9%)	18 (7%)	1	4
2	B	98/100 (98%)	92 (94%)	5 (5%)	1 (1%)	19	52
2	D	98/100 (98%)	89 (91%)	7 (7%)	2 (2%)	9	30
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%)	656 (87%)	67 (9%)	31 (4%)	3	11

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	29	ASP
1	A	40	ALA
1	A	177	GLU
1	C	15	PRO
1	C	29	ASP
1	C	40	ALA
1	C	224	HIS
1	C	225	THR
1	C	227	ASP
2	D	97	ARG
1	A	15	PRO
1	A	220	ASP
1	A	222	GLU
1	C	30	ASP
1	C	178	THR
1	C	223	GLY
1	A	2	SER
1	C	57	SER
1	C	163	THR
1	C	177	GLU
1	C	254	GLU
1	A	30	ASP
1	A	254	GLU
1	C	220	ASP
2	D	17	ASN
1	A	57	SER
2	B	97	ARG
1	C	219	GLN
1	C	251	SER
1	C	2	SER
1	C	194	ILE

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	230/234 (98%)	203 (88%)	27 (12%)	7 20

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	229/234 (98%)	200 (87%)	29 (13%)	5	16
2	B	93/95 (98%)	86 (92%)	7 (8%)	17	43
2	D	93/95 (98%)	84 (90%)	9 (10%)	10	29
3	P	8/8 (100%)	6 (75%)	2 (25%)	1	2
3	Q	8/8 (100%)	7 (88%)	1 (12%)	6	17
All	All	661/674 (98%)	586 (89%)	75 (11%)	7	22

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	21	ARG
1	A	29	ASP
1	A	35	ARG
1	A	48	ARG
1	A	53	GLU
1	A	62	ARG
1	A	137	ASP
1	A	138	THR
1	A	145	GLN
1	A	147	SER
1	A	151	SER
1	A	155	HIS
1	A	165	VAL
1	A	178	THR
1	A	179	LEU
1	A	186	LYS
1	A	187	THR
1	A	198	GLU
1	A	227	ASP
1	A	229	GLU
1	A	244	TRP
1	A	251	SER
1	A	258	THR
1	A	272	LEU
1	A	273	ARG
2	B	4	THR
2	B	9	VAL
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	21	ARG
1	C	29	ASP
1	C	35	ARG
1	C	48	ARG
1	C	53	GLU
1	C	65	ARG
1	C	89	GLU
1	C	121	LYS
1	C	137	ASP
1	C	138	THR
1	C	145	GLN
1	C	147	SER
1	C	151	SER
1	C	155	HIS
1	C	186	LYS
1	C	187	THR
1	C	198	GLU
1	C	219	GLN
1	C	220	ASP
1	C	228	THR
1	C	230	LEU
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	12	ARG
2	D	14	PRO
2	D	45	ARG
2	D	48	LYS
2	D	58	LYS
2	D	70	PHE
2	D	71	THR
2	D	73	THR

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Mol	Chain	Res	Type
3	P	1	VAL
3	P	5	ARG
3	Q	1	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) 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	148	ASN
1	A	181	HIS
1	A	224	HIS
1	A	255	GLN
1	A	263	HIS
2	B	42	ASN
2	B	89	GLN
1	C	72	GLN
1	C	87	GLN
1	C	115	GLN
1	C	127	ASN
1	C	141	GLN
1	C	145	GLN
1	C	219	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

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	0.45	0	6,6,6	0.31	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SO4	A	275	-	-	0/0/0/0	0/0/0/0

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

There are no such residues in this entry.

5.8 Polymer linkage issues

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 will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.