



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:57 PM UTC

PDB ID : 1KJ2 / pdb_00001kj2
Title : Murine Alloreactive ScFv TCR-Peptide-MHC Class I Molecule Complex
Authors : Reiser, J.-B.; Gregoire, C.; Darnault, C.; Mosser, T.; Guimezanes, A.; Schmitt-Verhulst, A.-M.; Fontecilla-Camps, J.C.; Mazza, G.; Malissen, B.; Housset, D.
Deposited on : 2001-12-04
Resolution : 2.71 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)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

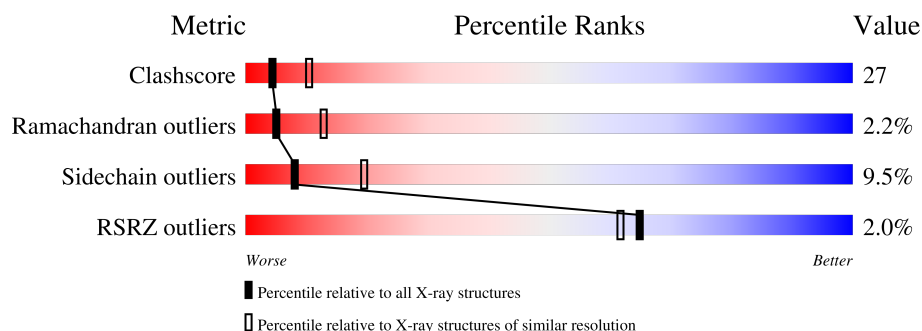
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	H	277	% 61% 35% .
1	I	277	4% 56% 38% 6% .
2	P	8	38% 50% 12%
2	Q	8	25% 50% 25%
3	L	99	57% 39% .
3	M	99	% 49% 43% 7%

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Mol	Chain	Length	Quality of chain
4	A	111	
4	D	111	
5	B	117	
5	E	117	
6	C	8	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	MAN	C	3	X	-	-	-
6	MAN	C	8	X	-	-	-
7	NAG	E	401	X	-	-	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 10144 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 Allogeneic H-2Kb MHC Class I Molecule.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	276	Total	C	N	O	S	0	0	0
			2247	1418	395	425	9			
1	I	277	Total	C	N	O	S	0	0	0
			2254	1423	396	426	9			

- Molecule 2 is a protein called Naturally processed octapeptide PKB1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	8	Total	C	N	O	0	0	0
			67	46	9	12			
2	Q	8	Total	C	N	O	0	0	0
			67	46	9	12			

- Molecule 3 is a protein called Beta-2 microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	99	Total	C	N	O	S	0	0	0
			821	524	138	152	7			
3	M	99	Total	C	N	O	S	0	0	0
			821	524	138	152	7			

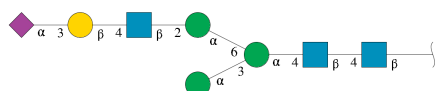
- Molecule 4 is a protein called KB5-C20 T-Cell receptor alpha-chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	A	111	Total	C	N	O	S	0	0	0
			881	560	149	170	2			
4	D	111	Total	C	N	O	S	0	0	0
			881	560	149	170	2			

- Molecule 5 is a protein called KB5-C20 T-Cell receptor beta-chain.

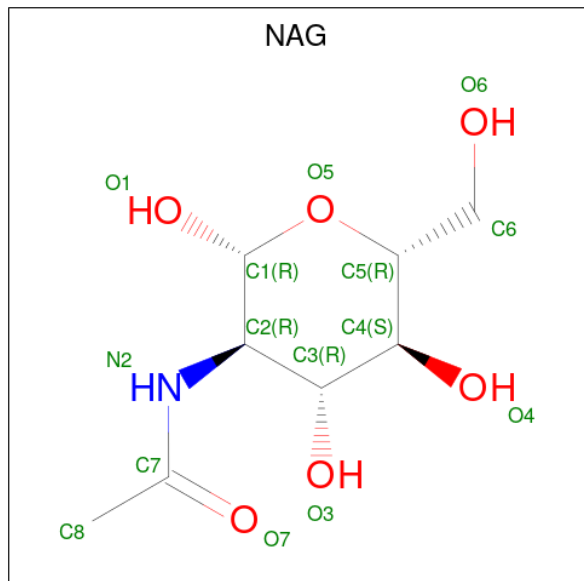
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	B	117	Total	C	N	O	S	0	0	0
			940	596	167	172	5			
5	E	117	Total	C	N	O	S	0	0	0
			940	596	167	172	5			

- Molecule 6 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	C	8	Total	C	N	O		0	0	0
			106	59	4	43				

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O		0	0
			14	8	1	5			
7	E	1	Total	C	N	O		0	0
			14	8	1	5			

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	H	22	Total O 22 22	0	0
8	P	2	Total O 2 2	0	0
8	L	12	Total O 12 12	0	0
8	A	5	Total O 5 5	0	0
8	B	10	Total O 10 10	0	0
8	I	18	Total O 18 18	0	0
8	Q	3	Total O 3 3	0	0
8	M	6	Total O 6 6	0	0
8	D	6	Total O 6 6	0	0
8	E	7	Total O 7 7	0	0

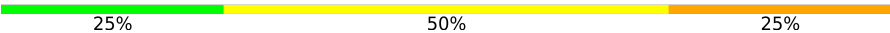
3 Residue-property plots [i](#)

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 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.

• Molecule 1: Allogeneic H-2Kb MHC Class I Molecule



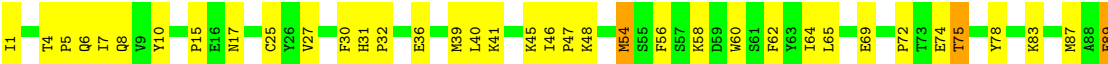
- Molecule 2: Naturally processed octapeptide PKB1

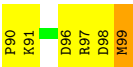
Chain Q: 



- Molecule 3: Beta-2 microglobulin

Chain L: 

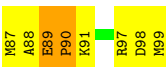




- Molecule 3: Beta-2 microglobulin

Chain M: 





- Molecule 4: KB5-C20 T-Cell receptor alpha-chain

Chain A: 

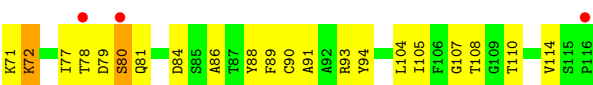




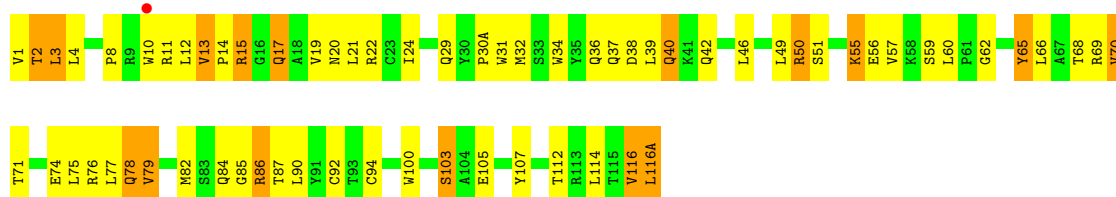
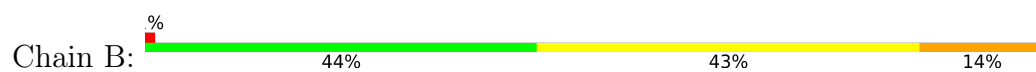
- Molecule 4: KB5-C20 T-Cell receptor alpha-chain

Chain D: 

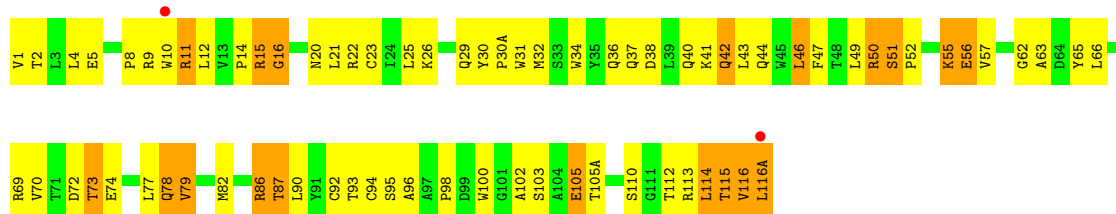




- Molecule 5: KB5-C20 T-Cell receptor beta-chain



- Molecule 5: KB5-C20 T-Cell receptor beta-chain



- Molecule 6: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)] alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.20Å 77.92Å 132.96Å 90.00° 108.23° 90.00°	Depositor
Resolution (Å)	12.00 – 2.71 12.00 – 2.71	Depositor EDS
% Data completeness (in resolution range)	(Not available) (12.00-2.71) 94.4 (12.00-2.71)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.19 (at 2.73Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.220 , 0.278 0.230 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.355	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10144	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.74 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.9477e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GAL, NAG, SIA, MAN

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	H	0.55	0/2309	1.01	9/3137 (0.3%)
1	I	0.54	0/2317	1.04	15/3149 (0.5%)
2	P	0.76	0/67	1.08	0/88
2	Q	0.67	0/67	0.97	0/88
3	L	0.58	0/847	1.05	3/1148 (0.3%)
3	M	0.49	0/847	1.01	3/1148 (0.3%)
4	A	0.49	0/904	0.93	1/1225 (0.1%)
4	D	0.47	0/904	0.99	3/1225 (0.2%)
5	B	0.54	0/962	1.09	7/1310 (0.5%)
5	E	0.52	0/962	1.15	8/1310 (0.6%)
All	All	0.53	0/10186	1.03	49/13828 (0.4%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	3	LEU	N-CA-C	-9.42	102.29	113.88
5	E	110	SER	N-CA-C	-9.02	102.05	113.23
5	E	44	GLN	N-CA-C	7.64	120.53	107.93
1	H	249	VAL	N-CA-C	7.55	117.54	109.01
1	H	168	LEU	N-CA-C	-7.14	103.58	111.36
1	I	249	VAL	CA-C-N	6.92	127.30	119.83
1	I	249	VAL	C-N-CA	6.92	127.30	119.83
5	E	51	SER	CA-C-N	6.90	126.87	119.76
5	E	51	SER	C-N-CA	6.90	126.87	119.76
3	M	54	MET	N-CA-C	6.82	118.65	110.19
1	H	236	ALA	N-CA-C	-6.62	104.69	112.89
4	A	84	ASP	N-CA-C	-6.60	105.09	113.01
3	M	2	GLN	N-CA-C	-6.57	96.26	107.23
1	H	82	LEU	N-CA-C	-6.51	104.22	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	257	TYR	N-CA-C	6.50	120.87	109.96
5	E	116	VAL	N-CA-C	6.41	119.24	109.12
1	I	249	VAL	N-CA-C	6.41	116.14	108.45
1	I	226	GLN	N-CA-C	6.38	119.04	111.71
5	E	51	SER	N-CA-C	6.21	118.46	110.39
1	I	163	THR	N-CA-C	6.19	118.57	111.02
4	D	84	ASP	N-CA-C	-6.18	104.90	112.88
5	B	51	SER	N-CA-C	6.06	119.37	110.10
4	D	104	LEU	N-CA-C	-6.03	99.44	109.46
1	I	174	ASN	N-CA-C	5.89	117.70	111.28
1	I	168	LEU	N-CA-C	-5.87	104.80	111.14
3	L	75	THR	N-CA-C	5.83	120.56	112.45
4	D	17	THR	N-CA-C	5.77	118.07	108.20
1	I	130	LEU	N-CA-C	-5.75	104.88	112.24
5	B	103	SER	N-CA-C	-5.73	105.49	112.88
5	E	16	GLY	N-CA-C	5.56	120.82	113.37
1	I	44	ARG	N-CA-C	-5.50	101.76	109.69
1	H	178	THR	N-CA-C	5.45	116.90	111.07
1	H	61	GLU	N-CA-C	-5.43	105.27	111.14
1	H	186	LYS	N-CA-C	-5.38	99.75	108.52
3	L	89	GLU	CA-C-N	5.35	126.53	119.84
3	L	89	GLU	C-N-CA	5.35	126.53	119.84
1	I	75	ARG	N-CA-C	-5.35	105.56	111.71
1	H	75	ARG	N-CA-C	-5.34	105.54	111.36
1	H	94	THR	N-CA-C	5.32	117.76	108.76
5	B	39	LEU	N-CA-C	5.32	119.25	112.34
1	I	113	TYR	N-CA-C	5.24	116.93	108.34
5	B	13	VAL	N-CA-C	5.16	120.03	108.88
1	I	39	ASP	N-CA-C	5.12	118.68	112.23
5	B	70	VAL	N-CA-C	5.08	116.53	111.00
5	B	65	TYR	N-CA-C	5.08	117.39	109.52
5	E	105	GLU	N-CA-C	5.08	116.81	111.28
1	I	275	GLU	N-CA-C	5.04	117.56	109.15
1	I	28	VAL	N-CA-C	-5.03	99.29	107.28
3	M	51	MET	N-CA-C	5.03	119.61	112.93

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	2247	0	2136	88	0
1	I	2254	0	2143	102	0
2	P	67	0	77	4	0
2	Q	67	0	77	6	0
3	L	821	0	796	34	0
3	M	821	0	796	44	0
4	A	881	0	844	56	0
4	D	881	0	845	75	0
5	B	940	0	936	74	0
5	E	940	0	936	82	0
6	C	106	0	89	7	0
7	B	14	0	13	0	0
7	E	14	0	13	0	0
8	A	5	0	0	1	0
8	B	10	0	0	1	0
8	D	6	0	0	2	0
8	E	7	0	0	4	0
8	H	22	0	0	1	0
8	I	18	0	0	3	0
8	L	12	0	0	0	0
8	M	6	0	0	0	0
8	P	2	0	0	0	0
8	Q	3	0	0	0	0
All	All	10144	0	9701	526	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 27.

All (526) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:108:THR:HA	5:E:42:GLN:HE22	1.22	1.02
3:M:37:ILE:O	3:M:51:MET:HE1	1.61	1.00
4:D:13:TRP:HB2	4:D:16:GLU:HG3	1.46	0.96
1:I:138:MET:HA	1:I:141:LEU:HD12	1.45	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:36:GLN:HB3	5:E:46:LEU:HD11	1.49	0.93
3:M:38:GLN:NE2	3:M:45:LYS:HD3	1.85	0.90
1:I:121:CYS:SG	3:M:1:ILE:HG13	2.13	0.88
4:A:5:GLN:HE21	4:A:107:GLY:HA3	1.37	0.87
1:H:157:ARG:O	1:H:161:GLU:HG3	1.74	0.87
1:H:234:ARG:HH11	3:L:8:GLN:NE2	1.73	0.86
1:I:138:MET:HA	1:I:141:LEU:CD1	2.04	0.86
4:D:61:ARG:HG2	4:D:78:THR:O	1.76	0.86
1:H:144:LYS:O	1:H:148:GLU:HG3	1.74	0.86
4:D:28:THR:HG22	4:D:94:TYR:HB2	1.58	0.86
1:I:194:ARG:HG3	1:I:195:PRO:HD2	1.55	0.85
4:A:47:ILE:HG23	4:A:64:ILE:HD11	1.57	0.85
4:A:13:TRP:O	4:A:16:GLU:HB2	1.77	0.85
1:H:218:GLN:HE21	1:H:223:GLU:HG2	1.42	0.84
5:B:15:ARG:H	5:B:116(A):LEU:HD23	1.39	0.84
5:E:1:VAL:HG12	5:E:2:THR:H	1.39	0.84
5:B:57:VAL:HG12	5:B:66:LEU:HD13	1.57	0.84
1:I:234:ARG:HH11	3:M:8:GLN:NE2	1.76	0.83
5:B:8:PRO:HG3	5:B:11:ARG:HD3	1.57	0.82
5:B:86:ARG:HH11	5:B:86:ARG:HG2	1.43	0.81
3:M:83:LYS:HG2	3:M:90:PRO:HG3	1.64	0.80
1:I:64:THR:O	1:I:68:LYS:HG2	1.83	0.79
3:L:48:LYS:NZ	3:L:69:GLU:H	1.80	0.79
5:B:15:ARG:H	5:B:116(A):LEU:CD2	1.94	0.79
4:D:28:THR:HG22	4:D:94:TYR:CB	2.11	0.79
4:A:30:ASN:HD22	4:A:30:ASN:H	1.29	0.78
1:I:263:HIS:CD2	1:I:265:GLY:H	2.02	0.77
1:H:234:ARG:HH11	3:L:8:GLN:HE22	1.31	0.77
4:A:7:PRO:HD3	6:C:1:NAG:H82	1.69	0.74
4:D:30:ASN:HD22	4:D:31:TYR:N	1.86	0.74
5:B:30(A):PRO:HB3	5:B:50:ARG:NH2	2.01	0.74
5:E:32:MET:HE2	5:E:69:ARG:CZ	2.16	0.74
5:B:31:TRP:CZ3	5:B:50:ARG:HB2	2.22	0.74
5:B:21:LEU:HD11	5:B:112:THR:HG21	1.70	0.73
4:D:61:ARG:HD2	4:D:79:ASP:HB3	1.70	0.73
4:D:4:ARG:CB	4:D:4:ARG:HH11	2.01	0.72
5:E:1:VAL:HG12	5:E:2:THR:N	2.03	0.72
1:I:263:HIS:HD2	1:I:265:GLY:H	1.36	0.72
5:E:8:PRO:HG3	5:E:11:ARG:NH1	2.05	0.72
1:H:160:LEU:O	1:H:165:VAL:HG23	1.89	0.72
5:B:14:PRO:HA	5:B:116(A):LEU:CD2	2.19	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:234:ARG:HH11	3:M:8:GLN:HE22	1.37	0.71
1:H:218:GLN:HG3	1:H:260:HIS:HD2	1.56	0.71
6:C:3:MAN:H61	6:C:4:MAN:H3	1.72	0.71
1:H:13:SER:HB3	1:H:78:LEU:HD13	1.73	0.70
1:I:114:GLN:HB2	1:I:156:LEU:HD21	1.73	0.70
1:I:29:ASP:O	1:I:30:ASP:HB2	1.90	0.70
1:H:218:GLN:HG3	1:H:260:HIS:CD2	2.27	0.70
5:B:32:MET:HE3	5:B:69:ARG:CZ	2.22	0.70
5:E:82:MET:HE1	5:E:114:LEU:HD12	1.74	0.70
5:B:36:GLN:HB3	5:B:46:LEU:HD11	1.72	0.69
4:A:5:GLN:HE21	4:A:107:GLY:CA	2.05	0.69
4:D:30:ASN:HD22	4:D:30:ASN:C	2.00	0.69
5:E:30:TYR:OH	5:E:103:SER:HB2	1.92	0.69
5:B:70:VAL:HG21	5:B:76:ARG:HD2	1.73	0.69
5:E:90:LEU:HD12	5:E:90:LEU:N	2.07	0.69
5:B:11:ARG:HG2	5:B:11:ARG:HH11	1.58	0.69
3:L:7:ILE:HD12	3:L:91:LYS:HD2	1.74	0.69
4:A:20:LEU:N	4:A:20:LEU:HD23	2.08	0.68
5:E:34:TRP:CZ3	5:E:92:CYS:HB2	2.28	0.68
1:H:103:VAL:HG12	1:H:109:LEU:HA	1.75	0.68
1:H:194:ARG:CD	1:H:195:PRO:HD2	2.23	0.68
4:D:28:THR:CG2	4:D:94:TYR:HB2	2.23	0.68
1:H:108:ARG:HH21	1:I:216:THR:HG21	1.58	0.68
1:I:138:MET:CA	1:I:141:LEU:HD12	2.24	0.68
1:I:194:ARG:CG	1:I:195:PRO:HD2	2.24	0.67
5:E:36:GLN:HB3	5:E:46:LEU:CD1	2.23	0.67
3:M:38:GLN:HE22	3:M:45:LYS:HD3	1.58	0.67
4:D:72:LYS:HD3	4:D:72:LYS:C	2.19	0.67
1:I:160:LEU:O	1:I:165:VAL:HG23	1.93	0.67
5:E:78:GLN:O	5:E:78:GLN:HG3	1.96	0.66
1:H:15:PRO:HB3	1:H:90:GLY:O	1.95	0.66
5:B:82:MET:HE2	5:B:86:ARG:HD2	1.78	0.66
4:A:5:GLN:HE22	4:A:90:CYS:H	1.42	0.66
5:B:82:MET:HE1	5:B:114:LEU:CB	2.26	0.66
4:D:5:GLN:HE22	4:D:90:CYS:H	1.44	0.66
4:D:47:ILE:HD13	4:D:64:ILE:HG13	1.77	0.66
5:E:112:THR:HG22	5:E:114:LEU:HD21	1.76	0.66
5:E:86:ARG:HG2	5:E:86:ARG:HH11	1.59	0.65
4:D:45:LEU:HD22	5:E:105(A):THR:CG2	2.27	0.65
4:D:108:THR:HA	5:E:42:GLN:NE2	2.05	0.65
4:D:47:ILE:HG23	4:D:64:ILE:HD11	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:38:PHE:CE1	4:D:86:ALA:HB2	2.31	0.65
5:B:24:ILE:CD1	5:B:74:GLU:HG3	2.27	0.65
4:D:14:GLU:O	4:D:16:GLU:HG2	1.97	0.64
5:B:22:ARG:NH2	5:B:74:GLU:HG2	2.12	0.64
5:B:14:PRO:HA	5:B:116(A):LEU:HD22	1.80	0.64
1:I:157:ARG:O	1:I:161:GLU:HG3	1.98	0.64
5:E:10:TRP:CZ2	5:E:12:LEU:HG	2.31	0.64
4:D:18:ALA:HB3	4:D:77:ILE:HB	1.80	0.64
4:D:38:PHE:CD1	4:D:86:ALA:HB2	2.33	0.64
1:H:50:ARG:HA	1:H:53:GLU:OE2	1.98	0.64
5:E:14:PRO:HA	5:E:116(A):LEU:HD22	1.79	0.64
5:B:36:GLN:NE2	5:B:65:TYR:OH	2.31	0.63
5:E:15:ARG:H	5:E:116(A):LEU:HD23	1.62	0.63
4:D:61:ARG:HD2	4:D:79:ASP:O	1.98	0.63
6:C:2:NAG:H83	6:C:5:NAG:H83	1.79	0.63
5:B:31:TRP:O	5:B:94:CYS:HA	1.99	0.63
1:I:135:ALA:HB1	1:I:140:ALA:HB3	1.81	0.62
3:L:25:CYS:HB2	3:L:39:MET:HE3	1.81	0.62
3:L:36:GLU:HB3	3:L:83:LYS:HB2	1.81	0.62
1:H:117:ALA:HB2	3:L:60:TRP:CE2	2.34	0.62
4:D:37:GLN:O	4:D:86:ALA:HB1	1.98	0.62
1:I:167:TRP:CE2	2:Q:1:LYS:HD2	2.33	0.62
1:H:108:ARG:HH11	1:H:108:ARG:HG2	1.64	0.62
5:E:8:PRO:HG3	5:E:11:ARG:HH11	1.63	0.61
4:A:28:THR:HG22	4:A:94:TYR:CB	2.31	0.61
4:D:13:TRP:HB2	4:D:16:GLU:CG	2.25	0.61
1:I:121:CYS:SG	3:M:1:ILE:HG21	2.40	0.61
3:M:7:ILE:HD12	3:M:91:LYS:HD2	1.81	0.61
4:D:4:ARG:HH11	4:D:4:ARG:HB3	1.64	0.61
5:E:50:ARG:HD3	5:E:50:ARG:C	2.25	0.61
4:D:45:LEU:HD22	5:E:105(A):THR:HG21	1.83	0.61
1:I:24:GLU:OE1	2:Q:2:VAL:HG11	2.01	0.60
1:I:224:LEU:CD2	1:I:247:VAL:HG21	2.31	0.60
1:H:194:ARG:CG	1:H:195:PRO:HD2	2.32	0.60
4:A:5:GLN:NE2	4:A:107:GLY:HA3	2.15	0.60
1:I:166:GLU:HG2	1:I:169:ARG:HH12	1.66	0.60
5:B:24:ILE:HD12	5:B:74:GLU:HG3	1.82	0.60
1:H:259:CYS:HB3	1:H:272:LEU:HB2	1.84	0.59
5:B:55:LYS:HB3	5:B:68:THR:HG22	1.84	0.59
1:I:259:CYS:HB3	1:I:272:LEU:HB2	1.83	0.59
3:M:88:ALA:HB3	3:M:89:GLU:OE2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:1:VAL:CG1	5:E:2:THR:H	2.14	0.59
5:B:82:MET:HE1	5:B:114:LEU:HB2	1.84	0.59
1:H:54:GLN:NE2	1:H:174:ASN:HB3	2.18	0.59
5:B:86:ARG:HH11	5:B:86:ARG:CG	2.13	0.59
3:M:1:ILE:O	3:M:1:ILE:HG22	2.02	0.59
5:E:86:ARG:HH11	5:E:86:ARG:CG	2.15	0.59
3:L:17:ASN:HA	3:L:72:PRO:O	2.02	0.59
4:D:2:GLN:OE1	4:D:25:GLU:HB3	2.02	0.59
4:D:36:GLN:HG3	4:D:88:TYR:CE1	2.37	0.59
4:D:77:ILE:HG22	8:D:122:HOH:O	2.01	0.59
1:H:16:GLY:C	1:H:18:GLY:H	2.11	0.59
5:E:95:SER:HB2	5:E:100:TRP:O	2.03	0.59
5:B:77:LEU:HD23	5:B:90:LEU:HD23	1.84	0.59
1:I:111:ARG:HD2	8:I:285:HOH:O	2.03	0.59
5:E:55:LYS:NZ	5:E:55:LYS:HB3	2.17	0.59
5:B:32:MET:CE	5:B:69:ARG:CZ	2.81	0.58
3:M:31:HIS:ND1	3:M:32:PRO:HA	2.18	0.58
1:H:47:PRO:HB3	1:H:60:TRP:CH2	2.38	0.58
3:L:96:ASP:HB3	3:L:99:MET:HG3	1.84	0.58
4:A:12:VAL:CG1	4:A:114:VAL:HG22	2.34	0.58
5:B:15:ARG:N	5:B:116(A):LEU:HD23	2.15	0.58
4:D:20:LEU:HD12	4:D:20:LEU:N	2.19	0.58
3:L:1:ILE:HG22	3:L:1:ILE:O	2.04	0.58
4:D:5:GLN:HE21	4:D:107:GLY:HA3	1.68	0.58
5:E:25:LEU:O	5:E:73:THR:HG22	2.02	0.58
4:D:3:VAL:CG1	4:D:105:ILE:HG22	2.34	0.58
3:M:5:PRO:HB3	3:M:30:PHE:HB3	1.86	0.57
4:D:26:ASP:OD2	4:D:28:THR:HB	2.04	0.57
4:A:10:LEU:HD12	4:A:11:THR:N	2.20	0.57
4:A:81:GLN:HG2	4:A:82:PRO:HD2	1.85	0.57
4:A:19:ILE:C	4:A:20:LEU:HD23	2.28	0.57
1:I:4:SER:HB2	1:I:101:CYS:O	2.05	0.57
4:A:57:GLU:HB3	1:I:256:TYR:CE1	2.40	0.57
5:B:70:VAL:HG12	5:B:71:THR:HG23	1.85	0.57
1:I:126:LEU:HD22	1:I:156:LEU:HD22	1.86	0.57
5:E:15:ARG:H	5:E:116(A):LEU:CD2	2.17	0.57
5:E:82:MET:HE1	5:E:114:LEU:CB	2.35	0.57
5:E:113:ARG:C	5:E:114:LEU:HD23	2.29	0.57
1:I:126:LEU:HD11	1:I:130:LEU:HA	1.87	0.57
1:I:47:PRO:HB3	1:I:60:TRP:CH2	2.40	0.57
1:I:123:TYR:HD2	1:I:124:ILE:HG22	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:16:GLU:OE1	3:M:19:LYS:HD2	2.06	0.56
4:A:30:ASN:HD22	4:A:30:ASN:N	1.96	0.56
1:I:167:TRP:O	1:I:171:TYR:CD2	2.58	0.56
3:M:7:ILE:CD1	3:M:91:LYS:HD2	2.34	0.56
4:D:54:ASP:HA	4:D:66:PHE:HB3	1.86	0.56
1:H:108:ARG:HH11	1:H:108:ARG:CG	2.18	0.56
1:H:194:ARG:HG3	1:H:195:PRO:HD2	1.85	0.56
4:A:17:THR:OG1	4:A:78:THR:HA	2.05	0.56
5:E:34:TRP:NE1	5:E:77:LEU:HB2	2.20	0.56
1:H:103:VAL:HG22	1:H:168:LEU:CD2	2.35	0.56
4:D:5:GLN:NE2	4:D:90:CYS:H	2.04	0.56
5:E:22:ARG:NH2	5:E:74:GLU:OE2	2.38	0.56
1:H:271:THR:O	1:H:272:LEU:HD23	2.05	0.56
5:B:84:GLN:HG3	5:B:85:GLY:H	1.69	0.55
4:D:81:GLN:C	4:D:114:VAL:HG11	2.30	0.55
1:I:166:GLU:CG	1:I:169:ARG:HH12	2.18	0.55
4:D:80:SER:O	4:D:81:GLN:HG3	2.06	0.55
1:H:263:HIS:CD2	1:H:265:GLY:H	2.24	0.55
1:I:21:ARG:NE	1:I:23:MET:HE2	2.21	0.55
8:H:296:HOH:O	1:I:108:ARG:HD3	2.06	0.55
5:E:82:MET:HE1	5:E:114:LEU:CD1	2.36	0.55
1:H:181:ARG:HH11	1:H:181:ARG:HG2	1.72	0.54
1:H:218:GLN:HE21	1:H:223:GLU:CG	2.19	0.54
5:B:38:ASP:OD1	5:B:40:GLN:HG2	2.07	0.54
5:E:112:THR:HG22	5:E:114:LEU:CD2	2.37	0.54
5:B:37:GLN:HB3	5:B:87:THR:CG2	2.38	0.54
1:I:21:ARG:CZ	1:I:23:MET:HE2	2.38	0.54
5:B:14:PRO:HG2	5:B:17:GLN:HB2	1.90	0.54
5:B:32:MET:HE3	5:B:69:ARG:NE	2.21	0.54
4:D:51:SER:HB2	4:D:68:LYS:NZ	2.22	0.54
4:D:77:ILE:CG2	8:D:122:HOH:O	2.55	0.54
4:A:61:ARG:O	4:A:78:THR:HG22	2.08	0.54
3:M:73:THR:HG1	3:M:76:ASP:CG	2.15	0.54
1:H:99:SER:OG	1:H:114:GLN:NE2	2.41	0.54
5:B:1:VAL:HG22	5:B:107:TYR:OH	2.08	0.54
1:I:210:PRO:O	1:I:263:HIS:HE1	1.89	0.54
4:A:50:ARG:HH11	5:B:105:GLU:CD	2.15	0.54
3:L:48:LYS:HZ1	3:L:69:GLU:H	1.55	0.53
5:E:31:TRP:HD1	5:E:96:ALA:O	1.92	0.53
5:B:11:ARG:HG2	5:B:11:ARG:NH1	2.20	0.53
1:H:266:LEU:HD13	1:H:270:LEU:HG	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:28:THR:CG2	4:A:94:TYR:HB2	2.38	0.53
1:H:234:ARG:HE	1:H:242:GLN:HE21	1.56	0.52
1:I:182:THR:HG22	1:I:182:THR:O	2.08	0.52
4:D:13:TRP:CE3	4:D:16:GLU:HG3	2.45	0.52
5:B:59:SER:O	5:B:60:LEU:HD12	2.09	0.52
1:I:135:ALA:HB1	1:I:140:ALA:CB	2.39	0.52
4:A:105:ILE:N	4:A:105:ILE:HD12	2.24	0.52
1:I:127:ASN:HB3	8:I:293:HOH:O	2.08	0.52
1:I:271:THR:C	1:I:272:LEU:HD23	2.35	0.52
5:E:47:PHE:CE2	5:E:56:GLU:HB3	2.45	0.52
1:H:70:ASN:O	1:H:73:SER:HB3	2.10	0.52
5:E:4:LEU:HD22	5:E:23:CYS:SG	2.50	0.52
1:I:99:SER:HA	1:I:113:TYR:O	2.10	0.52
1:I:234:ARG:HE	1:I:242:GLN:HE21	1.58	0.52
5:E:32:MET:HE2	5:E:69:ARG:NH2	2.25	0.52
1:I:13:SER:HB3	1:I:78:LEU:HD13	1.92	0.52
5:E:14:PRO:O	5:E:16:GLY:N	2.43	0.52
1:H:251:LEU:HD12	1:H:252:GLY:N	2.25	0.52
5:B:19:VAL:HB	5:B:79:VAL:HG13	1.92	0.52
4:D:30:ASN:C	4:D:30:ASN:ND2	2.68	0.52
5:E:82:MET:CE	5:E:114:LEU:HD12	2.38	0.52
4:A:61:ARG:NH1	4:A:84:ASP:OD2	2.41	0.51
5:B:70:VAL:CG2	5:B:76:ARG:HD2	2.40	0.51
1:H:55:GLU:CD	1:H:170:ARG:HH21	2.18	0.51
1:I:78:LEU:HD23	1:I:95:ILE:CD1	2.41	0.51
5:B:66:LEU:HB3	5:B:78:GLN:HG2	1.91	0.51
1:I:114:GLN:NE2	1:I:114:GLN:HA	2.25	0.51
1:I:172:LEU:O	1:I:173:LYS:C	2.53	0.51
1:I:224:LEU:HB3	1:I:227:ASP:OD2	2.10	0.51
1:H:3:HIS:HB3	1:H:29:ASP:OD1	2.10	0.51
5:B:10:TRP:CH2	5:B:12:LEU:HD21	2.45	0.51
3:M:37:ILE:HB	3:M:51:MET:CE	2.41	0.51
3:M:51:MET:O	3:M:64:ILE:HD11	2.10	0.51
5:E:2:THR:O	5:E:26:LYS:HD2	2.11	0.51
3:M:7:ILE:HG22	3:M:8:GLN:N	2.25	0.51
4:D:61:ARG:HD2	4:D:79:ASP:CB	2.40	0.51
1:H:133:TRP:HB2	1:H:144:LYS:HD2	1.92	0.51
4:A:30:ASN:H	4:A:30:ASN:ND2	2.03	0.51
1:I:42:ASN:N	1:I:43:PRO:HD3	2.26	0.51
3:L:36:GLU:OE2	3:L:83:LYS:HD2	2.11	0.50
3:M:37:ILE:HB	3:M:51:MET:HE2	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:28:THR:HG22	4:D:94:TYR:HB3	1.93	0.50
4:A:108:THR:HA	5:B:42:GLN:HE21	1.75	0.50
5:E:30(A):PRO:HB2	5:E:50:ARG:CZ	2.41	0.50
1:I:124:ILE:HA	1:I:134:THR:O	2.11	0.50
1:I:274:TRP:N	8:I:286:HOH:O	2.43	0.50
3:M:49:VAL:HG23	3:M:49:VAL:O	2.10	0.50
4:A:28:THR:HG22	4:A:94:TYR:HB2	1.92	0.50
5:E:32:MET:CE	5:E:69:ARG:CZ	2.88	0.50
1:H:224:LEU:N	1:H:224:LEU:HD22	2.27	0.50
1:I:230:LEU:HD23	1:I:230:LEU:H	1.76	0.50
4:D:13:TRP:HE3	4:D:16:GLU:CB	2.25	0.50
4:D:15:GLY:HA2	4:D:79:ASP:C	2.36	0.50
3:L:40:LEU:HD23	3:L:45:LYS:HA	1.94	0.50
4:A:2:GLN:HB2	4:A:105:ILE:HG21	1.94	0.50
5:B:13:VAL:O	5:B:13:VAL:HG13	2.12	0.50
1:I:182:THR:HG21	1:I:265:GLY:HA2	1.93	0.50
4:D:7:PRO:O	4:D:110:THR:HG23	2.11	0.50
4:A:28:THR:HG22	4:A:94:TYR:HB3	1.93	0.49
1:I:19:GLU:HB3	1:I:75:ARG:NH2	2.25	0.49
5:E:21:LEU:N	5:E:21:LEU:HD22	2.27	0.49
1:H:244:TRP:CZ2	3:L:99:MET:HE2	2.46	0.49
1:H:273:ARG:HG3	1:H:273:ARG:HH11	1.78	0.49
5:B:82:MET:HE2	5:B:86:ARG:CG	2.42	0.49
4:D:70:GLU:O	4:D:71:LYS:HB2	2.12	0.49
1:H:176:ASN:C	1:H:176:ASN:OD1	2.55	0.49
4:A:54:ASP:HA	4:A:66:PHE:HB3	1.93	0.49
1:H:226:GLN:HG2	1:H:227:ASP:N	2.27	0.49
4:D:13:TRP:CB	4:D:16:GLU:HG3	2.31	0.49
5:B:2:THR:O	5:B:2:THR:HG22	2.12	0.49
1:I:75:ARG:HH11	1:I:75:ARG:HG3	1.78	0.49
4:D:57:GLU:HG2	4:D:63:THR:HG23	1.95	0.49
3:M:39:MET:HB2	3:M:49:VAL:HG11	1.95	0.49
1:H:185:PRO:HD3	1:H:263:HIS:CD2	2.48	0.49
1:H:234:ARG:HE	1:H:242:GLN:NE2	2.10	0.49
1:H:222:GLU:OE1	1:H:223:GLU:C	2.56	0.48
3:L:41:LYS:HG3	3:L:78:TYR:CE2	2.49	0.48
1:I:255:GLN:HA	1:I:255:GLN:OE1	2.13	0.48
4:D:13:TRP:HE3	4:D:16:GLU:HG3	1.77	0.48
1:H:29:ASP:O	1:H:30:ASP:HB2	2.13	0.48
1:H:111:ARG:NE	1:H:128:GLU:OE2	2.46	0.48
1:H:16:GLY:O	1:H:18:GLY:N	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:117:ALA:HB2	3:L:60:TRP:CZ2	2.48	0.48
1:H:224:LEU:HD22	1:H:224:LEU:H	1.78	0.48
3:L:48:LYS:O	3:L:48:LYS:HG2	2.13	0.48
4:A:66:PHE:HD1	4:A:73:LEU:HG	1.79	0.48
1:I:224:LEU:HD22	1:I:247:VAL:HG21	1.95	0.48
5:E:31:TRP:NE1	5:E:98:PRO:HD3	2.29	0.48
4:A:26:ASP:OD2	4:A:28:THR:HB	2.14	0.48
4:A:47:ILE:HD13	4:A:64:ILE:HG13	1.95	0.48
1:I:275:GLU:HA	1:I:276:PRO:HD2	1.71	0.48
4:D:31:TYR:CD2	4:D:93:ARG:CZ	2.96	0.48
4:D:31:TYR:HE2	5:E:105(A):THR:HB	1.78	0.48
4:A:73:LEU:HD12	4:A:73:LEU:N	2.29	0.48
1:I:3:HIS:ND1	1:I:29:ASP:OD2	2.36	0.48
5:E:30(A):PRO:CB	5:E:50:ARG:NH2	2.76	0.48
1:H:234:ARG:NH1	3:L:8:GLN:NE2	2.53	0.48
5:E:1:VAL:CG1	5:E:2:THR:N	2.74	0.48
5:E:29:GLN:HG3	8:E:403:HOH:O	2.13	0.48
6:C:2:NAG:C7	6:C:5:NAG:H83	2.44	0.48
1:I:121:CYS:HG	3:M:1:ILE:HG13	1.78	0.47
1:H:108:ARG:CG	1:H:108:ARG:NH1	2.76	0.47
4:D:50:ARG:NH1	5:E:105:GLU:OE2	2.47	0.47
5:E:82:MET:HE1	5:E:114:LEU:HB2	1.96	0.47
4:D:54:ASP:O	4:D:55:LYS:HB2	2.15	0.47
3:M:16:GLU:O	3:M:17:ASN:C	2.56	0.47
3:M:82:VAL:O	3:M:87:MET:HE1	2.15	0.47
1:I:224:LEU:N	1:I:224:LEU:HD12	2.29	0.47
5:E:21:LEU:CD1	5:E:112:THR:HG21	2.45	0.47
1:I:155:ARG:NH2	2:Q:6:ILE:HG23	2.30	0.47
4:A:24:TYR:O	4:A:71:LYS:HD2	2.15	0.47
1:I:5:LEU:O	1:I:100:GLY:HA3	2.15	0.47
1:I:121:CYS:SG	3:M:1:ILE:CG1	2.97	0.47
1:I:182:THR:CG2	1:I:265:GLY:HA2	2.45	0.47
5:E:38:ASP:OD2	5:E:42:GLN:HB2	2.15	0.47
6:C:2:NAG:C8	6:C:5:NAG:H83	2.45	0.47
1:I:223:GLU:C	1:I:224:LEU:HD12	2.40	0.47
1:H:268:GLU:HG3	1:H:269:PRO:HD2	1.96	0.46
4:D:13:TRP:HE3	4:D:16:GLU:CG	2.29	0.46
4:A:57:GLU:OE1	1:I:255:GLN:HB2	2.15	0.46
1:I:96:GLN:NE2	3:M:62:PHE:CZ	2.84	0.46
1:I:250:PRO:HB2	1:I:253:LYS:HB2	1.97	0.46
5:E:37:GLN:HB3	5:E:87:THR:HG23	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:55:LYS:HZ2	1:I:255:GLN:NE2	2.12	0.46
5:B:29:GLN:C	5:B:30(A):PRO:HD3	2.40	0.46
1:I:85:TYR:HB2	1:I:87:GLN:HG3	1.96	0.46
5:E:31:TRP:CZ3	5:E:50:ARG:HB2	2.50	0.46
5:E:40:GLN:NE2	8:E:408:HOH:O	2.49	0.46
5:E:57:VAL:HG12	5:E:66:LEU:HD13	1.98	0.46
1:H:103:VAL:CG2	1:H:168:LEU:HD23	2.46	0.46
2:P:6:ILE:HG13	2:P:7:ASP:N	2.29	0.46
1:I:145:HIS:NE2	1:I:149:GLN:NE2	2.63	0.46
4:D:5:GLN:HE21	4:D:107:GLY:CA	2.29	0.46
4:D:12:VAL:HG22	4:D:16:GLU:HB2	1.97	0.46
5:E:114:LEU:HD23	5:E:114:LEU:N	2.29	0.46
1:H:8:PHE:HB3	3:L:56:PHE:CE2	2.51	0.46
1:H:194:ARG:HG3	1:H:195:PRO:CD	2.45	0.46
4:D:49:ILE:HD11	4:D:53:SER:HB3	1.98	0.46
1:H:176:ASN:OD1	1:H:177:ALA:N	2.49	0.46
1:H:251:LEU:HD12	1:H:252:GLY:H	1.80	0.46
4:A:4:ARG:NH1	4:A:23:SER:OG	2.49	0.46
4:A:62:PHE:CD1	4:A:62:PHE:N	2.84	0.46
4:A:80:SER:HB3	4:A:114:VAL:HG21	1.97	0.46
5:B:13:VAL:CG1	5:B:116:VAL:HA	2.45	0.46
3:M:21:ASN:HB3	3:M:70:PHE:CE1	2.51	0.46
1:H:68:LYS:O	1:H:72:GLN:HG3	2.16	0.45
4:A:57:GLU:HG3	4:A:63:THR:HG23	1.99	0.45
1:I:187:ALA:HA	1:I:204:TRP:O	2.16	0.45
5:E:30:TYR:N	5:E:30(A):PRO:HD3	2.31	0.45
4:A:28:THR:HG23	8:A:209:HOH:O	2.16	0.45
4:A:108:THR:O	5:B:42:GLN:NE2	2.50	0.45
1:I:234:ARG:NH1	3:M:8:GLN:NE2	2.56	0.45
4:D:26:ASP:HB3	4:D:29:PHE:CE2	2.51	0.45
4:D:33:PRO:HG2	4:D:91:ALA:HB3	1.97	0.45
1:H:36:PHE:HB2	1:H:45:TYR:CD1	2.51	0.45
1:H:41:GLU:HG3	1:H:42:ASN:N	2.31	0.45
5:B:31:TRP:CE3	5:B:50:ARG:HB2	2.51	0.45
1:I:114:GLN:HA	1:I:114:GLN:HE21	1.82	0.45
1:H:121:CYS:SG	3:L:1:ILE:HG13	2.56	0.45
4:D:16:GLU:N	4:D:80:SER:H	2.14	0.45
5:E:31:TRP:O	5:E:94:CYS:HA	2.17	0.45
5:B:50:ARG:HD3	5:B:50:ARG:C	2.41	0.45
1:I:129:ASP:O	1:I:131:LYS:HG3	2.17	0.45
1:I:78:LEU:HD23	1:I:95:ILE:HD11	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:71:THR:HG1	5:B:74:GLU:CD	2.25	0.45
1:I:232:GLU:HA	1:I:232:GLU:OE1	2.17	0.45
5:E:36:GLN:NE2	5:E:65:TYR:OH	2.50	0.45
2:Q:3:ILE:HG12	2:Q:4:THR:H	1.81	0.45
4:A:108:THR:CA	5:B:42:GLN:HE21	2.30	0.45
4:D:3:VAL:HG12	4:D:105:ILE:HG22	1.99	0.45
5:B:32:MET:HB3	5:B:75:LEU:HD22	1.98	0.45
3:M:14:PRO:HA	3:M:15:PRO:HD2	1.76	0.45
5:E:93:THR:HA	8:E:407:HOH:O	2.17	0.45
1:H:175:GLY:O	1:H:176:ASN:C	2.60	0.44
1:I:16:GLY:C	1:I:18:GLY:H	2.26	0.44
1:I:263:HIS:HD2	1:I:265:GLY:N	2.09	0.44
5:E:86:ARG:HB3	5:E:86:ARG:NH1	2.32	0.44
5:B:59:SER:C	5:B:60:LEU:HD12	2.42	0.44
5:B:82:MET:HE2	5:B:86:ARG:CD	2.44	0.44
4:D:78:THR:O	4:D:79:ASP:C	2.60	0.44
1:H:181:ARG:HG2	1:H:181:ARG:NH1	2.32	0.44
1:I:194:ARG:CD	1:I:195:PRO:HD2	2.48	0.44
1:I:225:ILE:HG22	1:I:226:GLN:N	2.32	0.44
3:M:38:GLN:HE22	3:M:45:LYS:CD	2.28	0.44
5:B:84:GLN:HG3	5:B:85:GLY:N	2.32	0.44
1:I:230:LEU:HD23	1:I:230:LEU:N	2.33	0.44
4:D:36:GLN:HG3	4:D:88:TYR:HE1	1.81	0.44
1:H:16:GLY:C	1:H:18:GLY:N	2.76	0.44
4:A:3:VAL:HG23	4:A:105:ILE:CG2	2.47	0.44
4:A:65:PHE:O	4:A:73:LEU:HA	2.17	0.44
5:B:15:ARG:H	5:B:116(A):LEU:HD21	1.80	0.44
5:B:62:GLY:HA3	5:B:86:ARG:HH22	1.82	0.44
1:I:224:LEU:N	1:I:224:LEU:CD1	2.81	0.44
1:H:41:GLU:O	1:H:42:ASN:HB3	2.18	0.44
1:H:204:TRP:HZ2	3:L:98:ASP:O	2.01	0.44
1:H:235:PRO:HG2	3:L:65:LEU:HD22	1.99	0.44
4:A:11:THR:HA	4:A:113:SER:O	2.17	0.44
1:H:54:GLN:HE22	1:H:174:ASN:HB3	1.81	0.44
5:B:34:TRP:NE1	5:B:77:LEU:HB2	2.33	0.44
5:B:34:TRP:CE2	5:B:77:LEU:HB2	2.53	0.43
1:H:97:VAL:HA	1:H:115:GLN:O	2.18	0.43
3:L:31:HIS:ND1	3:L:32:PRO:HA	2.33	0.43
4:A:70:GLU:O	4:A:72:LYS:HG3	2.18	0.43
4:A:81:GLN:N	4:A:114:VAL:HG11	2.32	0.43
1:I:152:GLU:CD	2:Q:6:ILE:HD13	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:29:PHE:HB3	4:D:32:PHE:CZ	2.53	0.43
1:H:99:SER:HA	1:H:113:TYR:O	2.17	0.43
5:B:30(A):PRO:CB	5:B:50:ARG:NH2	2.77	0.43
1:I:97:VAL:HA	1:I:115:GLN:O	2.18	0.43
1:I:227:ASP:OD1	1:I:227:ASP:N	2.51	0.43
3:M:51:MET:HB2	3:M:51:MET:HE3	1.55	0.43
3:M:73:THR:OG1	3:M:76:ASP:OD2	2.34	0.43
5:B:31:TRP:CZ2	5:B:50:ARG:HG3	2.54	0.43
3:M:28:THR:HG22	3:M:63:TYR:HB2	2.00	0.43
3:M:73:THR:OG1	3:M:76:ASP:HB2	2.18	0.43
4:D:3:VAL:HG11	4:D:105:ILE:HG22	1.98	0.43
5:E:30:TYR:CD1	5:E:102:ALA:HB1	2.54	0.43
3:L:10:TYR:CD2	3:L:10:TYR:N	2.86	0.43
4:A:48:SER:O	4:A:56:LYS:HG3	2.18	0.43
5:B:86:ARG:CG	5:B:86:ARG:NH1	2.78	0.43
5:B:100:TRP:HZ2	8:B:302:HOH:O	2.01	0.43
1:I:219:LEU:O	1:I:220:ASN:C	2.60	0.43
1:H:147:TRP:CZ2	2:P:8:LEU:HD23	2.54	0.43
5:E:62:GLY:O	5:E:63:ALA:HB2	2.18	0.43
4:D:12:VAL:HG23	4:D:13:TRP:CE3	2.53	0.43
5:E:31:TRP:CE3	5:E:50:ARG:HB2	2.53	0.43
1:H:23:MET:HE3	1:H:23:MET:HB2	1.86	0.43
4:D:31:TYR:CE2	5:E:105(A):THR:HB	2.54	0.43
4:D:61:ARG:CG	4:D:78:THR:O	2.57	0.43
5:E:14:PRO:O	5:E:15:ARG:C	2.62	0.43
5:E:90:LEU:N	5:E:90:LEU:CD1	2.80	0.43
5:E:30:TYR:CE1	5:E:102:ALA:C	2.97	0.42
5:E:30:TYR:CG	5:E:102:ALA:HB1	2.53	0.42
3:L:6:GLN:O	3:L:27:VAL:HA	2.20	0.42
4:A:17:THR:O	4:A:17:THR:HG22	2.19	0.42
1:I:44:ARG:HH22	1:I:61:GLU:HG3	1.84	0.42
5:E:55:LYS:HB3	5:E:55:LYS:HZ2	1.83	0.42
1:I:45:TYR:O	1:I:60:TRP:CE3	2.72	0.42
3:M:51:MET:HG3	3:M:64:ILE:HD11	2.00	0.42
3:L:74:GLU:OE1	3:L:74:GLU:HA	2.20	0.42
4:D:38:PHE:HB2	4:D:41:GLU:OE2	2.19	0.42
5:E:30(A):PRO:CB	5:E:50:ARG:CZ	2.97	0.42
3:M:98:ASP:O	3:M:99:MET:HG3	2.20	0.42
1:H:63:GLU:OE2	2:P:1:LYS:HE2	2.19	0.42
2:Q:6:ILE:HG13	2:Q:7:ASP:N	2.35	0.42
1:H:81:LEU:C	1:H:83:GLY:N	2.75	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:219:LEU:O	1:H:219:LEU:HG	2.19	0.42
4:A:50:ARG:HD3	5:B:105:GLU:OE2	2.19	0.42
1:I:14:ARG:HA	1:I:92:SER:HB2	2.01	0.42
1:I:58:GLU:H	1:I:58:GLU:HG3	1.45	0.42
1:I:80:THR:CG2	1:I:84:TYR:CE1	3.03	0.42
1:I:271:THR:O	1:I:272:LEU:HD23	2.20	0.42
4:A:5:GLN:HE21	4:A:107:GLY:C	2.28	0.42
5:B:4:LEU:HD13	5:B:92:CYS:SG	2.60	0.42
1:I:241:PHE:CD1	1:I:241:PHE:N	2.87	0.42
3:M:5:PRO:HA	3:M:28:THR:O	2.19	0.42
3:M:22:ILE:CD1	3:M:69:GLU:HG3	2.50	0.42
4:A:37:GLN:HE21	4:A:89:PHE:HE1	1.68	0.42
3:M:84:HIS:O	3:M:86:SER:N	2.52	0.42
1:H:201:LEU:O	1:H:246:SER:HA	2.20	0.42
3:L:56:PHE:HA	3:L:62:PHE:HA	2.02	0.42
4:A:83:GLY:C	4:A:85:SER:N	2.76	0.42
5:B:13:VAL:HA	5:B:14:PRO:HD2	1.90	0.42
3:M:19:LYS:H	3:M:71:THR:HG23	1.84	0.42
4:D:15:GLY:C	4:D:79:ASP:HA	2.45	0.42
5:E:112:THR:CG2	5:E:114:LEU:HD21	2.48	0.42
1:I:226:GLN:HG3	1:I:227:ASP:OD1	2.20	0.41
4:D:31:TYR:CZ	5:E:105:GLU:HG2	2.55	0.41
5:E:31:TRP:CE2	5:E:98:PRO:HD3	2.55	0.41
1:H:230:LEU:HD23	1:H:230:LEU:O	2.20	0.41
4:A:80:SER:O	4:A:81:GLN:HG3	2.19	0.41
5:B:82:MET:HE1	5:B:114:LEU:HD12	2.01	0.41
3:M:30:PHE:O	3:M:31:HIS:HB2	2.20	0.41
5:E:79:VAL:O	5:E:79:VAL:HG22	2.19	0.41
1:H:181:ARG:O	1:H:181:ARG:HG3	2.19	0.41
5:B:34:TRP:O	5:B:46:LEU:HB2	2.20	0.41
5:B:82:MET:HE1	5:B:114:LEU:CD1	2.50	0.41
1:I:209:TYR:CD2	1:I:209:TYR:C	2.98	0.41
1:H:203:CYS:O	1:H:244:TRP:HB2	2.20	0.41
5:B:55:LYS:CB	5:B:68:THR:HG22	2.50	0.41
4:D:63:THR:HG22	4:D:65:PHE:CE2	2.55	0.41
4:D:89:PHE:HZ	5:E:41:LYS:HB3	1.85	0.41
1:H:228:MET:HE3	1:H:230:LEU:HD13	2.01	0.41
3:L:7:ILE:CD1	3:L:91:LYS:HD2	2.46	0.41
3:M:24:ASN:OD1	3:M:67:HIS:HB3	2.21	0.41
4:D:13:TRP:HE3	4:D:16:GLU:HB3	1.85	0.41
5:E:51:SER:HA	5:E:52:PRO:HD3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:3:HIS:ND1	1:H:29:ASP:OD2	2.29	0.41
3:L:5:PRO:HB3	3:L:30:PHE:HB3	2.03	0.41
5:B:37:GLN:HB3	5:B:87:THR:HG22	2.02	0.41
1:H:20:PRO:HD2	1:H:75:ARG:HH11	1.86	0.41
1:I:21:ARG:CD	1:I:23:MET:HE2	2.51	0.41
1:I:215:LEU:CD2	1:I:261:VAL:HG22	2.51	0.41
1:I:232:GLU:O	1:I:233:THR:C	2.61	0.41
5:E:10:TRP:HE1	5:E:115:THR:HG23	1.85	0.41
1:I:145:HIS:CE1	1:I:149:GLN:HE22	2.39	0.41
1:I:172:LEU:O	1:I:176:ASN:N	2.51	0.41
1:H:13:SER:O	1:H:92:SER:OG	2.25	0.40
1:H:103:VAL:HG22	1:H:168:LEU:HD23	2.03	0.40
3:L:15:PRO:CG	3:L:97:ARG:HG3	2.51	0.40
3:L:46:ILE:HA	3:L:47:PRO:HD3	1.90	0.40
4:D:79:ASP:O	4:D:80:SER:O	2.39	0.40
5:E:15:ARG:NH2	8:E:404:HOH:O	2.55	0.40
6:C:3:MAN:H61	6:C:4:MAN:C3	2.30	0.40
1:H:42:ASN:N	1:H:43:PRO:CD	2.84	0.40
1:H:202:ARG:NH1	3:L:99:MET:HE1	2.35	0.40
5:B:29:GLN:O	5:B:30(A):PRO:HD3	2.21	0.40
1:H:224:LEU:H	1:H:224:LEU:CD2	2.34	0.40
4:A:55:LYS:NZ	1:I:255:GLN:NE2	2.70	0.40
1:H:66:LYS:NZ	2:P:2:VAL:O	2.51	0.40
3:L:54:MET:HG2	3:L:62:PHE:HD2	1.87	0.40
3:L:87:MET:SD	3:L:91:LYS:HE2	2.61	0.40
4:A:83:GLY:C	4:A:85:SER:H	2.29	0.40
5:B:19:VAL:HG12	5:B:20:ASN:N	2.37	0.40
3:M:69:GLU:O	3:M:70:PHE:HB3	2.21	0.40
5:E:38:ASP:OD1	5:E:38:ASP:C	2.64	0.40
6:C:2:NAG:N2	6:C:5:NAG:H83	2.37	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	H	274/277 (99%)	247 (90%)	21 (8%)	6 (2%)	5	13
1	I	275/277 (99%)	248 (90%)	23 (8%)	4 (2%)	8	20
2	P	6/8 (75%)	4 (67%)	1 (17%)	1 (17%)	0	0
2	Q	6/8 (75%)	4 (67%)	1 (17%)	1 (17%)	0	0
3	L	97/99 (98%)	90 (93%)	5 (5%)	2 (2%)	5	13
3	M	97/99 (98%)	81 (84%)	10 (10%)	6 (6%)	1	1
4	A	109/111 (98%)	97 (89%)	11 (10%)	1 (1%)	14	33
4	D	109/111 (98%)	88 (81%)	19 (17%)	2 (2%)	6	16
5	B	115/117 (98%)	106 (92%)	7 (6%)	2 (2%)	7	17
5	E	115/117 (98%)	104 (90%)	10 (9%)	1 (1%)	14	33
All	All	1203/1224 (98%)	1069 (89%)	108 (9%)	26 (2%)	5	13

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	17	LEU
1	H	226	GLN
4	A	80	SER
1	I	225	ILE
3	M	53	ASP
4	D	80	SER
5	E	15	ARG
3	L	90	PRO
5	B	2	THR
1	I	16	GLY
1	I	91	GLY
3	M	85	ASP
1	H	176	ASN
1	H	220	ASN
3	L	58	LYS
5	B	15	ARG
2	Q	6	ILE
3	M	90	PRO
4	D	13	TRP
1	H	224	LEU
1	I	14	ARG
3	M	50	GLU

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Mol	Chain	Res	Type
3	M	55	SER
2	P	6	ILE
1	H	91	GLY
3	M	31	HIS

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	H	234/235 (100%)	222 (95%)	12 (5%)	21	46
1	I	235/235 (100%)	216 (92%)	19 (8%)	11	26
2	P	8/8 (100%)	8 (100%)	0	100	100
2	Q	8/8 (100%)	7 (88%)	1 (12%)	4	10
3	L	94/94 (100%)	88 (94%)	6 (6%)	16	36
3	M	94/94 (100%)	88 (94%)	6 (6%)	16	36
4	A	96/96 (100%)	84 (88%)	12 (12%)	4	10
4	D	96/96 (100%)	85 (88%)	11 (12%)	5	13
5	B	102/102 (100%)	89 (87%)	13 (13%)	4	10
5	E	102/102 (100%)	80 (78%)	22 (22%)	1	3
All	All	1069/1070 (100%)	967 (90%)	102 (10%)	8	19

All (102) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	21	ARG
1	H	45	TYR
1	H	108	ARG
1	H	144	LYS
1	H	156	LEU
1	H	181	ARG
1	H	196	GLU
1	H	222	GLU
1	H	224	LEU

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Mol	Chain	Res	Type
1	H	230	LEU
1	H	258	THR
1	H	272	LEU
3	L	4	THR
3	L	54	MET
3	L	64	ILE
3	L	75	THR
3	L	89	GLU
3	L	99	MET
4	A	17	THR
4	A	20	LEU
4	A	23	SER
4	A	28	THR
4	A	30	ASN
4	A	41	GLU
4	A	50	ARG
4	A	56	LYS
4	A	57	GLU
4	A	73	LEU
4	A	110	THR
4	A	115	SER
5	B	3	LEU
5	B	17	GLN
5	B	40	GLN
5	B	49	LEU
5	B	50	ARG
5	B	55	LYS
5	B	56	GLU
5	B	78	GLN
5	B	79	VAL
5	B	86	ARG
5	B	103	SER
5	B	116	VAL
5	B	116(A)	LEU
1	I	10	THR
1	I	17	LEU
1	I	44	ARG
1	I	58	GLU
1	I	65	GLN
1	I	75	ARG
1	I	88	SER
1	I	103	VAL

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Mol	Chain	Res	Type
1	I	108	ARG
1	I	141	LEU
1	I	149	GLN
1	I	156	LEU
1	I	182	THR
1	I	196	GLU
1	I	200	THR
1	I	224	LEU
1	I	227	ASP
1	I	258	THR
1	I	272	LEU
2	Q	4	THR
3	M	2	GLN
3	M	16	GLU
3	M	40	LEU
3	M	64	ILE
3	M	89	GLU
3	M	97	ARG
4	D	2	GLN
4	D	3	VAL
4	D	4	ARG
4	D	11	THR
4	D	13	TRP
4	D	20	LEU
4	D	27	SER
4	D	28	THR
4	D	30	ASN
4	D	56	LYS
4	D	72	LYS
5	E	5	GLU
5	E	9	ARG
5	E	11	ARG
5	E	20	ASN
5	E	42	GLN
5	E	43	LEU
5	E	46	LEU
5	E	49	LEU
5	E	50	ARG
5	E	55	LYS
5	E	56	GLU
5	E	70	VAL
5	E	72	ASP

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Mol	Chain	Res	Type
5	E	73	THR
5	E	78	GLN
5	E	79	VAL
5	E	86	ARG
5	E	87	THR
5	E	114	LEU
5	E	115	THR
5	E	116	VAL
5	E	116(A)	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	42	ASN
1	H	54	GLN
1	H	65	GLN
1	H	72	GLN
1	H	114	GLN
1	H	127	ASN
1	H	218	GLN
1	H	220	ASN
1	H	226	GLN
1	H	242	GLN
1	H	260	HIS
1	H	263	HIS
1	H	264	GLN
3	L	6	GLN
3	L	8	GLN
3	L	13	HIS
3	L	29	GLN
3	L	67	HIS
4	A	5	GLN
4	A	30	ASN
4	A	37	GLN
5	B	7	ASN
5	B	20	ASN
5	B	29	GLN
5	B	36	GLN
5	B	37	GLN
5	B	42	GLN
5	B	44	GLN
5	B	78	GLN

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Mol	Chain	Res	Type
1	I	96	GLN
1	I	114	GLN
1	I	115	GLN
1	I	127	ASN
1	I	149	GLN
1	I	242	GLN
1	I	255	GLN
1	I	260	HIS
1	I	263	HIS
3	M	8	GLN
3	M	38	GLN
4	D	5	GLN
4	D	30	ASN
4	D	37	GLN
5	E	36	GLN
5	E	37	GLN
5	E	42	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 ⓘ

8 monosaccharides are 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	$\# Z > 2$
6	NAG	C	1	6,4	14,14,15	0.65	0	17,19,21	0.79	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	C	2	6	14,14,15	0.74	1 (7%)	17,19,21	0.71	0
6	MAN	C	3	6	11,11,12	0.59	0	15,15,17	1.45	2 (13%)
6	MAN	C	4	6	11,11,12	0.77	0	15,15,17	0.63	0
6	NAG	C	5	6	14,14,15	0.65	0	17,19,21	1.20	1 (5%)
6	GAL	C	6	6	11,11,12	0.66	0	15,15,17	0.40	0
6	SIA	C	7	6	20,20,21	0.76	0	21,28,31	0.90	1 (4%)
6	MAN	C	8	6	11,11,12	0.60	0	15,15,17	1.37	2 (13%)

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
6	NAG	C	1	6,4	-	0/6/23/26	0/1/1/1
6	NAG	C	2	6	-	0/6/23/26	0/1/1/1
6	MAN	C	3	6	1/1/4/5	1/2/19/22	0/1/1/1
6	MAN	C	4	6	-	1/2/19/22	0/1/1/1
6	NAG	C	5	6	-	4/6/23/26	0/1/1/1
6	GAL	C	6	6	-	2/2/19/22	0/1/1/1
6	SIA	C	7	6	-	3/18/34/38	0/1/1/1
6	MAN	C	8	6	1/1/4/5	1/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	2	NAG	C1-C2	2.18	1.55	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	3	MAN	C1-C2-C3	4.62	116.37	109.64
6	C	8	MAN	C1-O5-C5	3.66	117.10	112.19
6	C	5	NAG	C4-C3-C2	3.24	115.76	111.02
6	C	8	MAN	C1-C2-C3	3.22	114.34	109.64
6	C	1	NAG	C2-N2-C7	-2.52	119.52	122.90
6	C	7	SIA	O1B-C1-C2	2.35	118.81	112.71
6	C	3	MAN	O3-C3-C2	-2.12	105.73	110.05

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	C	3	MAN	C1
6	C	8	MAN	C1

All (12) torsion outliers are listed below:

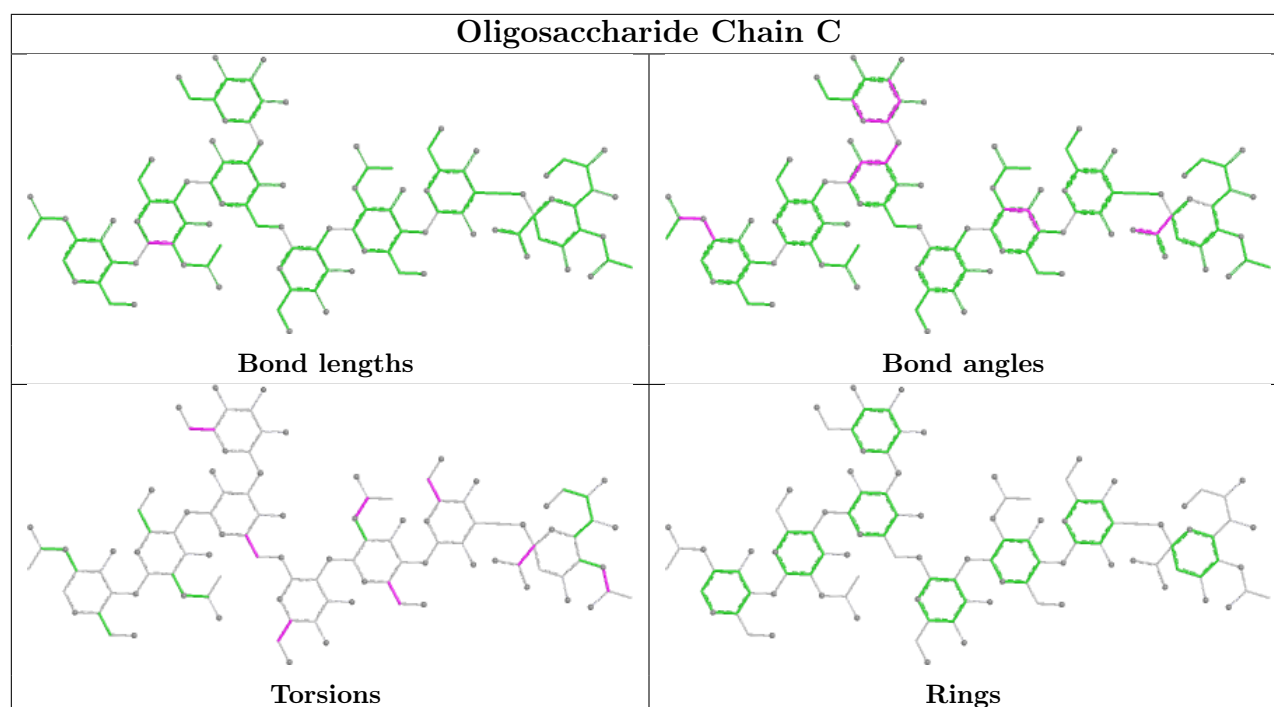
Mol	Chain	Res	Type	Atoms
6	C	7	SIA	C11-C10-N5-C5
6	C	7	SIA	O10-C10-N5-C5
6	C	5	NAG	O5-C5-C6-O6
6	C	5	NAG	C4-C5-C6-O6
6	C	6	GAL	O5-C5-C6-O6
6	C	5	NAG	C8-C7-N2-C2
6	C	6	GAL	C4-C5-C6-O6
6	C	5	NAG	O7-C7-N2-C2
6	C	8	MAN	O5-C5-C6-O6
6	C	4	MAN	O5-C5-C6-O6
6	C	3	MAN	O5-C5-C6-O6
6	C	7	SIA	O1A-C1-C2-O6

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	1	NAG	1	0
6	C	4	MAN	2	0
6	C	3	MAN	2	0
6	C	2	NAG	4	0
6	C	5	NAG	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

2 ligands are 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	$\# Z > 2$
7	NAG	B	301	5	14,14,15	0.63	0	17,19,21	0.86	1 (5%)
7	NAG	E	401	5	14,14,15	1.23	1 (7%)	17,19,21	1.09	2 (11%)

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
7	NAG	E	401	5	1/1/5/7	2/6/23/26	0/1/1/1
7	NAG	B	301	5	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	E	401	NAG	C1-C2	4.09	1.57	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	E	401	NAG	O5-C1-C2	2.63	115.36	111.29
7	B	301	NAG	C2-N2-C7	-2.59	119.42	122.90
7	E	401	NAG	C3-C4-C5	-2.25	106.15	110.23

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
7	E	401	NAG	C1

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	301	NAG	C8-C7-N2-C2
7	B	301	NAG	O7-C7-N2-C2
7	E	401	NAG	C8-C7-N2-C2
7	E	401	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	H	276/277 (99%)	-0.22	3 (1%) 78 76	13, 35, 75, 101	7 (2%)
1	I	277/277 (100%)	-0.05	11 (3%) 42 39	23, 40, 76, 102	10 (3%)
2	P	8/8 (100%)	-0.90	0 100 100	16, 19, 29, 31	0
2	Q	8/8 (100%)	-0.72	0 100 100	23, 29, 38, 42	0
3	L	99/99 (100%)	-0.31	0 100 100	22, 36, 56, 88	2 (2%)
3	M	99/99 (100%)	0.28	1 (1%) 79 78	26, 54, 86, 98	3 (3%)
4	A	111/111 (100%)	-0.06	1 (0%) 81 80	17, 43, 71, 82	1 (0%)
4	D	111/111 (100%)	0.43	5 (4%) 38 35	23, 58, 89, 104	1 (0%)
5	B	117/117 (100%)	-0.03	1 (0%) 81 80	20, 44, 77, 86	1 (0%)
5	E	117/117 (100%)	0.13	2 (1%) 69 66	23, 45, 81, 92	3 (2%)
All	All	1223/1224 (99%)	-0.03	24 (1%) 65 62	13, 42, 79, 104	28 (2%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	13	TRP	4.6
4	D	116	PRO	4.3
1	I	17	LEU	4.2
1	I	15	PRO	4.2
1	H	223	GLU	3.5
1	I	225	ILE	3.4
1	I	14	ARG	3.4
5	E	116(A)	LEU	3.2
1	I	256	TYR	3.1
3	M	49	VAL	3.0
5	B	10	TRP	2.8
1	I	91	GLY	2.7
1	I	194	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	I	16	GLY	2.6
5	E	10	TRP	2.5
4	D	80	SER	2.4
1	H	225	ILE	2.3
1	H	197	ASP	2.3
1	I	18	GLY	2.3
4	A	79	ASP	2.3
4	D	78	THR	2.2
1	I	197	ASP	2.2
4	D	10	LEU	2.2
1	I	221	GLY	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

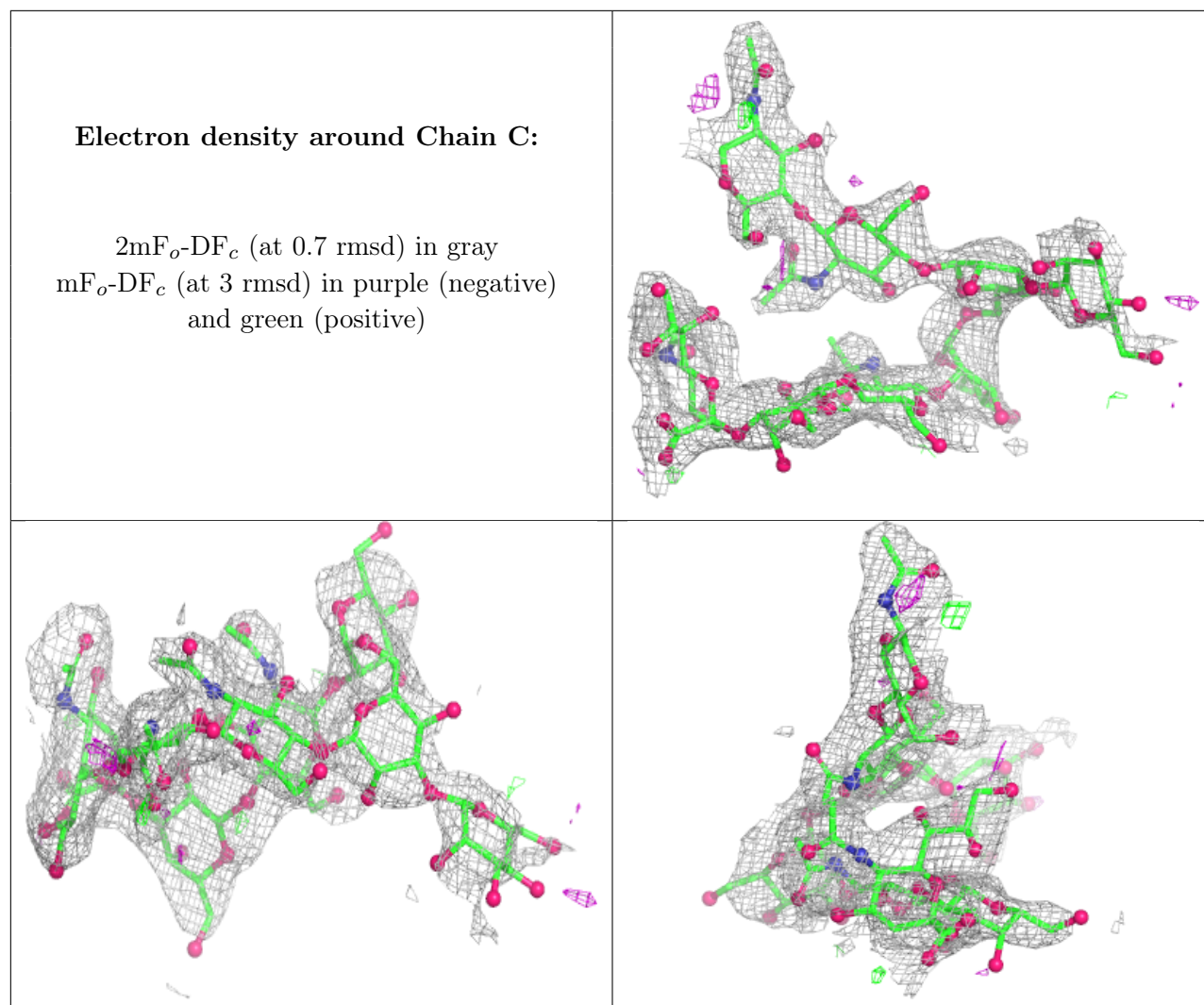
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	NAG	C	1	14/15	-	-	62,72,74,76	0
6	NAG	C	2	14/15	-	-	80,83,85,88	0
6	MAN	C	3	11/12	-	-	92,94,96,100	0
6	MAN	C	4	11/12	-	-	96,98,99,100	0
6	NAG	C	5	14/15	-	-	96,98,100,100	0
6	GAL	C	6	11/12	-	-	97,101,102,104	0
6	SIA	C	7	20/21	-	-	99,102,104,104	0
6	MAN	C	8	11/12	-	-	102,104,105,106	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	NAG	B	301	14/15	0.64	0.13	68,70,72,72	0
7	NAG	E	401	14/15	0.64	0.12	63,67,70,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.