



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:51 AM UTC

PDB ID : 1KPR / pdb_00001kpr
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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	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

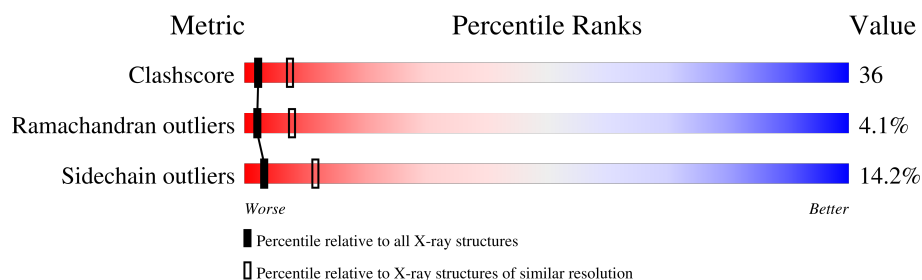
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (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 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	34% 43% 18% 5%
1	C	274	34% 42% 20% .
2	B	100	43% 44% 11% .
2	D	100	32% 55% 12% .
3	P	9	44% 33% 22%
3	Q	9	44% 33% 22%

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 mutation	UNP P13747
A	256	ALA	ARG	engineered mutation	UNP P13747
C	107	GLY	ARG	engineered mutation	UNP P13747
C	256	ALA	ARG	engineered mutation	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			

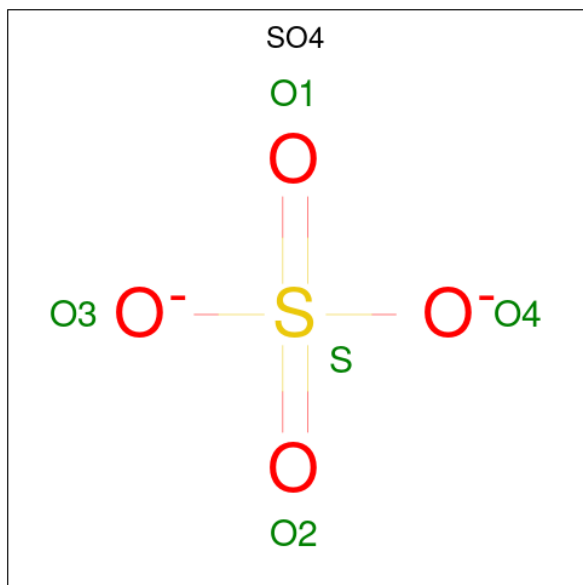
There are 2 discrepancies between the modelled and reference sequences:

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 (CCD ID: SO4) (formula: O₄S).



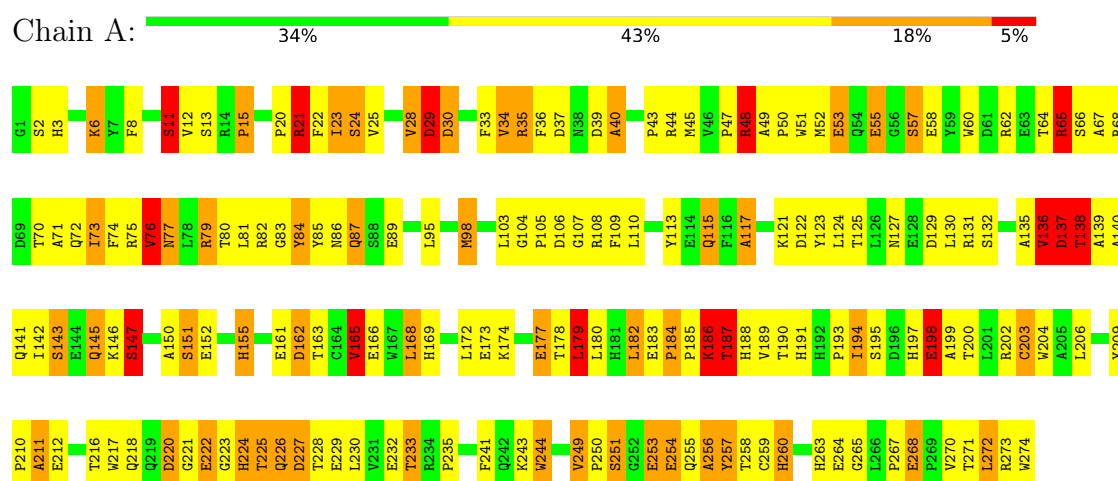
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

3 Residue-property plots

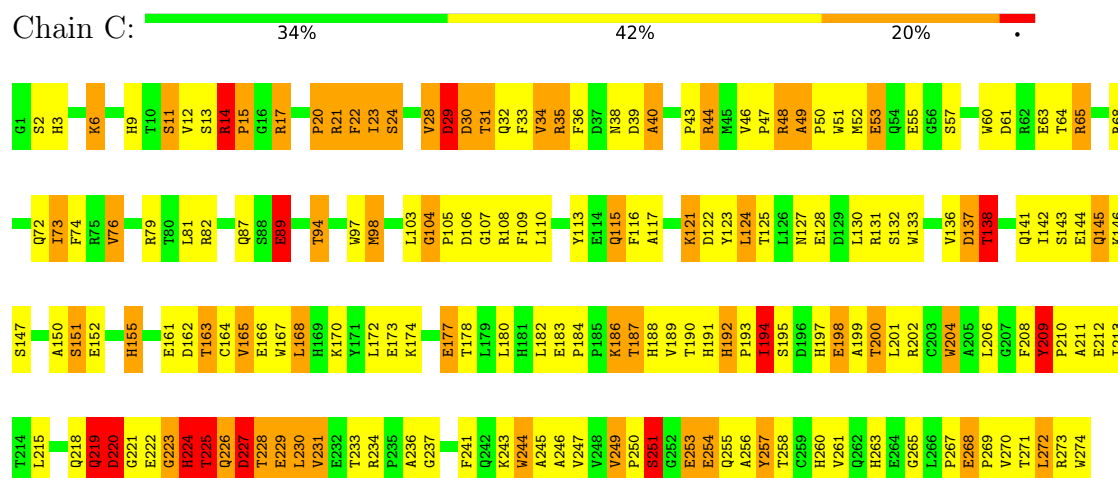
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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



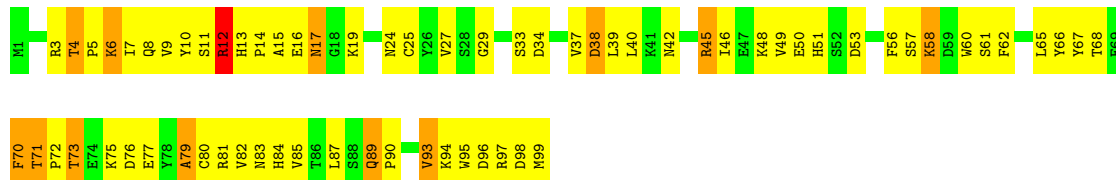
• Molecule 2: BETA-2-MICROGLOBULIN





• Molecule 2: BETA-2-MICROGLOBULIN

Chain D: 32% 55% 12%



• Molecule 3: Peptide VMAPRTVLL

Chain P: 44% 33% 22%



• Molecule 3: Peptide VMAPRTVLL

Chain Q: 44% 33% 22%



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	178.70Å 178.70Å 88.20Å 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

5.1 Standard geometry

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	1.61	34/2272 (1.5%)	1.73	59/3092 (1.9%)
1	C	1.65	29/2271 (1.3%)	1.77	62/3089 (2.0%)
2	B	1.82	16/851 (1.9%)	1.68	12/1152 (1.0%)
2	D	1.40	3/851 (0.4%)	1.63	17/1152 (1.5%)
3	P	1.82	1/69 (1.4%)	1.60	1/92 (1.1%)
3	Q	1.52	0/69	1.69	3/92 (3.3%)
All	All	1.63	83/6383 (1.3%)	1.72	154/8669 (1.8%)

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 (83) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	MET	SD-CE	11.38	2.08	1.79
1	C	256	ALA	CA-CB	10.41	1.72	1.53
2	B	10	TYR	C-O	-9.53	1.12	1.23
1	C	224	HIS	CA-CB	9.39	1.69	1.53
1	C	29	ASP	CA-C	-8.62	1.41	1.52
1	C	224	HIS	CE1-NE2	8.36	1.41	1.32
1	C	224	HIS	ND1-CE1	8.06	1.40	1.32
1	A	136	VAL	CA-CB	-7.68	1.44	1.55
1	A	142	ILE	CA-CB	-7.63	1.45	1.54
2	D	7	ILE	CA-CB	-7.41	1.44	1.54
1	A	224	HIS	CA-C	-7.38	1.42	1.53
2	B	27	VAL	CA-CB	-7.23	1.45	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	137	ASP	CA-C	-7.08	1.46	1.53
2	B	7	ILE	C-O	-7.07	1.16	1.24
1	A	139	ALA	CA-CB	-7.00	1.42	1.53
1	C	270	VAL	CA-CB	-7.00	1.46	1.54
1	C	224	HIS	N-CA	6.96	1.55	1.46
1	C	142	ILE	CA-CB	-6.74	1.46	1.54
1	C	224	HIS	CD2-NE2	6.73	1.45	1.37
1	A	117	ALA	CA-C	-6.31	1.45	1.52
1	A	12	VAL	C-O	-6.31	1.17	1.24
1	C	183	GLU	CA-CB	-6.29	1.45	1.53
1	A	147	SER	CA-C	-6.28	1.45	1.52
1	C	98	MET	CG-SD	6.26	1.96	1.80
1	C	34	VAL	CB-CG1	6.20	1.73	1.52
1	C	224	HIS	CG-CD2	6.17	1.42	1.35
2	B	35	ILE	C-O	-6.14	1.16	1.23
1	A	98	MET	SD-CE	6.12	1.94	1.79
2	B	11	SER	N-CA	6.06	1.52	1.45
1	A	115	GLN	CA-CB	-6.06	1.43	1.53
2	B	54	LEU	C-O	-6.06	1.16	1.23
1	A	249	VAL	CA-C	6.04	1.57	1.52
1	C	52	MET	SD-CE	-5.93	1.64	1.79
1	A	117	ALA	CA-CB	-5.91	1.41	1.54
1	A	65	ARG	CA-C	5.91	1.60	1.52
2	B	26	TYR	CA-CB	-5.90	1.45	1.53
1	A	147	SER	C-O	-5.84	1.17	1.24
1	C	186	LYS	C-O	-5.83	1.17	1.24
1	C	224	HIS	CG-ND1	5.81	1.44	1.38
1	A	115	GLN	CA-C	-5.77	1.45	1.52
1	C	49	ALA	N-CA	-5.74	1.39	1.45
3	P	6	THR	CA-CB	5.73	1.63	1.53
2	B	54	LEU	CA-C	-5.67	1.45	1.52
1	C	192	HIS	CA-C	5.67	1.59	1.52
1	A	76	VAL	CA-CB	5.66	1.61	1.54
2	B	86	THR	CA-CB	-5.65	1.44	1.53
1	A	67	ALA	CA-CB	-5.64	1.44	1.53
1	C	23	ILE	CA-C	-5.62	1.45	1.52
1	C	73	ILE	CG1-CD1	-5.61	1.29	1.51
2	D	57	SER	C-O	-5.59	1.17	1.24
1	A	179	LEU	CA-C	-5.58	1.45	1.52
2	B	48	LYS	CA-C	5.55	1.60	1.53
1	A	84	TYR	C-O	-5.51	1.17	1.24
1	C	44	ARG	CA-C	5.50	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	75	ARG	CA-C	5.47	1.60	1.52
1	A	22	PHE	N-CA	-5.45	1.39	1.46
2	B	46	ILE	CA-CB	-5.42	1.47	1.54
1	A	115	GLN	CB-CG	-5.41	1.36	1.52
1	A	48	ARG	CA-C	-5.39	1.45	1.52
1	C	31	THR	CA-C	5.39	1.59	1.52
2	B	41	LYS	N-CA	5.31	1.52	1.46
1	A	34	VAL	CA-CB	-5.30	1.47	1.55
1	A	29	ASP	C-O	-5.24	1.17	1.24
1	C	34	VAL	CA-C	-5.21	1.46	1.52
1	A	77	ASN	C-O	5.21	1.30	1.24
1	C	231	VAL	CA-CB	5.20	1.60	1.54
2	D	33	SER	CA-C	-5.19	1.45	1.52
1	C	29	ASP	C-O	-5.18	1.17	1.24
1	C	116	PHE	C-O	-5.16	1.17	1.24
1	C	209	TYR	CA-C	5.16	1.57	1.52
2	B	9	VAL	C-O	-5.15	1.18	1.24
1	A	95	LEU	N-CA	-5.13	1.40	1.46
1	A	21	ARG	C-O	-5.12	1.17	1.23
2	B	35	ILE	CA-C	-5.09	1.46	1.52
1	A	23	ILE	CA-CB	-5.08	1.47	1.55
1	A	165	VAL	CA-CB	5.08	1.60	1.54
2	B	7	ILE	CA-CB	-5.08	1.47	1.55
1	C	208	PHE	CA-C	5.08	1.60	1.53
1	A	185	PRO	CA-C	-5.04	1.47	1.52
1	A	203	CYS	CA-C	5.04	1.59	1.52
1	A	187	THR	C-O	-5.03	1.17	1.23
1	A	233	THR	C-O	-5.03	1.17	1.23
1	C	213	ILE	CA-CB	-5.02	1.50	1.55

All (154) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	249	VAL	N-CA-CB	-17.41	99.82	111.83
1	A	224	HIS	N-CA-C	-12.94	91.38	110.52
1	A	28	VAL	N-CA-C	-12.89	88.94	107.75
1	C	224	HIS	N-CA-C	11.16	134.57	110.80
1	A	76	VAL	N-CA-C	-10.60	100.57	110.53
1	C	28	VAL	N-CA-C	-10.04	92.52	107.37
1	C	17	ARG	N-CA-C	9.85	124.69	112.87
1	A	138	THR	N-CA-C	-9.69	101.59	113.50
1	C	76	VAL	N-CA-C	-9.58	101.53	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	107	GLY	N-CA-C	9.55	127.55	115.42
1	C	3	HIS	N-CA-C	9.48	123.56	109.24
1	C	257	TYR	N-CA-C	8.69	123.06	109.07
1	C	225	THR	N-CA-C	8.67	129.26	110.80
1	C	224	HIS	CA-C-N	8.62	137.99	121.54
1	C	224	HIS	C-N-CA	8.62	137.99	121.54
1	A	168	LEU	N-CA-C	-8.59	101.91	111.28
1	C	223	GLY	CA-C-N	8.56	137.89	121.54
1	C	223	GLY	C-N-CA	8.56	137.89	121.54
1	A	107	GLY	N-CA-C	8.55	126.69	115.40
1	C	138	THR	N-CA-C	-8.47	103.08	113.50
1	C	186	LYS	N-CA-C	-8.46	96.31	109.25
1	C	188	HIS	N-CA-C	8.45	121.42	108.42
1	C	49	ALA	CA-C-N	-8.37	111.19	119.64
1	C	49	ALA	C-N-CA	-8.37	111.19	119.64
1	C	168	LEU	N-CA-C	-8.36	102.17	111.28
1	A	224	HIS	CA-C-N	-8.13	109.94	122.62
1	A	224	HIS	C-N-CA	-8.13	109.94	122.62
1	A	139	ALA	N-CA-C	-7.89	102.68	111.28
1	A	186	LYS	N-CA-C	-7.86	98.65	110.28
1	C	226	GLN	N-CA-C	-7.79	100.76	111.28
1	A	257	TYR	N-CA-C	7.65	121.72	109.24
2	B	89	GLN	CA-C-N	-7.62	112.12	119.89
2	B	89	GLN	C-N-CA	-7.62	112.12	119.89
2	D	87	LEU	N-CA-C	7.60	121.20	108.20
1	C	143	SER	N-CA-C	-7.58	102.96	111.07
1	C	89	GLU	N-CA-C	-7.56	103.95	113.02
1	C	253	GLU	N-CA-C	7.50	122.11	113.19
1	A	225	THR	N-CA-C	-7.41	97.79	109.50
2	D	34	ASP	CA-C-N	-7.33	110.74	122.25
2	D	34	ASP	C-N-CA	-7.33	110.74	122.25
1	C	267	PRO	N-CA-C	-7.12	103.64	113.53
1	A	49	ALA	CA-C-N	-7.09	110.98	119.84
1	A	49	ALA	C-N-CA	-7.09	110.98	119.84
1	A	179	LEU	N-CA-C	7.00	118.99	111.36
1	C	256	ALA	N-CA-C	-6.86	104.77	113.01
1	A	6	LYS	N-CA-C	6.84	120.69	109.06
1	A	198	GLU	N-CA-C	6.84	118.91	108.52
1	C	184	PRO	CA-C-N	6.75	126.51	119.76
1	C	184	PRO	C-N-CA	6.75	126.51	119.76
2	B	92	ILE	CA-C-N	-6.55	114.56	123.14
2	B	92	ILE	C-N-CA	-6.55	114.56	123.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	143	SER	N-CA-C	-6.54	104.24	111.36
1	C	249	VAL	CA-C-N	6.52	130.05	120.46
1	C	249	VAL	C-N-CA	6.52	130.05	120.46
1	C	30	ASP	N-CA-C	6.44	124.51	110.80
1	C	204	TRP	N-CA-C	6.35	120.18	109.76
2	D	67	TYR	N-CA-C	6.35	118.50	108.79
2	D	13	HIS	CA-C-N	-6.29	113.28	119.76
2	D	13	HIS	C-N-CA	-6.29	113.28	119.76
2	B	33	SER	N-CA-C	6.25	120.76	113.20
1	C	227	ASP	CA-C-N	6.10	131.37	121.74
1	C	227	ASP	C-N-CA	6.10	131.37	121.74
1	A	24	SER	N-CA-CB	-6.09	100.55	110.59
1	C	115	GLN	N-CA-C	6.03	118.35	109.24
1	C	208	PHE	N-CA-C	6.03	119.44	110.52
1	C	146	LYS	N-CA-C	-6.02	104.72	111.28
1	A	197	HIS	N-CA-C	6.00	120.62	112.88
2	B	12	ARG	NH1-CZ-NH2	-6.00	111.50	119.30
1	A	180	LEU	CA-CB-CG	-5.99	95.35	116.30
1	C	227	ASP	N-CA-C	5.99	123.55	110.80
2	D	89	GLN	CA-C-N	-5.92	113.85	119.89
2	D	89	GLN	C-N-CA	-5.92	113.85	119.89
1	C	197	HIS	N-CA-C	5.84	120.11	112.92
2	D	83	ASN	N-CA-C	5.84	119.07	109.72
1	A	142	ILE	CB-CA-C	-5.84	104.26	112.14
1	A	267	PRO	N-CA-C	-5.80	105.82	113.65
1	A	225	THR	CB-CA-C	5.78	121.45	109.38
1	A	225	THR	N-CA-CB	-5.78	100.86	111.37
1	C	224	HIS	O-C-N	5.77	130.26	122.59
1	A	211	ALA	N-CA-C	5.75	120.03	112.89
1	C	172	LEU	N-CA-C	-5.73	105.04	111.28
1	C	6	LYS	N-CA-C	5.70	117.94	108.99
1	C	142	ILE	CB-CA-C	-5.66	104.61	112.02
1	A	162	ASP	O-C-N	5.64	126.95	120.58
1	A	87	GLN	N-CA-C	5.61	118.62	109.59
3	Q	8	LEU	N-CA-C	-5.57	100.10	108.96
1	A	268	GLU	CA-C-N	5.57	125.57	119.89
1	A	268	GLU	C-N-CA	5.57	125.57	119.89
1	A	260	HIS	CA-C-N	-5.56	115.70	123.10
1	A	260	HIS	C-N-CA	-5.56	115.70	123.10
1	A	162	ASP	CA-C-O	5.56	123.63	119.68
1	A	253	GLU	N-CA-C	5.56	119.92	113.20
2	D	12	ARG	NE-CZ-NH2	5.50	124.15	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	184	PRO	CA-C-N	5.46	126.95	120.23
1	A	184	PRO	C-N-CA	5.46	126.95	120.23
2	B	6	LYS	N-CA-C	-5.46	100.33	109.24
1	A	30	ASP	N-CA-C	5.44	122.39	110.80
1	A	162	ASP	CB-CA-C	-5.43	104.69	114.52
1	C	131	ARG	NE-CZ-NH1	-5.43	116.07	121.50
2	D	12	ARG	NH1-CZ-NH2	-5.41	112.27	119.30
2	D	45	ARG	NE-CZ-NH2	5.37	124.03	119.20
1	A	11	SER	CA-C-N	-5.36	116.11	123.14
1	A	11	SER	C-N-CA	-5.36	116.11	123.14
1	C	237	GLY	N-CA-C	-5.36	107.81	115.64
1	C	245	ALA	CA-C-N	-5.35	114.27	122.62
1	C	245	ALA	C-N-CA	-5.35	114.27	122.62
2	B	73	THR	CA-C-N	-5.34	111.80	121.14
2	B	73	THR	C-N-CA	-5.34	111.80	121.14
1	A	124	LEU	CA-C-N	-5.32	115.49	122.99
1	A	124	LEU	C-N-CA	-5.32	115.49	122.99
1	A	105	PRO	N-CA-C	-5.32	106.93	113.84
3	Q	3	ALA	CA-C-N	5.32	124.77	119.24
3	Q	3	ALA	C-N-CA	5.32	124.77	119.24
1	A	226	GLN	N-CA-C	-5.31	102.20	110.42
1	C	24	SER	N-CA-CB	-5.30	102.12	110.77
1	A	79	ARG	N-CA-C	-5.29	105.43	111.14
1	A	187	THR	N-CA-C	5.29	117.72	109.52
1	A	200	THR	CA-C-N	-5.29	115.70	123.00
1	A	200	THR	C-N-CA	-5.29	115.70	123.00
1	A	256	ALA	N-CA-C	-5.29	105.93	112.38
1	A	183	GLU	CA-C-N	-5.28	114.94	120.38
1	A	183	GLU	C-N-CA	-5.28	114.94	120.38
1	A	3	HIS	N-CA-C	5.27	117.69	109.52
1	C	144	GLU	N-CA-C	5.26	117.02	111.28
1	C	124	LEU	CA-C-N	-5.26	114.21	122.73
1	C	124	LEU	C-N-CA	-5.26	114.21	122.73
2	D	45	ARG	NE-CZ-NH1	-5.24	116.26	121.50
2	D	38	ASP	N-CA-C	5.23	117.92	109.40
1	A	137	ASP	CA-C-N	-5.21	112.91	121.92
1	A	137	ASP	C-N-CA	-5.21	112.91	121.92
1	C	229	GLU	N-CA-C	-5.21	98.53	107.23
1	C	79	ARG	N-CA-C	-5.20	105.26	111.03
1	C	14	ARG	CA-C-N	5.18	126.32	119.84
1	C	14	ARG	C-N-CA	5.18	126.32	119.84
2	B	46	ILE	O-C-N	-5.17	117.39	122.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3	ARG	NE-CZ-NH2	5.16	123.85	119.20
2	D	6	LYS	N-CA-C	-5.16	100.82	109.24
1	C	104	GLY	N-CA-C	-5.14	105.78	112.10
1	C	20	PRO	CA-C-N	5.14	128.62	121.02
1	C	20	PRO	C-N-CA	5.14	128.62	121.02
1	C	131	ARG	NE-CZ-NH2	5.13	123.82	119.20
1	C	105	PRO	N-CA-C	-5.13	107.17	113.84
1	C	94	THR	CA-C-N	-5.11	115.80	123.00
1	C	94	THR	C-N-CA	-5.11	115.80	123.00
2	D	79	ALA	N-CA-C	5.11	115.27	108.38
1	A	185	PRO	N-CA-C	5.10	118.50	110.50
2	D	58	LYS	N-CA-C	5.09	117.73	111.82
2	B	75	LYS	N-CA-C	5.08	119.36	112.04
3	P	2	MET	N-CA-C	-5.07	103.70	110.55
1	A	249	VAL	CA-C-N	5.06	126.16	119.84
1	A	249	VAL	C-N-CA	5.06	126.16	119.84
1	C	133	TRP	N-CA-C	5.05	117.80	109.72
1	C	249	VAL	N-CA-CB	-5.05	104.14	111.21
1	A	146	LYS	N-CA-C	-5.02	105.09	111.11

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 [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	2206	0	2028	167	0
1	C	2205	0	2027	179	0
2	B	828	0	780	55	0
2	D	828	0	780	62	0
3	P	69	0	83	12	0
3	Q	69	0	83	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	5	0	0	1	0
All	All	6210	0	5781	435	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 36.

All (435) 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	2.08	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:HH11	1:C:65:ARG:HB2	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:HH11	1:A:65:ARG:CB	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:138:THR:HA	1:A:141:GLN:HG3	1.55	0.87
1:A:218:GLN:CB	1:A:223:GLY:HA2	2.03	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
1:A:65:ARG:HB2	1:A:65:ARG:NH1	1.91	0.85
1:C:21:ARG:HG3	1:C:21:ARG:HH11	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:HH11	1:A:65:ARG:HB3	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HH11	1:A:65:ARG:HB2	1.43	0.78
2:D:42:ASN:ND2	2:D:77:GLU:H	1.84	0.76
1:A:35:ARG:HG2	1:A:35:ARG:HH11	1.49	0.76
2:D:42:ASN:HD21	2:D:77:GLU:H	1.32	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:HG3	1:A:21:ARG:HH11	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:A:177:GLU:CD	1:A:177:GLU:H	1.96	0.73
1:C:65:ARG:HH11	1:C:65:ARG:HB3	1.53	0.73
1:A:121:LYS:HG3	1:A:122:ASP:N	2.03	0.73
2:D:4:THR:HG22	2:D:5:PRO:HD2	1.71	0.73
2:D:73:THR:HG22	2:D:75:LYS:N	2.04	0.73
1:C:202:ARG:NH2	2:D:99:MET:HE3	2.05	0.72
1:A:194:ILE:HD12	1:A:194:ILE:N	2.05	0.72
2:D:9:VAL:HG23	2:D:93:VAL:CG2	2.19	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
1:A:98:MET:HE2	2:B:60:TRP:CZ2	2.28	0.69
1:A:211:ALA:HB2	1:A:241:PHE:CE1	2.27	0.69
1:C:230:LEU:CD1	1:C:243:LYS:HE3	2.22	0.69
1:A:98:MET:HE2	2:B:60:TRP:CH2	2.28	0.69
2:D:51:HIS:HA	2:D:65:LEU:O	1.92	0.69
1:A:202:ARG:HG3	1:A:202:ARG:HH11	1.57	0.69
1:A:55:GLU:HA	1:A:55:GLU:OE1	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:C:50:PRO:HA	1:C:53:GLU:OE2	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:50:PRO:HA	1:A:53:GLU:OE2	1.94	0.68
1:A:65:ARG:CB	1:A:65:ARG:NH1	2.47	0.68
1:A:271:THR:C	1:A:272:LEU:HD23	2.19	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:HG3	1:C:202:ARG:HH11	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:A:255:GLN:CD	1:A:255:GLN:N	2.53	0.65
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
2:D:29:GLY:HA2	2:D:61:SER:OG	1.97	0.65
3:P:5:ARG:NH2	3:Q:5:ARG:HD3	2.12	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: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:A:8:PHE:HB3	2:B:56:PHE:CE2	2.34	0.63
1:C:98:MET:HE2	2:D:60:TRP:CH2	2.34	0.63
1:C:249:VAL:HG22	1:C:257:TYR:CE1	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
2:D:16:GLU:O	2:D:17:ASN:C	2.42	0.62
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:VAL:CG1	1:A:190:THR:N	2.62	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:204:TRP:CE3	1:C:206:LEU:HD21	2.34	0.62
1:C:263:HIS:HD2	1:C:265:GLY:N	1.90	0.61
1:A:220:ASP:O	1:A:220:ASP:OD2	2.17	0.61
1:C:21:ARG:HH11	1:C:21:ARG:CG	2.11	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
2:D:98:ASP:C	2:D:99:MET:HG3	2.24	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:C:221:GLY:C	1:C:222:GLU:HG3	2.27	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
1:A:150:ALA:O	1:A:151:SER:CB	2.50	0.59
2:B:51:HIS:HA	2:B:65:LEU:O	2.03	0.59
1:C:150:ALA:O	1:C:151:SER:CB	2.48	0.59
1:C:152:GLU:CD	3:P:5:ARG:HH12	2.11	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:209:TYR:HA	1:C:210:PRO:O	2.03	0.58
1:C:229:GLU:OE2	1:C:244:TRP:HH2	1.86	0.58
1:C:138:THR:HA	1:C:141:GLN:HE21	1.68	0.58
1:C:209:TYR:HA	1:C:210:PRO:C	2.27	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:35:ARG:HH11	1:C:46:VAL:HG21	1.67	0.57
1:C:121:LYS:HG3	1:C:122:ASP:N	2.19	0.57
1:A:65:ARG:NH2	1:C:145:GLN:NE2	2.45	0.57
1:A:145:GLN:CD	1:C:65:ARG:HH21	2.11	0.57
1:C:128:GLU:O	1:C:130:LEU:HG	2.03	0.57
1:A:47:PRO:HB3	1:A:60:TRP:CZ2	2.40	0.57
1:A:194:ILE:HD13	1:A:199:ALA:HA	1.86	0.57
1:A:177:GLU:CD	1:A:177:GLU:N	2.64	0.56
2:B:49:VAL:CG1	2:B:50:GLU:N	2.68	0.56
1:A:222:GLU:O	1:A:223:GLY:C	2.47	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:47:PRO:HB3	1:C:60:TRP:CZ2	2.41	0.56
2:B:17:ASN:OD1	2:B:73:THR:C	2.48	0.56
1:C:82:ARG:NE	1:C:89:GLU:HB2	2.19	0.56
1:A:194:ILE:N	1:A:194:ILE:CD1	2.69	0.56
1:A:13:SER:HA	1:A:20:PRO:HB3	1.89	0.55
1:C:44:ARG:NH2	1:C:61:ASP:OD1	2.39	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:N	1:C:194:ILE:CD1	2.70	0.54
1:C:166:GLU:O	1:C:166:GLU:HG2	2.06	0.54
1:A:270:VAL:HG12	1:A:272:LEU:CD2	2.37	0.54
2:B:9:VAL:HG23	2:B:93:VAL:CG2	2.36	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:103:LEU:HA	1:A:108:ARG:O	2.07	0.53
1:A:193:PRO:HA	1:A:199:ALA:HA	1.88	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:C:244:TRP:C	1:C:244:TRP:HE3	2.15	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:85:TYR:O	1:A:86:ASN:C	2.50	0.53
1:A:193:PRO:HA	1:A:199:ALA:CA	2.39	0.53
1:C:202:ARG:HG3	1:C:202:ARG:NH1	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: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
1:A:48:ARG:NH1	2:B:53:ASP:OD2	2.42	0.52
2:B:4:THR:CG2	2:B:5:PRO:HD2	2.40	0.52
1:A:37:ASP:C	1:A:39:ASP:H	2.16	0.52
2:D:40:LEU:HD11	2:D:81:ARG:HB2	1.91	0.52
1:A:23:ILE:HG22	1:A:24:SER:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:203:CYS:HB2	1:A:217:TRP:CZ2	2.45	0.51
2:D:9:VAL:HG21	2:D:93:VAL:CG2	2.37	0.51
2:B:29:GLY:HA2	2:B:61:SER:OG	2.10	0.51
2:B:73:THR:CG2	2:B:76:ASP:H	2.23	0.51
1:C:162:ASP:O	1:C:163:THR:C	2.52	0.51
1:A:70:THR:O	1:A:71:ALA:C	2.52	0.51
1:C:44:ARG:HA	1:C:64:THR:CG2	2.41	0.51
2:B:16:GLU:O	2:B:17:ASN:C	2.53	0.51
2:D:9:VAL:CG2	2:D:93:VAL:HG23	2.41	0.51
1:C:55:GLU:CD	1:C:174:LYS:NZ	2.69	0.51
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
3:P:5:ARG:NH2	3:Q:5:ARG:HE	2.07	0.50
1:C:193:PRO:HA	1:C:199:ALA:HA	1.93	0.50
1:A:104:GLY:C	1:A:106:ASP:N	2.69	0.50
1:A:243:LYS:HG2	1:A:244:TRP:H	1.74	0.50
2:B:98:ASP:C	2:B:99:MET:HG3	2.36	0.50
1:C:215:LEU:CD2	1:C:261:VAL:HG22	2.42	0.50
1:C:250:PRO:O	1:C:251:SER:C	2.55	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:A:51:TRP:HB2	1:A:174:LYS:O	2.12	0.49
1:C:35:ARG:HH12	1:C:46:VAL:HG21	1.76	0.49
1:C:98:MET:HE2	2:D:60:TRP:CZ2	2.46	0.49
1:C:138:THR:HA	1:C:141:GLN:HG3	1.93	0.49
2:D:29:GLY:HA2	2:D:61:SER:CB	2.42	0.49
1:A:202:ARG:HG3	1:A:202:ARG:NH1	2.25	0.49
1:A:79:ARG:HD2	1:C:72:GLN:NE2	2.27	0.49
2:B:73:THR:HG22	2:B:76:ASP:H	1.76	0.49
1:C:244:TRP:C	1:C:244:TRP:CE3	2.90	0.49
2:D:9:VAL:HG21	2:D:93:VAL:O	2.13	0.49
1:A:82:ARG:HG3	1:A:87:GLN:HB2	1.94	0.49
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.47	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:37:ASP:C	1:A:39:ASP:N	2.69	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
2:D:9:VAL:HA	2:D:24:ASN:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:C:218:GLN:HE21	1:C:260:HIS:CE1	2.31	0.49
1:A:254:GLU:OE1	1:A:254:GLU:N	2.42	0.48
1:C:219:GLN:NE2	1:C:220:ASP:OD2	2.44	0.48
1:A:127:ASN:ND2	1:A:132:SER:OG	2.44	0.48
1:A:202:ARG:HD3	1:A:244:TRP:CD2	2.49	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:55:GLU:HA	1:C:55:GLU:OE1	2.12	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
1:C:53:GLU:C	1:C:55:GLU:H	2.21	0.48
1:A:76:VAL:HG11	3:P:8:LEU:HD11	1.96	0.48
1:A:232:GLU:O	1:A:233:THR:C	2.54	0.48
2:D:73:THR:HG22	2:D:76:ASP:H	1.77	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: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:195:SER:OG	1:A:198:GLU:HG3	2.14	0.48
1:A:23:ILE:CG2	1:A:24:SER:N	2.75	0.47
1:A:51:TRP:CZ3	1:A:52:MET:SD	3.07	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
2:B:89:GLN:O	2:B:90:PRO:C	2.55	0.47
1:C:33:PHE:C	1:C:48:ARG:HB2	2.40	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:224:HIS:C	1:A:225:THR:HG23	2.40	0.47
1:A:216:THR:O	1:A:259:CYS:HA	2.15	0.47
2:B:42:ASN:ND2	2:B:77:GLU:HB2	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
1:A:28:VAL:O	1:A:29:ASP:O	2.34	0.46
1:A:226:GLN:O	1:A:227:ASP:C	2.56	0.46
1:A:255:GLN:C	1:A:257:TYR:N	2.72	0.46
1:C:127:ASN:O	1:C:130:LEU:N	2.47	0.46
1:C:212:GLU:HA	1:C:212:GLU:OE1	2.15	0.46
3:P:5:ARG:HH21	3:Q:5:ARG:NE	2.08	0.46
1:A:255:GLN:NE2	1:A:255:GLN:H	2.10	0.46
2:B:31:HIS:ND1	2:B:32:PRO:HA	2.31	0.46
1:C:109:PHE:CE2	1:C:161:GLU:HA	2.50	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:A:79:ARG:O	1:A:80:THR:C	2.57	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:129:ASP:OD2	1:A:131:ARG:HG3	2.16	0.46
1:A:155:HIS:CE1	1:C:155:HIS:ND1	2.83	0.46
1:A:103:LEU:HG	1:A:168:LEU:HD23	1.96	0.45
1:A:198:GLU:HB3	1:A:250:PRO:HA	1.98	0.45
1:A:202:ARG:NH2	2:B:99:MET:HE3	2.30	0.45
2:B:73:THR:HB	2:B:76:ASP:HB2	1.98	0.45
1:A:202:ARG:HD3	1:A:244:TRP:CE2	2.51	0.45
2:B:16:GLU:HG2	2:B:19:LYS:CB	2.47	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:C:219:GLN:HE21	1:C:220:ASP:CG	2.23	0.45
2:D:70:PHE:CD1	2:D:70:PHE:C	2.93	0.45
1:A:121:LYS:HE2	2:B:1(A):ILE:HD11	1.99	0.45
1:C:202:ARG:HH22	2:D:99:MET:HE3	1.81	0.45
1:A:55:GLU:CD	1:A:174:LYS:NZ	2.75	0.45
1:C:115:GLN:CG	1:C:125:THR:HG23	2.34	0.45
1:C:180:LEU:HD23	1:C:180:LEU:HA	1.66	0.45
2:D:51:HIS:HB3	2:D:66:TYR:CD2	2.52	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:A:79:ARG:NH1	1:C:72:GLN:HE22	2.15	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:21:ARG:HD3	1:C:38:ASN:OD1	2.17	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
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:A:162:ASP:HB2	1:A:163:THR:H	1.40	0.44
1:C:253:GLU:O	1:C:255:GLN:N	2.51	0.44
2:B:96:ASP:OD1	2:B:98:ASP:HB2	2.18	0.44
1:C:104:GLY:C	1:C:106:ASP:N	2.75	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:145:GLN:CD	1:C:65:ARG:NH2	2.75	0.44
1:A:166:GLU:O	1:A:169:HIS:HB2	2.18	0.44
1:A:81:LEU:HD21	1:A:123:TYR:CZ	2.53	0.43
1:A:191:HIS:ND1	1:A:274:TRP:CH2	2.83	0.43
1:C:29:ASP:O	1:C:31:THR:N	2.50	0.43
1:C:82:ARG:HG3	1:C:87:GLN:O	2.18	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: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:N	1:C:33:PHE:CD1	2.86	0.43
1:C:109:PHE:CD1	1:C:109:PHE:C	2.97	0.43
1:A:89:GLU:O	1:A:89:GLU:HG2	2.19	0.43
1:A:145:GLN:OE1	1:C:65:ARG:NH2	2.46	0.43
1:A:250:PRO:O	1:A:251:SER:C	2.59	0.43
1:A:272:LEU:HD23	1:A:272:LEU:N	2.34	0.43
1:C:166:GLU:OE2	1:C:170:LYS:HE3	2.19	0.43
1:A:209:TYR:CD1	1:A:209:TYR:C	2.97	0.43
1:C:243:LYS:HG2	1:C:244:TRP:N	2.34	0.43
1:A:44:ARG:HA	1:A:64:THR:HG23	2.01	0.43
1:A:83:GLY:O	1:A:84:TYR:C	2.61	0.43
1:C:128:GLU:H	1:C:128:GLU:HG2	1.58	0.43
1:C:103:LEU:HG	1:C:168:LEU:HD23	2.01	0.42
3:P:5:ARG:CZ	3:Q:5:ARG:NE	2.82	0.42
1:A:129:ASP:O	1:A:130:LEU:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:PHE:CD2	1:A:161:GLU:HA	2.54	0.42
2:B:6:LYS:HB2	2:B:6:LYS:HE3	1.89	0.42
2:B:21:ASN:CG	2:B:22:PHE:H	2.26	0.42
1:C:271:THR:C	1:C:272:LEU:HD23	2.45	0.42
1:C:194:ILE:HD13	1:C:199:ALA:CA	2.46	0.42
1:A:186:LYS:HD3	1:A:186:LYS:N	2.33	0.42
1:C:117:ALA:HB2	2:D:60:TRP:CZ2	2.54	0.42
1:C:162:ASP:O	1:C:164:CYS:N	2.53	0.42
1:C:249:VAL:HA	1:C:250:PRO:HD3	1.69	0.42
1:A:104:GLY:C	1:A:106:ASP:H	2.26	0.42
1:C:6:LYS:HD2	1:C:113:TYR:OH	2.20	0.42
1:C:22:PHE:H	1:C:38:ASN:CG	2.26	0.42
1:A:39:ASP:O	1:A:40:ALA:C	2.62	0.42
1:A:249:VAL:HA	1:A:250:PRO:HD3	1.77	0.42
2:D:6:LYS:O	2:D:27:VAL:HA	2.19	0.42
1:A:109:PHE:CD1	1:A:109:PHE:C	2.98	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:229:GLU:OE2	1:C:244:TRP:CH2	2.70	0.42
1:A:224:HIS:C	1:A:225:THR:CG2	2.91	0.42
1:C:191:HIS:ND1	1:C:274:TRP:CH2	2.77	0.42
1:C:209:TYR:CA	1:C:210:PRO:O	2.67	0.42
1:A:21:ARG:HH11	1:A:21:ARG:CG	2.24	0.42
1:A:235:PRO:O	2:B:10:TYR:OH	2.31	0.42
1:C:39:ASP:O	1:C:40:ALA:O	2.38	0.42
1:C:233:THR:HG22	1:C:234:ARG:N	2.34	0.42
1:A:184:PRO:HB3	1:A:265:GLY:C	2.44	0.41
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:A:76:VAL:O	1:A:77:ASN:C	2.64	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:C:230:LEU:HD11	1:C:243:LYS:HE3	1.99	0.41
2:D:89:GLN:O	2:D:90:PRO:C	2.60	0.41
1:A:33:PHE:CD1	1:A:33:PHE:N	2.89	0.41
1:A:45:MET:O	1:A:60:TRP:CE3	2.73	0.41
1:A:65:ARG:O	1:A:66:SER:C	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:71:THR:HA	2:B:72:PRO:HD2	1.69	0.41
1:A:143:SER:O	1:A:147:SER:HB2	2.20	0.41
1:A:191:HIS:CE1	1:A:254:GLU:OE2	2.74	0.41
1:C:48:ARG:HD2	1:C:48:ARG:HA	1.47	0.41
2:D:71:THR:HA	2:D:72:PRO:HD2	1.67	0.41
1:C:14:ARG:HD3	1:C:17:ARG:HH12	1.86	0.41
1:C:35:ARG:NH1	1:C:46:VAL:CG2	2.70	0.41
1:C:124:LEU:HA	1:C:124:LEU:HD12	1.83	0.41
1:C:194:ILE:CD1	1:C:199:ALA:HA	2.49	0.41
1:C:231:VAL:HG13	1:C:244:TRP:CZ3	2.56	0.41
2:D:9:VAL:HG23	2:D:93:VAL:HG23	1.99	0.41
2:B:29:GLY:HA2	2:B:61:SER:CB	2.51	0.41
2:B:23:LEU:HD12	2:B:23:LEU:HA	1.79	0.40
2:D:39:LEU:HD13	2:D:68:THR:HG22	2.02	0.40
1:A:127:ASN:HD22	1:A:132:SER:C	2.29	0.40
1:C:63:GLU:CD	3:Q:1:VAL:HG12	2.46	0.40
2:D:84:HIS:O	2:D:85:VAL:C	2.63	0.40
1:A:57:SER:O	1:A:58:GLU:C	2.64	0.40
2:B:49:VAL:CG1	2:B:50:GLU:H	2.34	0.40
1:C:115:GLN:HB3	2:D:60:TRP:CH2	2.55	0.40
2:D:3:ARG:HG2	2:D:3:ARG:HH11	1.86	0.40
2:D:98:ASP:O	2:D:99:MET:CG	2.49	0.40
1:C:109:PHE:CD2	1:C:161:GLU:HA	2.56	0.40
1:C:117:ALA:HA	1:C:121:LYS:O	2.21	0.40
1:C:192:HIS:O	1:C:200:THR:N	2.50	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:161:GLU:O	1:A:165:VAL:CG1	2.70	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
2:B:96:ASP:C	2:B:98:ASP:H	2.29	0.40
1:C:9:HIS:CE1	3:Q:2:MET:HE3	2.56	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

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	272/274 (99%)	231 (85%)	31 (11%)	10 (4%)	2	9
1	C	272/274 (99%)	230 (85%)	24 (9%)	18 (7%)	1	3
2	B	98/100 (98%)	92 (94%)	5 (5%)	1 (1%)	12	38
2	D	98/100 (98%)	89 (91%)	7 (7%)	2 (2%)	6	21
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%)	2	8

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

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Mol	Chain	Res	Type
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%)	194 (84%)	36 (16%)	2	9
1	C	229/234 (98%)	193 (84%)	36 (16%)	2	9
2	B	93/95 (98%)	85 (91%)	8 (9%)	10	31
2	D	93/95 (98%)	82 (88%)	11 (12%)	5	17
3	P	8/8 (100%)	6 (75%)	2 (25%)	0	2
3	Q	8/8 (100%)	7 (88%)	1 (12%)	4	15
All	All	661/674 (98%)	567 (86%)	94 (14%)	3	12

All (94) 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

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Mol	Chain	Res	Type
1	A	55	GLU
1	A	62	ARG
1	A	65	ARG
1	A	73	ILE
1	A	76	VAL
1	A	136	VAL
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	173	GLU
1	A	178	THR
1	A	179	LEU
1	A	182	LEU
1	A	186	LYS
1	A	187	THR
1	A	194	ILE
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	268	GLU
1	A	272	LEU
1	A	273	ARG
2	B	4	THR
2	B	9	VAL
2	B	12	ARG
2	B	48	LYS
2	B	58	LYS
2	B	71	THR
2	B	73	THR
2	B	93	VAL
1	C	11	SER
1	C	14	ARG
1	C	15	PRO
1	C	21	ARG
1	C	29	ASP

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Mol	Chain	Res	Type
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	136	VAL
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	165	VAL
1	C	173	GLU
1	C	177	GLU
1	C	186	LYS
1	C	187	THR
1	C	194	ILE
1	C	198	GLU
1	C	200	THR
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
1	C	273	ARG
2	D	4	THR
2	D	11	SER
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
2	D	93	VAL

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	87	GLN
1	A	115	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	141	GLN
1	C	145	GLN
1	C	188	HIS
1	C	218	GLN
1	C	219	GLN
1	C	255	GLN
1	C	260	HIS
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 [i](#)

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](#)

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	0.39	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 ⓘ

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.