



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:31 AM UTC

PDB ID : 5K9O / pdb_00005k9o
Title : Crystal structure of multidonor HV1-18+HD3-9 class broadly neutralizing Influenza A antibody 31.b.09 in complex with Hemagglutinin H1A/California/04/2009
Authors : Joyce, M.G.; Thomas, P.V.; Wheatley, A.K.; McDermott, A.B.; Mascola, J.R.; Kwong, P.D.
Deposited on : 2016-06-01
Resolution : 3.39 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

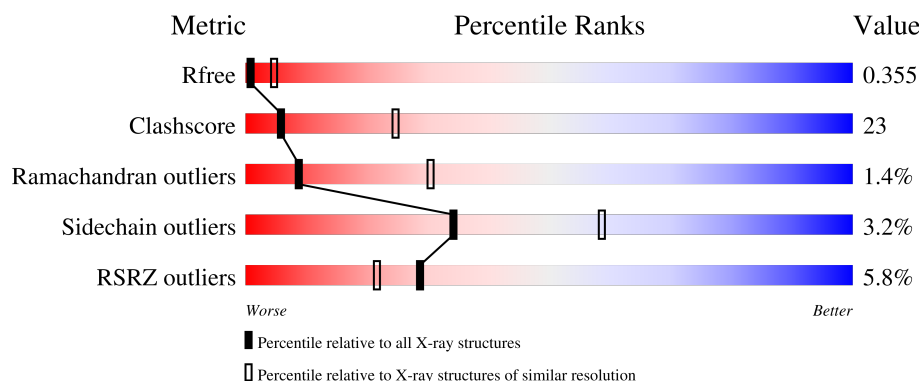
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.39 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))
R_{free}	180053	1047 (3.42-3.34)
Clashscore	190562	1067 (3.42-3.34)
Ramachandran outliers	187476	1056 (3.42-3.34)
Sidechain outliers	187428	1056 (3.42-3.34)
RSRZ outliers	180081	1047 (3.42-3.34)

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	A	227	7% 74% 20% • •
1	H	227	6% 71% 21% 5% •
2	B	219	7% 69% 23% 5% •
2	L	219	5% 74% 21% • •

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Mol	Chain	Length	Quality of chain
3	F	505	5% 78% 11% • • 8%
3	I	505	4% 75% 15% • 8%
4	C	3	33% 67%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14056 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 31.b.09 Heavy Fv.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1655	1043	277	327	8			
1	H	221	Total	C	N	O	S	0	0	0
			1661	1046	280	327	8			

- Molecule 2 is a protein called 31.b.09 Light Fv.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	0	0	0
			1654	1036	281	331	6			
2	L	217	Total	C	N	O	S	0	0	0
			1684	1055	285	337	7			

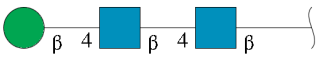
- Molecule 3 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	467	Total	C	N	O	S	0	0	0
			3685	2329	628	711	17			
3	I	466	Total	C	N	O	S	0	0	0
			3678	2317	629	715	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	506	SER	-	expression tag	UNP C3W5S1
F	507	GLY	-	expression tag	UNP C3W5S1
I	506	SER	-	expression tag	UNP C3W5S1
I	507	GLY	-	expression tag	UNP C3W5S1

- Molecule 4 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

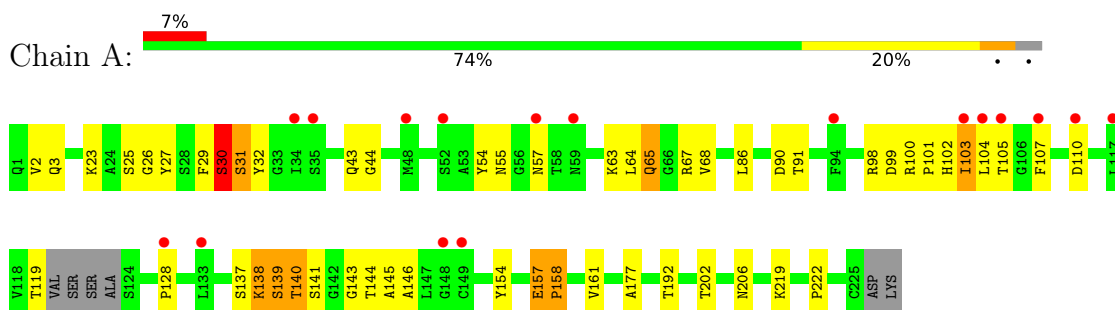


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	C	3	Total	C	N	O	0	0	0
			39	22	2	15			

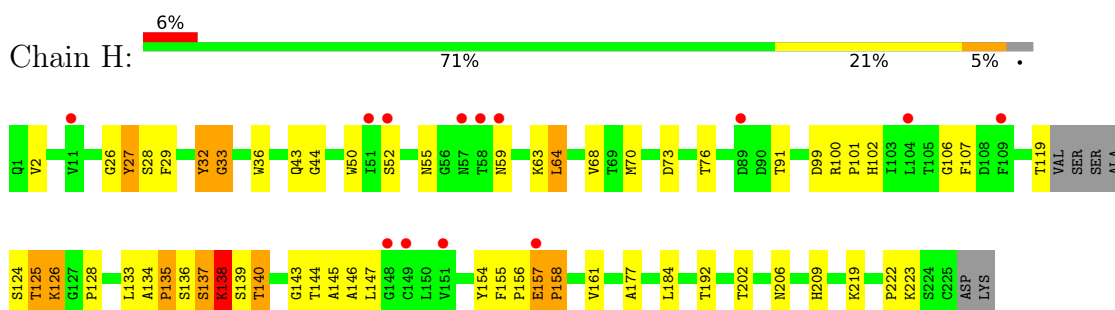
3 Residue-property plots [i](#)

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.

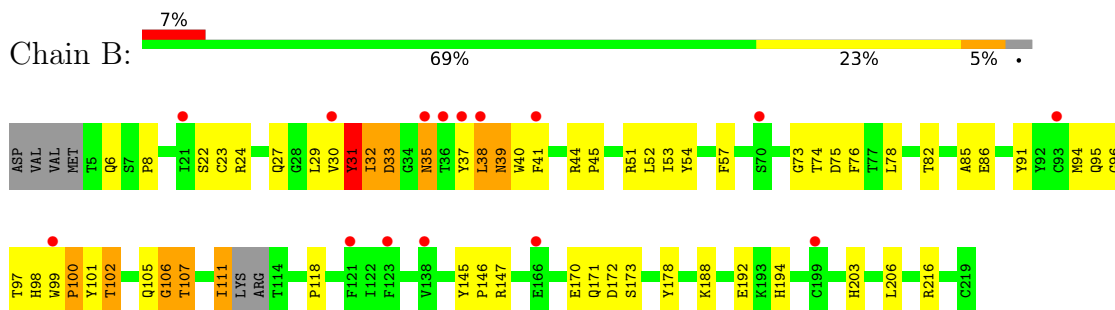
• Molecule 1: 31.b.09 Heavy Fv



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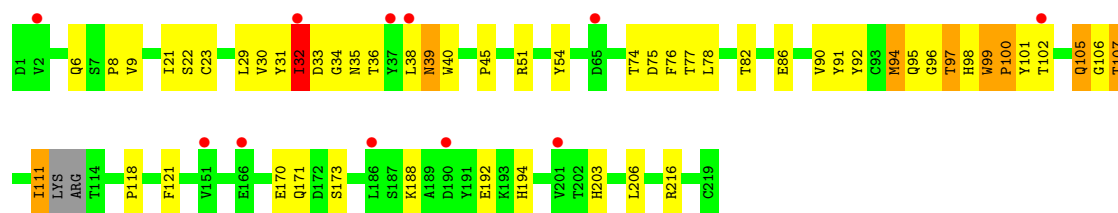


• Molecule 2: 31.b.09 Light Fv

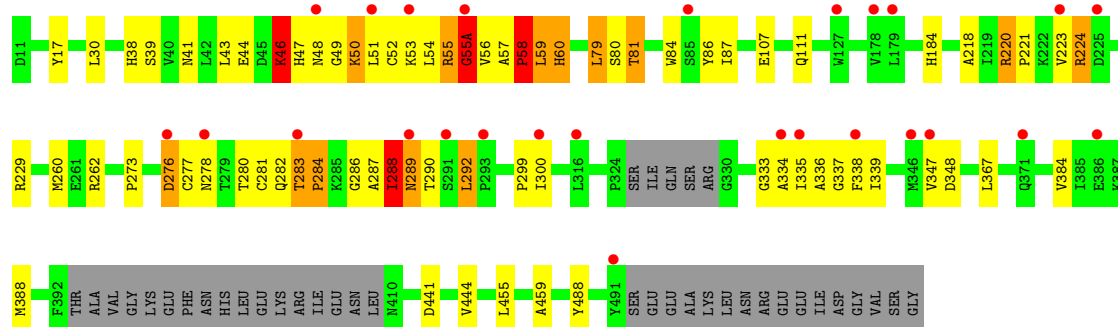
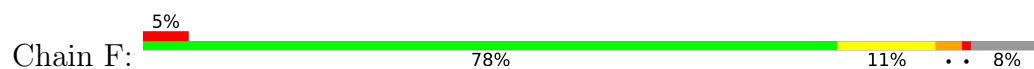


• Molecule 2: 31.b.09 Light Fv

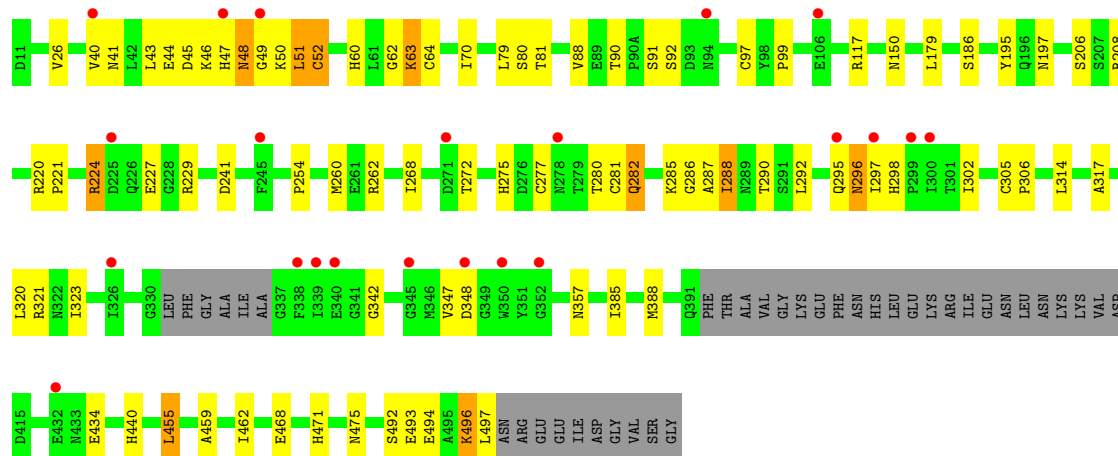
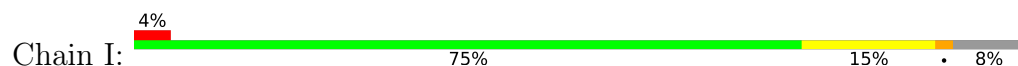




• Molecule 3: Hemagglutinin



• Molecule 3: Hemagglutinin



• Molecule 4: beta-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 6 2 2	Depositor
Cell constants a, b, c, α , β , γ	208.65Å 208.65Å 252.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.97 – 3.39 38.97 – 3.39	Depositor EDS
% Data completeness (in resolution range)	63.8 (38.97-3.39) 63.7 (38.97-3.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 3.40Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.291 , 0.346 0.306 , 0.355	Depositor DCC
R_{free} test set	1432 reflections (0.66%)	wwPDB-VP
Wilson B-factor (Å ²)	102.9	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 190.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	14056	wwPDB-VP
Average B, all atoms (Å ²)	207.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% of the height of the origin peak. No significant pseudotranslation is detected.*

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 [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.27	0/1695	0.57	3/2308 (0.1%)
1	H	0.28	0/1701	0.61	8/2315 (0.3%)
2	B	0.32	1/1692 (0.1%)	0.65	3/2299 (0.1%)
2	L	0.24	0/1722	0.55	0/2340
3	F	0.29	2/3774 (0.1%)	0.90	10/5117 (0.2%)
3	I	0.21	0/3765	0.46	1/5104 (0.0%)
All	All	0.27	3/14349 (0.0%)	0.66	25/19483 (0.1%)

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
3	F	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	55(A)	GLY	C-N	-7.93	1.22	1.33
3	F	220	ARG	C-N	5.36	1.40	1.33
2	B	100	PRO	N-CD	5.16	1.54	1.47

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	55(A)	GLY	CA-C-N	-35.35	58.34	121.97
3	F	55(A)	GLY	C-N-CA	-35.35	58.34	121.97
1	A	157	GLU	C-N-CD	8.68	139.68	120.60
1	H	157	GLU	C-N-CD	8.60	139.53	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	55(A)	GLY	O-C-N	7.79	132.83	122.70
3	I	280	THR	N-CA-C	7.68	122.36	112.92
1	H	134	ALA	C-N-CD	-7.49	94.30	125.00
3	F	220	ARG	CA-C-N	-6.46	113.12	119.78
3	F	220	ARG	C-N-CA	-6.46	113.12	119.78
2	B	39	ASN	N-CA-CB	-6.46	101.53	110.38
2	B	105	GLN	N-CA-C	-6.24	104.48	111.28
1	H	158	PRO	CA-N-CD	-5.83	103.33	111.50
1	H	27	TYR	CA-C-N	5.82	132.18	121.70
1	H	27	TYR	C-N-CA	5.82	132.18	121.70
1	H	33	GLY	N-CA-C	5.77	118.57	111.93
3	F	283	THR	CA-C-N	5.73	127.00	119.84
3	F	283	THR	C-N-CA	5.73	127.00	119.84
1	H	138	LYS	N-CA-C	-5.68	101.04	109.11
1	A	158	PRO	CA-N-CD	-5.67	103.56	111.50
3	F	46	LYS	N-CA-C	-5.60	102.46	110.59
3	F	57	ALA	CA-C-N	-5.38	113.12	119.84
3	F	57	ALA	C-N-CA	-5.38	113.12	119.84
2	B	106	GLY	N-CA-C	-5.35	105.94	112.68
1	A	30	SER	N-CA-C	-5.22	102.66	110.23
1	H	137	SER	CB-CA-C	-5.10	110.67	116.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	F	55(A)	GLY	Mainchain

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	1655	0	1611	77	0
1	H	1661	0	1621	120	0
2	B	1654	0	1584	121	0
2	L	1684	0	1621	117	0
3	F	3685	0	3557	144	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	I	3678	0	3546	106	0
4	C	39	0	34	0	0
All	All	14056	0	13574	637	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 23.

All (637) 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)
3:F:59:LEU:CD2	3:F:81:THR:HA	1.24	1.65
1:A:107:PHE:CZ	2:B:99:TRP:CH2	1.76	1.64
1:H:124:SER:CA	1:H:155:PHE:HE1	1.05	1.59
1:A:107:PHE:CE2	2:B:99:TRP:CZ3	1.90	1.57
1:H:124:SER:HA	1:H:155:PHE:CE1	0.98	1.51
1:A:107:PHE:CE2	2:B:99:TRP:CH2	1.93	1.50
2:B:57:PHE:HB2	3:F:289:ASN:ND2	1.19	1.47
1:H:124:SER:CA	1:H:155:PHE:CE1	1.84	1.43
3:F:54:LEU:O	3:F:56:VAL:CG2	1.71	1.38
2:L:8:PRO:O	2:L:107:THR:CG2	1.74	1.34
1:A:110:ASP:OD1	2:B:51:ARG:NH1	1.56	1.34
2:L:39:ASN:H	2:L:94:MET:CG	1.43	1.32
1:A:107:PHE:HE2	2:B:99:TRP:CZ3	1.34	1.31
1:A:102:HIS:NE2	3:F:347:VAL:O	1.66	1.29
2:B:57:PHE:CB	3:F:289:ASN:ND2	1.95	1.29
3:F:280:THR:OG1	3:F:288:ILE:HD11	1.17	1.29
3:I:51:LEU:CD2	3:I:272:THR:OG1	1.83	1.26
3:F:59:LEU:CD2	3:F:81:THR:CA	2.13	1.25
2:B:95:GLN:CG	2:B:101:TYR:HA	1.64	1.25
2:B:95:GLN:HG2	2:B:101:TYR:CA	1.51	1.24
2:L:39:ASN:N	2:L:94:MET:HG2	1.53	1.23
3:I:51:LEU:HD22	3:I:272:THR:OG1	1.34	1.22
2:L:91:TYR:O	2:L:106:GLY:HA2	1.36	1.21
2:B:57:PHE:CB	3:F:289:ASN:HD21	1.53	1.20
3:F:280:THR:OG1	3:F:288:ILE:CD1	1.90	1.19
1:A:141:SER:OG	1:A:144:THR:O	1.59	1.18
1:A:139:SER:O	1:A:141:SER:N	1.76	1.18
2:L:6:GLN:HE22	2:L:106:GLY:N	1.40	1.17
1:H:106:GLY:HA3	2:L:35:ASN:HD21	1.08	1.17
3:F:221:PRO:O	3:F:229:ARG:NH2	1.77	1.16
1:H:126:LYS:HE3	1:H:184:LEU:HD21	1.19	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:54:LEU:O	3:F:56:VAL:HG21	1.32	1.15
3:I:496:LYS:HB3	3:I:497:LEU:HB2	1.23	1.15
1:H:100:ARG:NH1	2:L:99:TRP:HH2	0.91	1.13
3:F:282:GLN:HB2	3:F:288:ILE:HG13	1.23	1.13
1:H:126:LYS:HD3	1:H:184:LEU:CD2	1.80	1.11
1:H:100:ARG:NH1	2:L:99:TRP:CH2	1.77	1.09
3:F:55(A):GLY:O	3:F:56:VAL:HG22	1.37	1.09
2:B:98:HIS:HB3	2:B:100:PRO:HD2	1.30	1.09
3:F:59:LEU:HD22	3:F:81:THR:CA	1.79	1.08
3:F:55(A):GLY:O	3:F:56:VAL:CG2	1.84	1.07
3:F:59:LEU:HD21	3:F:81:THR:HA	1.33	1.07
2:L:38:LEU:HB2	2:L:94:MET:HG3	1.37	1.06
1:H:124:SER:HA	1:H:155:PHE:CD1	1.89	1.06
1:A:30:SER:OG	1:A:101:PRO:HG3	1.58	1.04
2:B:39:ASN:HB3	2:B:54:TYR:HA	1.40	1.04
2:B:44:ARG:HH22	2:B:86:GLU:CD	1.66	1.04
1:A:107:PHE:HZ	2:B:99:TRP:CH2	1.28	1.03
3:F:48:ASN:HB3	3:F:51:LEU:HD13	1.34	1.03
1:H:36:TRP:CD1	1:H:70:MET:HE1	1.93	1.03
1:H:135:PRO:HG2	1:H:147:LEU:HB3	1.39	1.03
2:L:39:ASN:H	2:L:94:MET:HG2	0.87	1.02
3:I:224:ARG:HH21	3:I:224:ARG:HB2	1.22	1.02
2:B:44:ARG:NH2	2:B:86:GLU:OE2	1.91	1.01
1:A:107:PHE:CZ	2:B:99:TRP:HH2	1.36	1.01
1:H:106:GLY:CA	2:L:35:ASN:HD21	1.72	1.01
1:H:126:LYS:CE	1:H:184:LEU:HD21	1.90	1.00
2:B:37:TYR:HD1	2:B:76:PHE:CE2	1.74	1.00
2:L:9:VAL:C	2:L:107:THR:HG22	1.86	0.99
3:I:496:LYS:HG2	3:I:497:LEU:CD1	1.93	0.99
2:L:90:VAL:CG1	2:L:106:GLY:O	2.10	0.99
3:F:54:LEU:O	3:F:56:VAL:HG23	1.62	0.99
2:L:8:PRO:O	2:L:107:THR:HG21	0.82	0.98
1:A:107:PHE:HZ	2:B:99:TRP:CZ2	1.82	0.98
2:B:32:ILE:HG21	3:F:38:HIS:HB3	1.45	0.98
3:F:280:THR:HG1	3:F:288:ILE:HD11	1.23	0.98
3:F:287:ALA:HA	3:F:288:ILE:HG22	1.42	0.98
1:A:157:GLU:CG	1:A:158:PRO:HA	1.94	0.97
1:H:126:LYS:HE3	1:H:184:LEU:CD2	1.94	0.97
1:H:55:ASN:ND2	3:I:348:ASP:OD1	1.97	0.97
2:L:91:TYR:O	2:L:106:GLY:CA	2.13	0.97
1:H:36:TRP:CD1	1:H:70:MET:CE	2.47	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:90:VAL:HG13	2:L:106:GLY:O	1.64	0.97
3:I:496:LYS:HG2	3:I:497:LEU:HD12	1.46	0.97
2:B:37:TYR:CD1	2:B:76:PHE:CE2	2.39	0.96
3:I:496:LYS:HG3	3:I:497:LEU:HA	1.48	0.96
3:F:59:LEU:HD21	3:F:81:THR:CA	1.90	0.96
3:F:286:GLY:C	3:F:288:ILE:CG2	2.39	0.95
2:L:6:GLN:NE2	2:L:106:GLY:N	2.13	0.95
2:B:44:ARG:NH2	2:B:86:GLU:HG2	1.81	0.95
3:I:492:SER:OG	3:I:494:GLU:HG2	1.65	0.95
2:B:29:LEU:HB2	2:B:97:THR:HG21	1.46	0.95
1:H:124:SER:C	1:H:155:PHE:CE1	2.45	0.95
2:L:9:VAL:HA	2:L:107:THR:CG2	1.97	0.94
3:F:49:GLY:H	3:F:284:PRO:HG2	1.29	0.94
2:L:6:GLN:NE2	2:L:106:GLY:H	1.66	0.94
2:L:32:ILE:HG22	2:L:33:ASP:H	1.31	0.94
1:H:124:SER:CB	1:H:155:PHE:HE1	1.80	0.93
2:B:44:ARG:NH2	2:B:86:GLU:CG	2.30	0.93
1:H:126:LYS:HD3	1:H:184:LEU:HD22	1.50	0.93
2:B:95:GLN:HG2	2:B:101:TYR:HA	0.93	0.93
2:L:6:GLN:HE22	2:L:106:GLY:H	1.15	0.92
1:H:36:TRP:HD1	1:H:70:MET:HE1	1.33	0.92
3:I:51:LEU:HD22	3:I:272:THR:HG1	1.24	0.92
1:H:126:LYS:CD	1:H:184:LEU:CD2	2.47	0.91
2:B:98:HIS:HB3	2:B:100:PRO:CD	2.00	0.91
1:H:126:LYS:CE	1:H:184:LEU:CD2	2.48	0.91
3:F:55:ARG:C	3:F:56:VAL:HG23	1.95	0.91
2:B:75:ASP:O	2:B:76:PHE:CD1	2.23	0.90
3:F:288:ILE:H	3:F:290:THR:HG23	1.33	0.90
2:L:8:PRO:C	2:L:107:THR:HG21	1.94	0.90
3:F:218:ALA:O	3:F:220:ARG:NH2	2.03	0.90
1:H:140:THR:HG21	1:H:146:ALA:N	1.87	0.90
3:F:338:PHE:CD2	3:F:339:ILE:N	2.41	0.89
2:L:92:TYR:CE1	2:L:106:GLY:HA3	2.07	0.89
2:L:98:HIS:CD2	2:L:100:PRO:HD2	2.08	0.89
2:L:9:VAL:CA	2:L:107:THR:HG22	2.03	0.88
3:F:282:GLN:HB2	3:F:288:ILE:CG1	2.04	0.88
3:I:496:LYS:CG	3:I:497:LEU:HA	2.04	0.88
2:B:95:GLN:OE1	2:B:101:TYR:O	1.89	0.87
1:H:125:THR:N	1:H:155:PHE:HD1	1.71	0.87
1:H:126:LYS:HD3	1:H:155:PHE:HB3	1.57	0.87
1:H:126:LYS:CD	1:H:155:PHE:HB3	2.04	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:GLU:OE2	1:A:177:ALA:CB	2.22	0.86
3:I:224:ARG:HH21	3:I:224:ARG:CB	1.88	0.86
3:F:59:LEU:O	3:F:60:HIS:C	2.19	0.85
1:A:65:GLN:O	1:A:65:GLN:NE2	2.09	0.85
2:B:44:ARG:NH2	2:B:86:GLU:CD	2.32	0.85
1:A:139:SER:O	1:A:141:SER:O	1.93	0.85
3:I:496:LYS:HB3	3:I:497:LEU:CB	2.05	0.84
2:L:194:HIS:O	2:L:216:ARG:NH1	2.09	0.84
3:I:45:ASP:HB3	3:I:296:ASN:ND2	1.91	0.84
2:B:38:LEU:HB3	2:B:94:MET:CB	2.08	0.83
3:F:59:LEU:HD22	3:F:81:THR:HA	0.85	0.83
1:H:140:THR:CB	1:H:145:ALA:HA	2.08	0.83
1:H:136:SER:O	1:H:139:SER:N	2.12	0.83
1:A:107:PHE:CE2	2:B:99:TRP:HZ3	1.53	0.83
3:F:46:LYS:NZ	3:F:287:ALA:O	2.12	0.83
1:H:106:GLY:HA3	2:L:35:ASN:ND2	1.93	0.82
2:B:194:HIS:O	2:B:216:ARG:NH1	2.13	0.82
2:B:57:PHE:CG	3:F:289:ASN:ND2	2.47	0.82
1:A:157:GLU:OE2	1:A:177:ALA:HB1	1.80	0.82
2:B:57:PHE:HB2	3:F:289:ASN:HD21	0.70	0.82
1:H:124:SER:CA	1:H:155:PHE:CD1	2.57	0.82
3:I:97:CYS:O	3:I:224:ARG:HD2	1.79	0.82
1:H:126:LYS:CD	1:H:184:LEU:HD22	2.08	0.81
3:I:51:LEU:HD23	3:I:272:THR:OG1	1.78	0.81
1:H:106:GLY:CA	2:L:35:ASN:ND2	2.43	0.81
1:H:124:SER:C	1:H:155:PHE:CD1	2.58	0.81
1:H:140:THR:HG21	1:H:146:ALA:H	1.46	0.81
1:H:135:PRO:CG	1:H:147:LEU:HB3	2.09	0.81
1:H:136:SER:HB2	1:H:223:LYS:HB2	1.63	0.81
3:F:287:ALA:CA	3:F:288:ILE:HG22	2.11	0.81
3:F:58:PRO:HB3	3:F:84:TRP:CG	2.17	0.80
2:B:37:TYR:OH	2:B:74:THR:HA	1.80	0.80
3:I:320:LEU:HD12	3:I:440:HIS:HB3	1.63	0.79
2:B:8:PRO:O	2:B:107:THR:HG23	1.81	0.79
2:L:39:ASN:N	2:L:94:MET:CG	2.24	0.79
2:L:39:ASN:N	2:L:94:MET:SD	2.54	0.79
3:F:287:ALA:HB1	3:F:289:ASN:HB3	1.65	0.79
2:B:32:ILE:CG2	3:F:38:HIS:HB3	2.11	0.78
2:B:38:LEU:HB3	2:B:94:MET:HB3	1.66	0.78
1:A:54:TYR:N	1:A:55:ASN:HA	1.98	0.78
2:B:39:ASN:O	2:B:94:MET:HG3	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:9:VAL:HA	2:L:107:THR:HG22	1.62	0.78
2:L:32:ILE:HG22	2:L:33:ASP:N	1.98	0.78
3:F:286:GLY:C	3:F:288:ILE:HG22	2.08	0.78
3:I:357:ASN:ND2	3:I:475:ASN:OD1	2.11	0.78
2:L:95:GLN:HG3	2:L:101:TYR:HA	1.65	0.78
3:F:280:THR:CB	3:F:288:ILE:HD11	2.14	0.78
2:L:9:VAL:CA	2:L:107:THR:CG2	2.60	0.77
1:H:124:SER:C	1:H:155:PHE:HE1	1.86	0.77
3:F:58:PRO:CG	3:F:87:ILE:HA	2.15	0.77
1:H:124:SER:HA	1:H:155:PHE:CZ	2.06	0.77
1:A:157:GLU:HG2	1:A:158:PRO:HA	1.63	0.77
1:H:138:LYS:C	1:H:138:LYS:HD2	2.09	0.77
3:F:455:LEU:HD13	3:F:459:ALA:HB3	1.64	0.77
2:B:95:GLN:CG	2:B:101:TYR:CA	2.30	0.77
1:A:65:GLN:HE21	1:A:65:GLN:C	1.94	0.76
2:B:29:LEU:HB2	2:B:97:THR:CG2	2.16	0.76
3:F:55:ARG:O	3:F:56:VAL:N	2.17	0.76
3:F:48:ASN:OD1	3:F:51:LEU:HD21	1.85	0.76
2:L:92:TYR:CD1	2:L:106:GLY:HA3	2.20	0.76
1:H:125:THR:N	1:H:155:PHE:CD1	2.54	0.76
1:H:33:GLY:N	1:H:100:ARG:O	2.18	0.76
1:H:64:LEU:HD22	1:H:64:LEU:N	2.01	0.76
3:F:54:LEU:C	3:F:56:VAL:CG2	2.59	0.75
2:L:32:ILE:HD12	2:L:32:ILE:N	2.01	0.75
2:B:24:ARG:CB	2:B:75:ASP:OD2	2.34	0.75
3:I:47:HIS:CG	3:I:48:ASN:H	2.04	0.75
1:A:64:LEU:N	1:A:64:LEU:HD22	2.02	0.75
1:A:102:HIS:CE1	3:F:347:VAL:O	2.40	0.75
2:L:9:VAL:O	2:L:107:THR:HG22	1.86	0.75
3:F:58:PRO:HG3	3:F:87:ILE:HA	1.69	0.74
3:F:55:ARG:HB2	3:F:278:ASN:OD1	1.86	0.74
3:F:286:GLY:C	3:F:288:ILE:HG21	2.11	0.73
2:L:95:GLN:HB2	2:L:101:TYR:CD2	2.24	0.73
2:B:91:TYR:O	2:B:106:GLY:CA	2.36	0.73
1:H:140:THR:HB	1:H:145:ALA:HA	1.69	0.73
2:L:38:LEU:CB	2:L:94:MET:HG3	2.15	0.73
2:L:33:ASP:O	2:L:35:ASN:N	2.22	0.73
3:F:54:LEU:C	3:F:56:VAL:HG21	2.14	0.73
3:F:282:GLN:CB	3:F:288:ILE:HG13	2.14	0.73
1:A:30:SER:HB3	1:A:98:ARG:HH21	1.52	0.72
3:F:59:LEU:O	3:F:60:HIS:O	2.07	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:95:GLN:HG2	2:L:96:GLY:N	2.04	0.72
1:H:43:GLN:HG3	1:H:44:GLY:H	1.52	0.72
2:B:24:ARG:HB2	2:B:75:ASP:OD2	1.87	0.72
2:B:39:ASN:O	2:B:94:MET:CG	2.38	0.71
3:F:44:GLU:HG3	3:F:292:LEU:HD22	1.72	0.71
1:H:124:SER:O	1:H:125:THR:HG23	1.91	0.71
1:A:107:PHE:CE2	2:B:99:TRP:HH2	1.60	0.71
3:F:286:GLY:O	3:F:288:ILE:CG2	2.39	0.70
3:I:41:ASN:ND2	3:I:314:LEU:O	2.24	0.70
1:A:30:SER:OG	1:A:32:TYR:O	2.09	0.70
2:L:90:VAL:HG12	2:L:106:GLY:O	1.91	0.70
3:I:496:LYS:HG2	3:I:497:LEU:HD13	1.73	0.70
3:F:276:ASP:N	3:F:276:ASP:OD1	2.25	0.69
3:I:272:THR:HG21	3:I:285:LYS:HG3	1.74	0.69
3:F:48:ASN:CB	3:F:51:LEU:HD13	2.17	0.69
3:F:48:ASN:HB3	3:F:51:LEU:CD1	2.15	0.69
1:H:136:SER:CB	1:H:223:LYS:HB2	2.23	0.69
1:A:32:TYR:HB2	1:A:101:PRO:HB3	1.74	0.69
3:I:492:SER:CB	3:I:494:GLU:HG2	2.22	0.69
3:F:54:LEU:HB3	3:F:56:VAL:HG21	1.73	0.69
3:F:287:ALA:N	3:F:288:ILE:HB	2.07	0.69
3:F:287:ALA:C	3:F:289:ASN:HB3	2.18	0.69
2:L:98:HIS:HD2	2:L:99:TRP:H	1.39	0.69
3:F:287:ALA:CA	3:F:289:ASN:HB3	2.23	0.68
3:F:58:PRO:HB3	3:F:84:TRP:CB	2.23	0.68
2:L:92:TYR:CD1	2:L:106:GLY:CA	2.76	0.68
1:H:126:LYS:CD	1:H:184:LEU:HD21	2.18	0.68
2:L:9:VAL:HA	2:L:107:THR:HG23	1.75	0.68
1:H:63:LYS:HB2	1:H:64:LEU:HD22	1.76	0.68
3:I:296:ASN:N	3:I:296:ASN:HD22	1.91	0.68
1:H:135:PRO:HG2	1:H:147:LEU:CB	2.22	0.68
1:H:102:HIS:CD2	3:I:347:VAL:HG23	2.29	0.68
1:H:140:THR:HG22	2:L:121:PHE:CE2	2.30	0.68
1:A:139:SER:O	1:A:141:SER:C	2.37	0.67
3:I:45:ASP:HB3	3:I:296:ASN:HD21	1.59	0.67
3:I:455:LEU:HD13	3:I:459:ALA:HB3	1.76	0.67
2:B:38:LEU:HB3	2:B:94:MET:HB2	1.74	0.67
3:F:58:PRO:HG3	3:F:87:ILE:HG22	1.75	0.67
2:L:22:SER:HB2	2:L:76:PHE:CE1	2.29	0.67
2:B:91:TYR:O	2:B:106:GLY:HA2	1.94	0.67
2:L:98:HIS:CD2	2:L:99:TRP:H	2.13	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:50:LYS:HB3	3:I:275:HIS:HD2	1.59	0.67
2:B:98:HIS:O	2:B:101:TYR:CE1	2.49	0.66
2:L:38:LEU:HB2	2:L:94:MET:CG	2.19	0.66
2:B:98:HIS:O	2:B:101:TYR:CD1	2.48	0.66
3:F:79:LEU:HD12	3:F:81:THR:HB	1.77	0.66
1:H:126:LYS:O	1:H:209:HIS:CE1	2.49	0.66
1:A:107:PHE:CZ	2:B:99:TRP:CZ3	2.43	0.65
3:F:55:ARG:CG	3:F:55(A):GLY:H	2.10	0.65
2:L:99:TRP:HA	2:L:99:TRP:CE3	2.31	0.65
1:A:157:GLU:OE1	1:A:177:ALA:HB3	1.97	0.65
3:F:287:ALA:CB	3:F:289:ASN:HB3	2.26	0.65
3:I:282:GLN:NE2	3:I:282:GLN:HA	2.11	0.65
2:L:39:ASN:OD1	2:L:94:MET:SD	2.54	0.65
1:A:31:SER:HB2	1:A:32:TYR:CE2	2.31	0.65
3:I:492:SER:OG	3:I:494:GLU:CG	2.41	0.65
1:A:157:GLU:HG3	1:A:158:PRO:HA	1.77	0.65
1:A:157:GLU:CD	1:A:177:ALA:CB	2.70	0.65
2:B:98:HIS:CB	2:B:100:PRO:HD2	2.17	0.65
2:B:45:PRO:CG	2:B:170:GLU:HG3	2.26	0.65
3:F:49:GLY:N	3:F:284:PRO:HG2	2.09	0.65
2:L:95:GLN:HG3	2:L:101:TYR:CG	2.32	0.65
1:A:102:HIS:CE1	1:A:104:LEU:HB3	2.32	0.64
2:B:37:TYR:CD1	2:B:76:PHE:CD2	2.85	0.64
3:F:17:TYR:HE1	3:F:444:VAL:HG21	1.62	0.64
1:A:86:LEU:HD12	1:A:90:ASP:HB3	1.78	0.64
2:B:75:ASP:O	2:B:76:PHE:HD1	1.80	0.64
2:L:91:TYR:O	2:L:106:GLY:C	2.41	0.64
2:B:118:PRO:HD2	2:B:206:LEU:HD11	1.79	0.64
3:F:224:ARG:NH1	3:F:224:ARG:HG3	2.12	0.64
3:F:286:GLY:O	3:F:288:ILE:HG22	1.94	0.64
1:H:100:ARG:HD3	1:H:107:PHE:CD1	2.33	0.64
1:A:64:LEU:HD12	1:A:67:ARG:HH21	1.63	0.63
2:L:38:LEU:H	2:L:94:MET:CG	2.11	0.63
2:L:95:GLN:CG	2:L:101:TYR:HA	2.28	0.63
1:A:30:SER:HB3	1:A:98:ARG:NH2	2.13	0.63
3:F:17:TYR:CE1	3:F:444:VAL:HG21	2.33	0.63
3:I:90:THR:HG22	3:I:91:SER:H	1.63	0.63
3:F:59:LEU:HD21	3:F:81:THR:N	2.13	0.63
3:I:496:LYS:CB	3:I:497:LEU:HB2	2.16	0.63
2:B:6:GLN:HE22	2:B:106:GLY:HA2	1.64	0.62
3:F:55:ARG:HG2	3:F:55(A):GLY:H	1.62	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:497:LEU:HG	3:I:497:LEU:O	1.99	0.62
2:L:22:SER:HA	2:L:76:PHE:O	1.98	0.62
2:B:40:TRP:O	2:B:52:LEU:HD12	2.00	0.62
1:H:125:THR:HA	1:H:155:PHE:O	2.00	0.62
3:F:54:LEU:C	3:F:56:VAL:HG23	2.22	0.62
1:H:126:LYS:N	1:H:155:PHE:O	2.33	0.61
3:I:50:LYS:HB3	3:I:275:HIS:CD2	2.36	0.61
1:A:139:SER:C	1:A:141:SER:N	2.57	0.61
3:I:268:ILE:HD11	3:I:302:ILE:HD12	1.80	0.61
3:I:224:ARG:HB2	3:I:224:ARG:NH2	2.05	0.61
3:F:287:ALA:HA	3:F:288:ILE:CG2	2.25	0.61
1:H:157:GLU:HB3	1:H:158:PRO:HA	1.81	0.61
3:I:97:CYS:O	3:I:224:ARG:CD	2.49	0.61
3:I:52:CYS:SG	3:I:277:CYS:N	2.65	0.61
2:L:118:PRO:HD2	2:L:206:LEU:HD11	1.81	0.61
3:F:287:ALA:HB1	3:F:289:ASN:CB	2.30	0.61
1:A:43:GLN:HG3	1:A:44:GLY:H	1.65	0.60
2:L:38:LEU:CB	2:L:94:MET:HE2	2.31	0.60
1:H:33:GLY:H	1:H:100:ARG:C	2.09	0.60
1:H:126:LYS:CG	1:H:155:PHE:HB3	2.31	0.60
2:L:29:LEU:HD21	2:L:96:GLY:HA2	1.82	0.60
2:L:38:LEU:C	2:L:94:MET:HG2	2.27	0.60
2:B:74:THR:HG23	2:B:74:THR:O	2.01	0.60
3:F:46:LYS:HE3	3:F:283:THR:HG23	1.83	0.60
1:A:102:HIS:CD2	3:F:347:VAL:O	2.52	0.59
2:B:98:HIS:HB3	2:B:100:PRO:CG	2.32	0.59
1:A:31:SER:HB2	1:A:32:TYR:CZ	2.37	0.59
3:I:51:LEU:CD2	3:I:272:THR:HG1	1.91	0.59
2:B:203:HIS:HB3	2:B:206:LEU:HD13	1.83	0.59
1:A:29:PHE:CE2	1:A:31:SER:HA	2.37	0.59
3:I:50:LYS:HD3	3:I:275:HIS:NE2	2.17	0.59
2:B:32:ILE:HG21	3:F:38:HIS:CB	2.28	0.59
2:L:38:LEU:H	2:L:94:MET:HG3	1.67	0.59
2:B:27:GLN:NE2	2:B:98:HIS:NE2	2.50	0.59
3:F:59:LEU:CD2	3:F:81:THR:CB	2.80	0.59
3:F:184:HIS:HB3	3:F:220:ARG:NH1	2.17	0.58
2:B:40:TRP:O	2:B:52:LEU:HB2	2.03	0.58
3:F:335:ILE:O	3:F:336:ALA:HB3	2.03	0.58
1:H:156:PRO:O	1:H:209:HIS:NE2	2.31	0.58
3:F:224:ARG:HG3	3:F:224:ARG:HH11	1.68	0.58
1:H:140:THR:CG2	1:H:145:ALA:HA	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:51:LEU:HD13	3:I:286:GLY:H	1.68	0.58
1:A:107:PHE:HE1	2:B:31:TYR:CE1	2.21	0.58
2:L:35:ASN:O	2:L:36:THR:OG1	2.21	0.58
3:I:385:ILE:O	3:I:388:MET:HG3	2.04	0.58
3:F:58:PRO:CD	3:F:87:ILE:HA	2.33	0.58
3:I:224:ARG:HH21	3:I:224:ARG:CG	2.14	0.58
1:H:140:THR:HG21	1:H:145:ALA:HA	1.86	0.57
2:L:39:ASN:ND2	2:L:51:ARG:HG3	2.19	0.57
2:B:31:TYR:C	2:B:32:ILE:HD12	2.29	0.57
3:F:224:ARG:HH11	3:F:224:ARG:CG	2.15	0.57
3:I:320:LEU:HD23	3:I:321:ARG:N	2.19	0.57
3:I:62:GLY:O	3:I:63:LYS:HB3	2.04	0.57
3:F:43:LEU:N	3:F:292:LEU:HD23	2.20	0.57
1:H:73:ASP:HB3	1:H:76:THR:HG22	1.87	0.57
3:I:282:GLN:CA	3:I:282:GLN:HE21	2.17	0.57
2:B:24:ARG:HB3	2:B:75:ASP:OD2	2.04	0.57
2:L:8:PRO:C	2:L:107:THR:CG2	2.67	0.57
3:F:55:ARG:CG	3:F:55:ARG:HH21	2.18	0.57
1:H:135:PRO:CD	1:H:147:LEU:HB3	2.34	0.56
3:F:48:ASN:CB	3:F:51:LEU:CD1	2.80	0.56
3:F:286:GLY:O	3:F:288:ILE:HG21	2.05	0.56
2:L:95:GLN:HG2	2:L:96:GLY:H	1.68	0.56
2:L:98:HIS:CD2	2:L:99:TRP:N	2.73	0.56
1:H:124:SER:CB	1:H:155:PHE:CE1	2.67	0.56
1:A:140:THR:HG21	1:A:146:ALA:H	1.69	0.56
2:B:41:PHE:CZ	2:B:94:MET:SD	2.98	0.56
2:B:44:ARG:CZ	2:B:86:GLU:CG	2.84	0.56
3:F:50:LYS:HG3	3:F:273:PRO:O	2.06	0.56
1:H:126:LYS:CE	1:H:184:LEU:HD22	2.27	0.56
3:I:296:ASN:HD22	3:I:296:ASN:H	1.54	0.56
2:L:38:LEU:N	2:L:94:MET:CG	2.68	0.56
3:I:295:GLN:HB3	3:I:306:PRO:HG2	1.88	0.56
1:H:32:TYR:N	1:H:32:TYR:CD1	2.73	0.56
3:I:51:LEU:HD12	3:I:282:GLN:OE1	2.06	0.56
2:L:23:CYS:O	2:L:75:ASP:HA	2.06	0.56
1:H:126:LYS:HD3	1:H:155:PHE:CB	2.32	0.56
2:L:31:TYR:C	2:L:32:ILE:HD12	2.31	0.56
1:H:27:TYR:HA	1:H:28:SER:OG	2.06	0.56
1:H:106:GLY:HA2	2:L:35:ASN:ND2	2.19	0.56
3:I:220:ARG:HB3	3:I:221:PRO:HD2	1.86	0.55
1:A:63:LYS:HB2	1:A:64:LEU:HD22	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:32:ILE:HD12	2:B:32:ILE:N	2.21	0.55
3:F:58:PRO:HD3	3:F:86:TYR:O	2.05	0.55
2:B:57:PHE:HB2	3:F:289:ASN:HD22	1.53	0.55
2:L:45:PRO:CG	2:L:170:GLU:HG3	2.37	0.55
3:I:496:LYS:CB	3:I:497:LEU:CA	2.84	0.55
2:L:38:LEU:CB	2:L:94:MET:CG	2.82	0.55
1:A:107:PHE:HE2	2:B:99:TRP:HZ3	0.96	0.55
3:F:287:ALA:CA	3:F:289:ASN:CB	2.85	0.55
1:A:138:LYS:NZ	1:A:138:LYS:HB3	2.22	0.54
3:F:59:LEU:HD23	3:F:81:THR:HG23	1.88	0.54
1:H:59:ASN:ND2	2:L:99:TRP:NE1	2.56	0.54
1:A:157:GLU:HG2	1:A:158:PRO:CA	2.36	0.54
2:L:32:ILE:CG2	2:L:33:ASP:H	2.10	0.54
2:B:45:PRO:HG2	2:B:170:GLU:HG3	1.89	0.54
1:H:64:LEU:O	1:H:68:VAL:HG22	2.08	0.54
1:H:128:PRO:HB3	1:H:154:TYR:HB3	1.89	0.54
2:L:38:LEU:N	2:L:94:MET:HG2	2.22	0.54
3:F:338:PHE:CD2	3:F:338:PHE:C	2.85	0.54
1:A:107:PHE:HE2	2:B:99:TRP:CH2	1.65	0.54
2:L:96:GLY:O	2:L:98:HIS:N	2.41	0.54
2:B:95:GLN:HB3	2:B:101:TYR:CB	2.37	0.53
3:F:338:PHE:CG	3:F:339:ILE:N	2.75	0.53
2:L:188:LYS:O	2:L:192:GLU:HG2	2.08	0.53
2:L:203:HIS:HB3	2:L:206:LEU:HD13	1.89	0.53
2:B:39:ASN:CB	2:B:53:ILE:O	2.56	0.53
2:B:86:GLU:HA	2:B:173:SER:HA	1.90	0.53
3:F:280:THR:CA	3:F:288:ILE:HD11	2.38	0.53
1:A:138:LYS:HB3	1:A:138:LYS:HZ2	1.73	0.53
2:B:44:ARG:CZ	2:B:86:GLU:OE2	2.54	0.53
2:L:99:TRP:N	2:L:100:PRO:HD2	2.24	0.53
3:F:52:CYS:SG	3:F:277:CYS:N	2.81	0.53
2:L:6:GLN:NE2	2:L:105:GLN:N	2.57	0.53
1:A:128:PRO:HB3	1:A:154:TYR:HB3	1.91	0.53
2:B:35:ASN:C	2:B:35:ASN:HD22	2.17	0.53
3:F:184:HIS:HB3	3:F:220:ARG:HH12	1.73	0.52
1:H:126:LYS:CG	1:H:155:PHE:N	2.72	0.52
3:I:496:LYS:HZ2	3:I:497:LEU:HD13	1.74	0.52
3:I:323:ILE:O	3:I:342:GLY:N	2.41	0.52
2:L:38:LEU:HB2	2:L:94:MET:HE2	1.91	0.52
2:B:6:GLN:NE2	2:B:106:GLY:HA2	2.21	0.52
3:F:48:ASN:OD1	3:F:51:LEU:CD2	2.55	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:90:THR:O	3:I:92:SER:OG	2.19	0.52
2:L:32:ILE:N	2:L:32:ILE:CD1	2.73	0.52
2:B:32:ILE:N	2:B:32:ILE:CD1	2.73	0.52
1:H:157:GLU:CB	1:H:158:PRO:HA	2.39	0.52
3:I:47:HIS:CG	3:I:48:ASN:N	2.72	0.52
3:F:287:ALA:N	3:F:288:ILE:HG22	2.24	0.52
3:I:282:GLN:HA	3:I:282:GLN:HE21	1.75	0.52
1:H:140:THR:CG2	1:H:146:ALA:H	2.19	0.52
3:I:496:LYS:NZ	3:I:497:LEU:CD1	2.73	0.52
3:F:287:ALA:N	3:F:288:ILE:CG2	2.73	0.52
2:L:111:ILE:HG13	2:L:171:GLN:CD	2.35	0.52
2:L:99:TRP:O	2:L:101:TYR:N	2.43	0.51
1:H:126:LYS:HG3	1:H:155:PHE:N	2.25	0.51
1:H:157:GLU:OE2	1:H:177:ALA:HB3	2.10	0.51
1:A:23:LYS:HE3	1:A:25:SER:HB3	1.91	0.51
3:I:50:LYS:O	3:I:51:LEU:C	2.52	0.51
2:B:147:ARG:HB2	2:B:178:TYR:CE2	2.46	0.51
1:H:50:TRP:CZ2	1:H:52:SER:HB2	2.46	0.51
1:H:43:GLN:HG3	1:H:44:GLY:N	2.24	0.51
2:B:101:TYR:O	2:B:102:THR:OG1	2.28	0.51
3:I:282:GLN:NE2	3:I:282:GLN:CA	2.73	0.51
1:A:102:HIS:O	1:A:103:ILE:HG22	2.11	0.51
3:F:287:ALA:N	3:F:288:ILE:CB	2.73	0.51
3:I:206:SER:OG	3:I:241:ASP:OD2	2.28	0.51
3:I:496:LYS:CB	3:I:497:LEU:HA	2.41	0.51
1:H:100:ARG:HD2	1:H:102:HIS:O	2.12	0.50
3:I:224:ARG:NH2	3:I:224:ARG:CG	2.73	0.50
3:F:384:VAL:O	3:F:388:MET:HG2	2.12	0.50
2:L:38:LEU:CA	2:L:94:MET:CG	2.89	0.50
1:A:64:LEU:O	1:A:68:VAL:HG22	2.12	0.50
1:H:126:LYS:CB	1:H:126:LYS:NZ	2.73	0.50
2:L:95:GLN:HB2	2:L:101:TYR:HD2	1.74	0.50
1:A:107:PHE:HE1	2:B:31:TYR:HE1	1.59	0.50
1:H:135:PRO:CG	1:H:147:LEU:CB	2.85	0.50
1:A:138:LYS:NZ	1:A:138:LYS:CB	2.73	0.50
3:F:333:GLY:O	3:F:337:GLY:HA3	2.11	0.50
2:L:95:GLN:NE2	2:L:97:THR:O	2.44	0.50
1:H:126:LYS:CB	1:H:126:LYS:HZ3	2.24	0.50
3:I:40:VAL:HG12	2:L:33:ASP:HA	1.92	0.50
2:B:8:PRO:O	2:B:107:THR:CG2	2.56	0.50
2:B:188:LYS:O	2:B:192:GLU:HG2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:75:ASP:C	2:B:76:PHE:CD1	2.89	0.49
2:B:96:GLY:O	2:B:98:HIS:N	2.40	0.49
3:F:287:ALA:CA	3:F:288:ILE:CG2	2.86	0.49
3:I:47:HIS:ND1	3:I:48:ASN:N	2.60	0.49
1:A:64:LEU:N	1:A:64:LEU:CD2	2.73	0.49
1:A:157:GLU:CD	1:A:177:ALA:HB3	2.36	0.49
2:B:99:TRP:HZ3	2:B:101:TYR:HH	1.54	0.49
1:H:33:GLY:CA	1:H:100:ARG:O	2.60	0.49
1:H:33:GLY:HA3	1:H:100:ARG:O	2.12	0.49
2:B:33:ASP:OD2	3:F:39:SER:O	2.29	0.49
2:B:75:ASP:C	2:B:76:PHE:HD1	2.20	0.49
1:H:140:THR:HG21	1:H:145:ALA:C	2.38	0.49
3:I:47:HIS:ND1	3:I:49:GLY:N	2.58	0.49
2:B:30:VAL:O	2:B:31:TYR:CB	2.60	0.49
3:I:296:ASN:ND2	3:I:296:ASN:N	2.59	0.49
3:I:471:HIS:HB3	3:I:493:GLU:OE1	2.13	0.49
2:B:91:TYR:O	2:B:106:GLY:HA3	2.13	0.48
2:B:95:GLN:HB3	2:B:101:TYR:HB3	1.94	0.48
3:F:220:ARG:HG3	3:F:229:ARG:HG3	1.94	0.48
2:B:118:PRO:HD2	2:B:206:LEU:CD1	2.41	0.48
3:F:55:ARG:CG	3:F:55(A):GLY:N	2.73	0.48
3:F:58:PRO:CG	3:F:87:ILE:HG22	2.41	0.48
3:F:334:ALA:HA	3:F:338:PHE:CE1	2.48	0.48
2:B:23:CYS:O	2:B:76:PHE:N	2.46	0.48
2:L:86:GLU:HA	2:L:173:SER:HB2	1.95	0.48
2:B:111:ILE:C	2:B:171:GLN:HE22	2.22	0.48
3:F:287:ALA:HA	3:F:289:ASN:HB2	1.95	0.48
1:H:36:TRP:CD1	1:H:70:MET:HE2	2.42	0.48
3:I:496:LYS:HZ3	3:I:497:LEU:CD1	2.25	0.48
1:A:91:THR:HG23	1:A:119:THR:HA	1.95	0.48
3:F:51:LEU:HG	3:F:53:LYS:HE3	1.95	0.48
3:F:55:ARG:CG	3:F:55:ARG:NH2	2.76	0.48
1:H:27:TYR:HB3	1:H:29:PHE:HB2	1.95	0.48
3:I:297:ILE:HG22	3:I:298:HIS:CD2	2.48	0.48
2:L:76:PHE:CD2	2:L:77:THR:HB	2.49	0.48
1:H:140:THR:HG22	2:L:121:PHE:CD2	2.49	0.48
3:I:492:SER:HB2	3:I:494:GLU:HG2	1.95	0.48
3:F:287:ALA:HA	3:F:289:ASN:CB	2.44	0.47
3:F:288:ILE:N	3:F:289:ASN:C	2.72	0.47
1:H:133:LEU:O	1:H:135:PRO:HD3	2.13	0.47
2:L:118:PRO:HD2	2:L:206:LEU:CD1	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:SER:O	1:A:141:SER:CA	2.59	0.47
2:B:41:PHE:CE2	2:B:94:MET:SD	3.07	0.47
2:L:38:LEU:HB3	2:L:94:MET:HE2	1.95	0.47
2:L:95:GLN:CB	2:L:101:TYR:CD2	2.95	0.47
1:H:135:PRO:HD2	1:H:147:LEU:HB3	1.97	0.47
3:I:44:GLU:HB2	3:I:292:LEU:HD23	1.97	0.47
3:I:462:ILE:HD11	3:I:468:GLU:HB2	1.96	0.47
2:L:6:GLN:HE21	2:L:105:GLN:H	1.62	0.47
2:L:38:LEU:N	2:L:94:MET:HG3	2.29	0.47
1:A:107:PHE:CE1	2:B:31:TYR:HE1	2.32	0.47
1:H:135:PRO:CD	1:H:147:LEU:CB	2.92	0.47
1:H:136:SER:O	1:H:139:SER:CA	2.62	0.47
3:I:63:LYS:O	3:I:63:LYS:HG3	2.14	0.47
3:I:79:LEU:HD23	3:I:117:ARG:HD3	1.97	0.47
3:I:179:LEU:O	3:I:254:PRO:HB3	2.14	0.47
3:I:281:CYS:O	3:I:302:ILE:O	2.33	0.47
2:L:99:TRP:HA	2:L:99:TRP:HE3	1.76	0.47
1:H:140:THR:HG22	2:L:121:PHE:HE2	1.78	0.47
3:I:47:HIS:CE1	3:I:286:GLY:HA3	2.50	0.47
2:L:30:VAL:HG23	2:L:31:TYR:H	1.80	0.47
1:A:31:SER:HB2	1:A:32:TYR:CD2	2.49	0.47
2:B:97:THR:OG1	2:B:98:HIS:N	2.48	0.46
3:F:79:LEU:HG	3:F:80:SER:O	2.16	0.46
1:A:161:VAL:HA	1:A:206:ASN:O	2.16	0.46
3:F:260:MET:HE2	3:F:262:ARG:HG3	1.98	0.46
3:F:287:ALA:CA	3:F:288:ILE:CB	2.93	0.46
1:A:143:GLY:O	1:A:144:THR:OG1	2.29	0.46
2:B:35:ASN:C	2:B:35:ASN:ND2	2.73	0.46
2:B:96:GLY:O	2:B:97:THR:OG1	2.25	0.46
1:H:161:VAL:HA	1:H:206:ASN:O	2.14	0.46
3:I:282:GLN:NE2	3:I:287:ALA:HB2	2.31	0.46
2:L:30:VAL:HG23	2:L:31:TYR:N	2.31	0.46
2:B:57:PHE:CB	3:F:289:ASN:CG	2.83	0.46
1:H:59:ASN:ND2	2:L:99:TRP:HE1	2.14	0.46
2:L:38:LEU:CA	2:L:94:MET:HG2	2.45	0.46
2:B:98:HIS:HB3	2:B:100:PRO:HG2	1.97	0.46
3:I:195:TYR:O	3:I:197:ASN:N	2.48	0.46
3:I:79:LEU:CD2	3:I:117:ARG:HD3	2.46	0.46
2:B:44:ARG:NH1	2:B:86:GLU:OE2	2.49	0.45
1:H:100:ARG:HD3	1:H:107:PHE:HD1	1.81	0.45
3:I:496:LYS:CB	3:I:497:LEU:CB	2.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:99:TRP:O	2:L:100:PRO:C	2.59	0.45
1:H:64:LEU:N	1:H:64:LEU:CD2	2.73	0.45
1:H:126:LYS:HG2	1:H:155:PHE:N	2.31	0.45
2:L:6:GLN:HE21	2:L:105:GLN:N	2.13	0.45
2:L:95:GLN:HG3	2:L:101:TYR:CA	2.42	0.45
2:B:57:PHE:CD1	3:F:289:ASN:ND2	2.84	0.45
2:B:99:TRP:N	2:B:100:PRO:HD2	2.32	0.45
3:F:286:GLY:C	3:F:288:ILE:CB	2.90	0.45
1:H:135:PRO:CG	1:H:147:LEU:CA	2.95	0.45
2:B:40:TRP:HE3	2:B:78:LEU:CD2	2.29	0.45
2:B:98:HIS:O	2:B:101:TYR:HE1	1.99	0.45
3:F:58:PRO:HB3	3:F:84:TRP:HB2	1.99	0.45
2:B:98:HIS:O	2:B:101:TYR:HD1	1.97	0.45
1:H:140:THR:HG21	1:H:145:ALA:CA	2.47	0.45
2:B:40:TRP:CZ3	2:B:78:LEU:HB3	2.51	0.45
3:F:286:GLY:C	3:F:288:ILE:HB	2.42	0.45
3:F:455:LEU:O	3:F:455:LEU:HD12	2.17	0.45
2:L:111:ILE:C	2:L:171:GLN:HE22	2.24	0.45
3:F:55:ARG:O	3:F:56:VAL:HG23	2.16	0.45
2:L:45:PRO:HG2	2:L:170:GLU:HG3	1.97	0.45
1:A:140:THR:HB	1:A:145:ALA:HA	1.98	0.44
1:A:202:THR:HG23	1:A:219:LYS:HE2	1.99	0.44
2:B:38:LEU:HD13	2:B:38:LEU:HA	1.78	0.44
3:I:47:HIS:O	3:I:48:ASN:HB2	2.17	0.44
2:B:73:GLY:O	2:B:75:ASP:N	2.43	0.44
1:H:2:VAL:HG12	1:H:26:GLY:O	2.17	0.44
1:A:29:PHE:CZ	1:A:30:SER:O	2.70	0.44
2:B:85:ALA:HB1	2:B:172:ASP:O	2.18	0.44
2:B:96:GLY:C	2:B:97:THR:HG23	2.41	0.44
3:F:284:PRO:HA	3:F:286:GLY:H	1.82	0.44
3:I:50:LYS:CB	3:I:275:HIS:CD2	3.01	0.44
2:L:6:GLN:NE2	2:L:105:GLN:C	2.75	0.44
1:H:100:ARG:HD3	1:H:107:PHE:CE1	2.52	0.44
2:L:77:THR:OG1	2:L:78:LEU:N	2.51	0.44
1:A:29:PHE:CE2	1:A:30:SER:O	2.70	0.44
3:F:48:ASN:HA	3:F:51:LEU:HD11	2.00	0.43
3:I:268:ILE:HD11	3:I:302:ILE:CD1	2.46	0.43
3:I:496:LYS:NZ	3:I:497:LEU:HD13	2.32	0.43
1:A:57:ASN:HB2	3:F:367:LEU:HD22	2.00	0.43
1:A:157:GLU:OE1	1:A:177:ALA:CB	2.64	0.43
3:F:281:CYS:SG	3:F:282:GLN:N	2.91	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:32:ILE:CG2	2:L:33:ASP:N	2.69	0.43
1:A:26:GLY:HA2	1:A:27:TYR:HA	1.58	0.43
3:F:58:PRO:HG3	3:F:87:ILE:CG2	2.46	0.43
3:I:305:CYS:HB3	3:I:306:PRO:HD2	2.01	0.43
1:A:140:THR:HG21	1:A:146:ALA:N	2.34	0.43
1:H:32:TYR:N	1:H:32:TYR:HD1	2.15	0.43
3:I:47:HIS:HE1	3:I:49:GLY:O	2.01	0.43
3:I:455:LEU:HD12	3:I:455:LEU:O	2.19	0.43
2:L:38:LEU:O	2:L:54:TYR:HA	2.19	0.43
1:A:102:HIS:HB3	1:A:105:THR:HG22	2.01	0.43
3:F:54:LEU:HD12	3:F:54:LEU:HA	1.86	0.43
3:F:55(A):GLY:N	3:F:56:VAL:HG23	2.33	0.43
1:H:137:SER:HB3	1:H:140:THR:H	1.84	0.43
3:F:50:LYS:HE3	3:F:273:PRO:HB2	2.01	0.43
2:L:74:THR:O	2:L:75:ASP:CG	2.62	0.42
1:H:202:THR:HG23	1:H:219:LYS:HE2	2.00	0.42
3:I:288:ILE:HD12	3:I:290:THR:H	1.84	0.42
1:A:63:LYS:C	1:A:64:LEU:HD22	2.44	0.42
3:I:63:LYS:O	3:I:64:CYS:SG	2.77	0.42
2:L:40:TRP:CG	2:L:78:LEU:HD22	2.55	0.42
3:F:334:ALA:HB3	3:F:441:ASP:OD1	2.19	0.42
1:H:2:VAL:HA	1:H:26:GLY:HA3	2.02	0.42
1:H:126:LYS:HG2	1:H:155:PHE:HB3	1.98	0.42
2:L:96:GLY:C	2:L:97:THR:HG1	2.27	0.42
1:H:36:TRP:NE1	1:H:70:MET:HE2	2.35	0.42
1:H:63:LYS:HB2	1:H:64:LEU:CD2	2.47	0.42
1:H:143:GLY:O	1:H:144:THR:OG1	2.31	0.42
2:B:39:ASN:HB2	2:B:53:ILE:O	2.19	0.42
2:B:40:TRP:HE3	2:B:78:LEU:HD23	1.84	0.42
1:A:2:VAL:O	1:A:3:GLN:HG3	2.20	0.42
1:H:126:LYS:HG3	1:H:154:TYR:C	2.45	0.42
1:H:145:ALA:O	1:H:192:THR:HA	2.19	0.42
3:I:79:LEU:HA	3:I:80:SER:C	2.44	0.42
2:B:86:GLU:N	2:B:173:SER:O	2.53	0.42
3:I:320:LEU:HD23	3:I:321:ARG:H	1.84	0.42
3:F:59:LEU:HD21	3:F:81:THR:H	1.82	0.41
1:H:158:PRO:HD2	1:H:158:PRO:O	2.20	0.41
3:I:290:THR:OG1	3:I:306:PRO:HD3	2.20	0.41
3:I:26:VAL:HG21	3:I:317:ALA:HB2	2.02	0.41
3:I:41:ASN:OD1	3:I:43:LEU:O	2.38	0.41
3:I:60:HIS:HA	3:I:88:VAL:HG23	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:260:MET:HE2	3:I:262:ARG:HB3	2.03	0.41
2:B:99:TRP:N	2:B:100:PRO:CD	2.84	0.41
3:F:41:ASN:HB3	3:F:292:LEU:HD21	2.03	0.41
3:F:280:THR:OG1	3:F:288:ILE:HD13	2.04	0.41
1:A:145:ALA:O	1:A:192:THR:HA	2.21	0.41
1:H:33:GLY:O	1:H:99:ASP:N	2.46	0.41
3:I:60:HIS:HA	3:I:88:VAL:CG2	2.51	0.41
2:B:29:LEU:HD11	2:B:37:TYR:CD2	2.55	0.41
3:F:223:VAL:HB	3:F:229:ARG:NH1	2.34	0.41
3:F:79:LEU:HA	3:F:80:SER:C	2.46	0.41
1:H:135:PRO:HD2	1:H:147:LEU:CB	2.51	0.41
3:I:496:LYS:CG	3:I:497:LEU:CA	2.86	0.41
2:B:145:TYR:CG	2:B:146:PRO:HA	2.56	0.41
3:I:90:THR:HG22	3:I:91:SER:N	2.33	0.41
2:B:39:ASN:CB	2:B:54:TYR:HA	2.28	0.41
1:H:91:THR:CG2	1:H:119:THR:HA	2.50	0.41
1:H:138:LYS:HD2	1:H:138:LYS:O	2.21	0.41
3:I:99:PRO:HB2	3:I:229:ARG:HD3	2.03	0.41
3:I:496:LYS:HB3	3:I:497:LEU:CA	2.46	0.41
2:L:38:LEU:CA	2:L:94:MET:HG3	2.50	0.41
1:A:99:ASP:OD1	1:A:100:ARG:N	2.48	0.40
3:F:107:GLU:O	3:F:111:GLN:HG2	2.21	0.40
2:L:6:GLN:HE22	2:L:106:GLY:CA	2.25	0.40
2:L:21:ILE:O	2:L:77:THR:HA	2.21	0.40
1:H:63:LYS:C	1:H:64:LEU:HD13	2.46	0.40
3:I:70:ILE:O	3:I:150:ASN:ND2	2.53	0.40
3:F:50:LYS:HD2	3:F:273:PRO:HG2	2.04	0.40
3:I:47:HIS:CE1	3:I:49:GLY:O	2.75	0.40
3:F:30:LEU:HD21	3:I:434:GLU:HG3	2.02	0.40
3:F:299:PRO:HB2	3:F:300:ILE:HD12	2.03	0.40
1:H:157:GLU:OE2	1:H:177:ALA:CB	2.69	0.40
3:I:186:SER:HB3	3:I:227:GLU:HB3	2.04	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	217/227 (96%)	196 (90%)	18 (8%)	3 (1%)	9	30
1	H	217/227 (96%)	193 (89%)	21 (10%)	3 (1%)	9	30
2	B	209/219 (95%)	187 (90%)	19 (9%)	3 (1%)	9	30
2	L	213/219 (97%)	186 (87%)	20 (9%)	7 (3%)	3	17
3	F	461/505 (91%)	422 (92%)	33 (7%)	6 (1%)	9	32
3	I	460/505 (91%)	426 (93%)	32 (7%)	2 (0%)	30	58
All	All	1777/1902 (93%)	1610 (91%)	143 (8%)	24 (1%)	9	30

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	THR
2	B	82	THR
3	F	60	HIS
3	F	288	ILE
2	L	34	GLY
2	L	39	ASN
2	L	82	THR
3	F	50	LYS
3	F	81	THR
2	L	32	ILE
2	B	31	TYR
2	B	102	THR
1	H	101	PRO
3	I	81	THR
2	L	97	THR
1	A	222	PRO
3	F	58	PRO
1	H	135	PRO
1	H	222	PRO
3	I	48	ASN
2	L	100	PRO
2	L	102	THR
1	A	103	ILE
3	F	284	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	186/192 (97%)	180 (97%)	6 (3%)	34	59
1	H	187/192 (97%)	181 (97%)	6 (3%)	34	59
2	B	188/194 (97%)	180 (96%)	8 (4%)	26	51
2	L	192/194 (99%)	186 (97%)	6 (3%)	35	59
3	F	407/440 (92%)	394 (97%)	13 (3%)	34	59
3	I	408/440 (93%)	397 (97%)	11 (3%)	39	61
All	All	1568/1652 (95%)	1518 (97%)	50 (3%)	34	59

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	SER
1	A	31	SER
1	A	65	GLN
1	A	137	SER
1	A	138	LYS
1	A	139	SER
2	B	22	SER
2	B	31	TYR
2	B	32	ILE
2	B	33	ASP
2	B	35	ASN
2	B	38	LEU
2	B	107	THR
2	B	111	ILE
3	F	46	LYS
3	F	47	HIS
3	F	55	ARG
3	F	58	PRO
3	F	59	LEU
3	F	79	LEU
3	F	224	ARG
3	F	276	ASP

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Mol	Chain	Res	Type
3	F	288	ILE
3	F	289	ASN
3	F	292	LEU
3	F	348	ASP
3	F	488	TYR
1	H	32	TYR
1	H	64	LEU
1	H	125	THR
1	H	126	LYS
1	H	138	LYS
1	H	140	THR
3	I	46	LYS
3	I	51	LEU
3	I	52	CYS
3	I	63	LYS
3	I	208	ARG
3	I	224	ARG
3	I	282	GLN
3	I	288	ILE
3	I	296	ASN
3	I	455	LEU
3	I	496	LYS
2	L	32	ILE
2	L	94	MET
2	L	99	TRP
2	L	105	GLN
2	L	107	THR
2	L	111	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
2	B	27	GLN
2	B	35	ASN
2	B	95	GLN
2	B	157	ASN
2	B	194	HIS
3	F	133	ASN
3	F	192	GLN
3	F	275	HIS
3	F	289	ASN

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Mol	Chain	Res	Type
3	F	382	ASN
3	F	424	ASN
3	F	464	ASN
3	F	483	ASN
1	H	59	ASN
3	I	41	ASN
3	I	94	ASN
3	I	192	GLN
3	I	275	HIS
3	I	296	ASN
3	I	298	HIS
3	I	424	ASN
3	I	483	ASN
2	L	6	GLN
2	L	35	ASN
2	L	39	ASN
2	L	98	HIS
2	L	157	ASN
2	L	194	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

3 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
4	NAG	C	1	4,3	14,14,15	0.91	1 (7%)	17,19,21	1.19	1 (5%)
4	NAG	C	2	4	14,14,15	0.26	0	17,19,21	0.72	0
4	BMA	C	3	4	11,11,12	0.67	0	15,15,17	0.92	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	C	1	4,3	-	1/6/23/26	0/1/1/1
4	NAG	C	2	4	-	2/6/23/26	0/1/1/1
4	BMA	C	3	4	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1	NAG	O5-C1	3.11	1.48	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1	NAG	C1-O5-C5	4.69	118.47	112.19
4	C	3	BMA	C1-O5-C5	2.19	115.12	112.19

There are no chirality outliers.

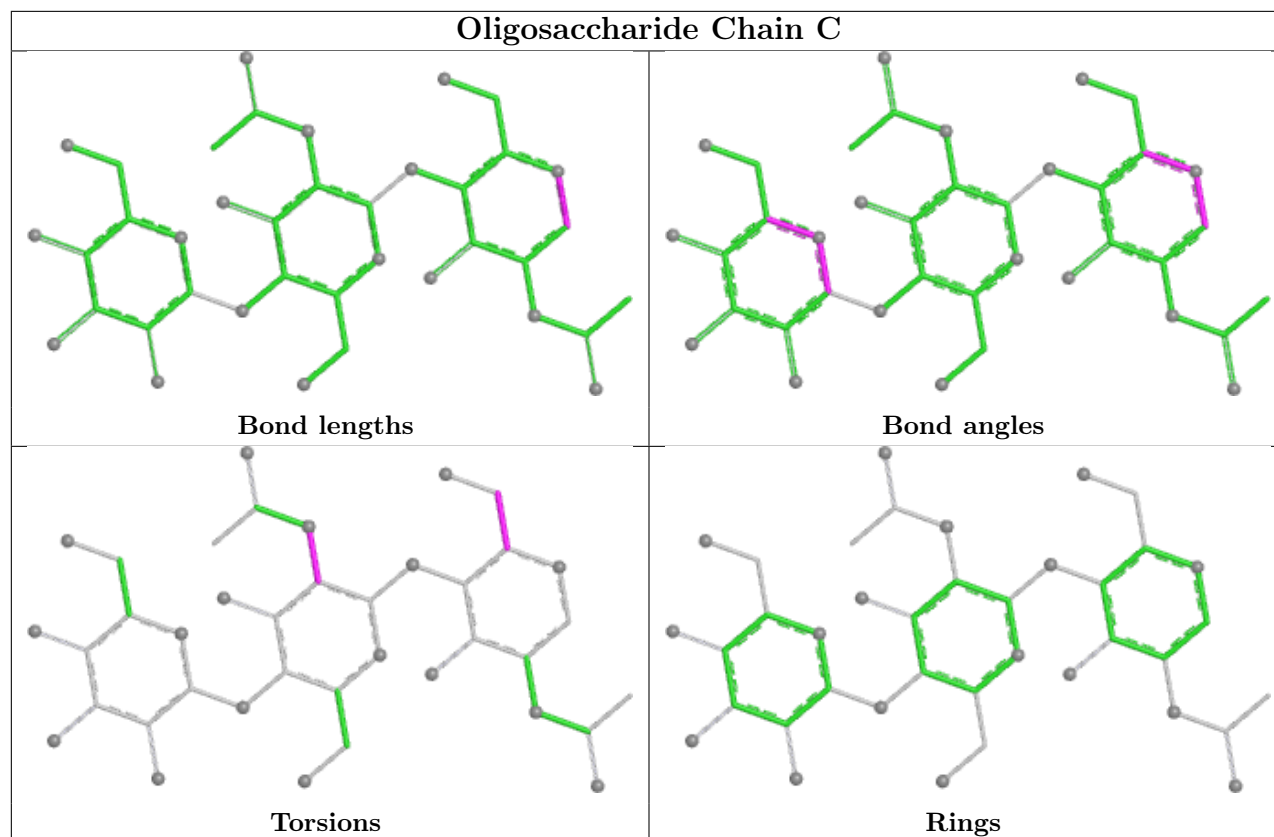
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	1	NAG	O5-C5-C6-O6
4	C	2	NAG	C3-C2-N2-C7
4	C	2	NAG	C1-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	A	221/227 (97%)	0.39	17 (7%)	19	16	110, 280, 411, 476	0
1	H	221/227 (97%)	0.33	13 (5%)	28	21	124, 248, 353, 409	3 (1%)
2	B	213/219 (97%)	0.38	15 (7%)	22	18	164, 286, 364, 434	2 (0%)
2	L	217/219 (99%)	0.42	11 (5%)	33	24	143, 256, 349, 411	3 (1%)
3	F	467/505 (92%)	0.34	26 (5%)	30	22	54, 155, 288, 362	9 (1%)
3	I	466/505 (92%)	0.17	22 (4%)	36	27	44, 122, 236, 322	9 (1%)
All	All	1805/1902 (94%)	0.31	104 (5%)	29	21	44, 208, 351, 476	26 (1%)

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	55(A)	GLY	6.3
3	F	291	SER	6.0
2	B	38	LEU	5.5
3	F	335	ILE	4.8
3	F	334	ALA	4.6
2	B	99	TRP	4.3
3	I	49	GLY	4.2
2	L	201	VAL	4.1
2	L	37	TYR	4.1
1	A	59	ASN	4.0
1	A	104	LEU	3.9
1	A	148	GLY	3.7
2	L	32	ILE	3.7
3	F	276	ASP	3.7
3	I	350	TRP	3.7
3	I	352	GLY	3.5
1	H	104	LEU	3.5
3	I	245	PHE	3.5
1	A	128	PRO	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	117	LEU	3.4
3	F	48	ASN	3.4
3	I	225	ASP	3.4
2	L	166	GLU	3.4
2	B	41	PHE	3.3
2	B	166	GLU	3.3
1	A	35	SER	3.2
1	A	110	ASP	3.2
3	I	47	HIS	3.2
2	B	37	TYR	3.2
2	L	102	THR	3.1
2	L	186	LEU	3.1
1	A	133	LEU	3.1
2	L	190	ASP	3.1
3	F	283	THR	3.0
3	F	371	GLN	3.0
2	B	199	CYS	3.0
3	I	340	GLU	3.0
3	I	271	ASP	2.9
3	F	225	ASP	2.9
2	B	35	ASN	2.8
3	F	127	TRP	2.8
1	A	103	ILE	2.8
2	B	36	THR	2.8
3	I	432	GLU	2.7
2	B	121	PHE	2.6
3	F	179	LEU	2.6
3	F	491	TYR	2.6
3	I	300	ILE	2.6
3	I	295	GLN	2.6
2	B	70	SER	2.6
2	B	30	VAL	2.6
1	A	34	ILE	2.5
3	F	316	LEU	2.5
3	I	345	GLY	2.5
3	I	106	GLU	2.5
2	L	65	ASP	2.4
3	I	339	ILE	2.4
3	F	53	LYS	2.4
1	A	105	THR	2.4
3	F	278	ASN	2.4
1	H	148	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	151	VAL	2.3
2	B	123	PHE	2.3
1	H	149	CYS	2.3
2	L	2	VAL	2.3
3	I	297	ILE	2.3
1	H	58	THR	2.3
3	F	51	LEU	2.3
1	H	11	VAL	2.3
3	I	40	VAL	2.3
3	F	178	VAL	2.2
3	F	289	ASN	2.2
1	H	157	GLU	2.2
2	L	38	LEU	2.2
1	A	94	PHE	2.2
1	A	52	SER	2.2
3	F	223	VAL	2.2
3	I	278	ASN	2.2
1	H	109	PHE	2.2
3	F	386	GLU	2.2
3	I	299	PRO	2.2
1	A	107	PHE	2.2
3	I	338	PHE	2.2
1	H	57	ASN	2.2
3	I	348	ASP	2.2
3	F	293	PRO	2.1
3	I	94	ASN	2.1
2	B	93	CYS	2.1
1	H	51	ILE	2.1
1	H	52	SER	2.1
1	H	59	ASN	2.1
2	B	138	VAL	2.1
3	F	347	VAL	2.1
1	A	48	MET	2.1
3	F	338	PHE	2.1
3	F	85	SER	2.1
3	F	346	MET	2.0
1	H	89	ASP	2.0
1	A	149	CYS	2.0
2	L	151	VAL	2.0
2	B	21	ILE	2.0
3	I	326	ILE	2.0
1	A	57	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
3	F	300	ILE	2.0

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

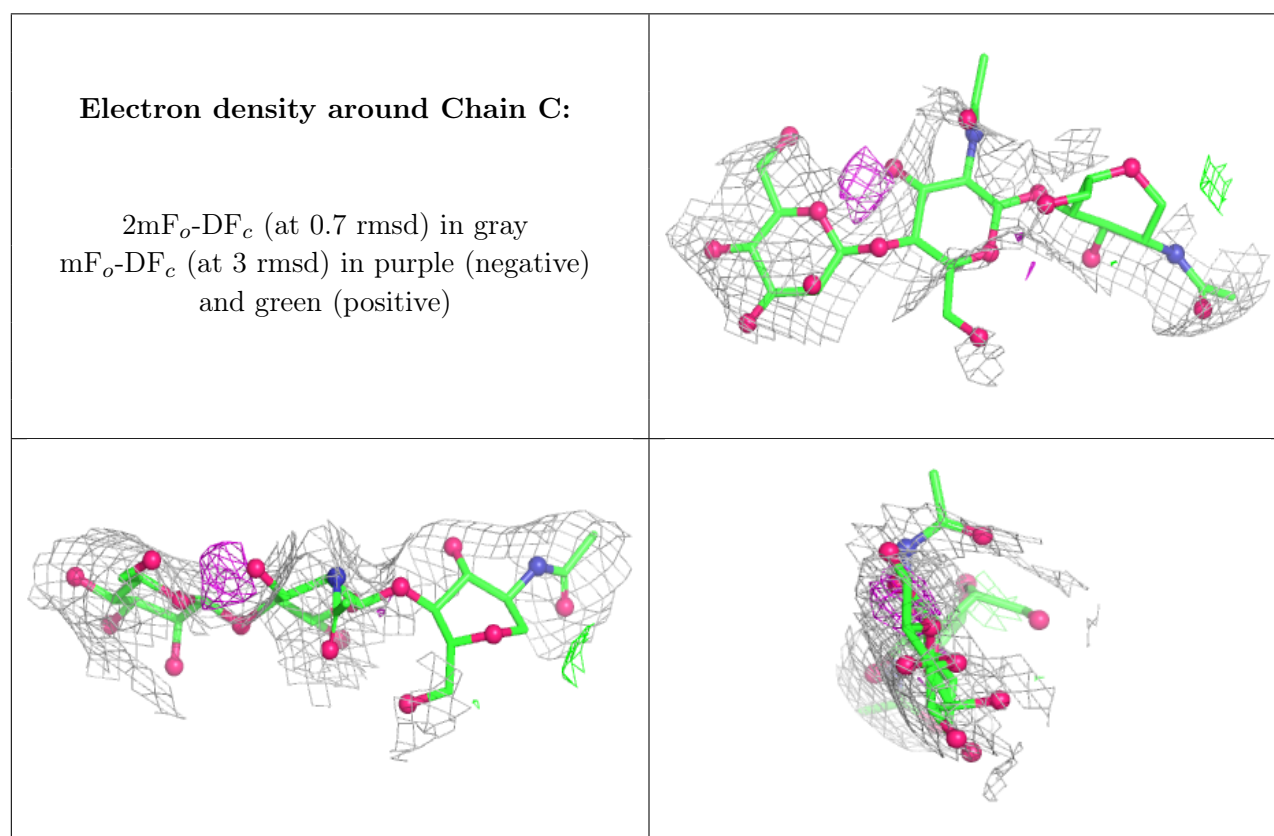
There are no non-standard protein/DNA/RNA residues in this entry.

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
4	NAG	C	1	14/15	-	-	212,224,243,250	0
4	NAG	C	2	14/15	-	-	250,259,274,277	0
4	BMA	C	3	11/12	-	-	247,256,270,274	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.