



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:11 AM UTC

PDB ID : 1HRN / pdb_00001hrn
Title : HIGH RESOLUTION CRYSTAL STRUCTURES OF RECOMBINANT HUMAN RENIN IN COMPLEX WITH POLYHYDROXYMONOAMIDE INHIBITORS
Authors : Tong, L.; Anderson, P.C.
Deposited on : 1995-03-31
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

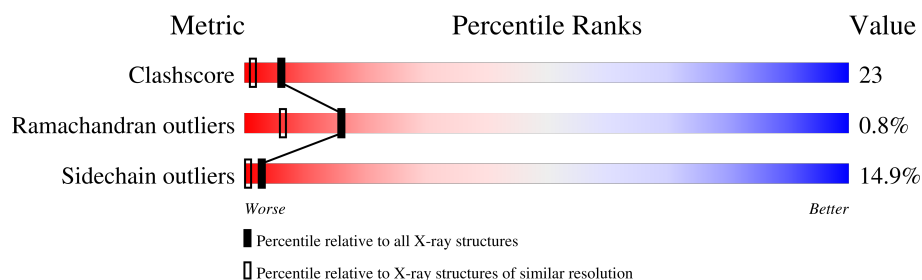
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	337	 65% 24% 9% ..
1	B	337	 52% 33% 10% ..

2 Entry composition [i](#)

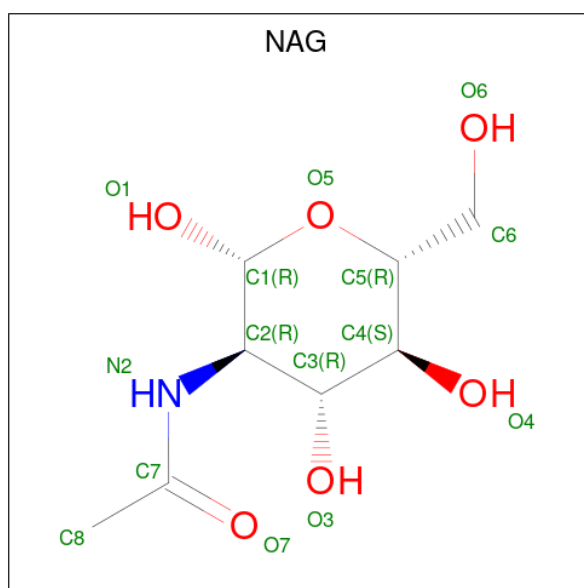
There are 4 unique types of molecules in this entry. The entry contains 5593 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 RENIN.

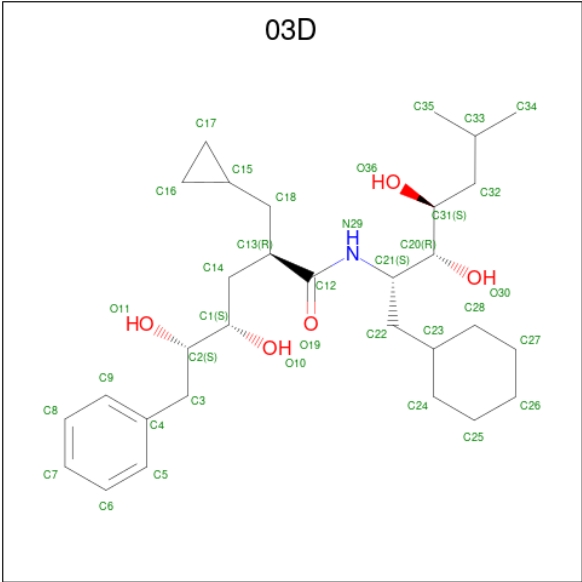
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	335	Total	C	N	O	S	0	0	1
			2576	1644	417	501	14			
1	B	334	Total	C	N	O	S	0	0	1
			2551	1630	409	498	14			

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



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

- Molecule 3 is (2R,4S,5S)-N-[(2S,3R,4S)-1-cyclohexyl-3,4-dihydroxy-6-methylheptan-2-yl]-2-(cyclopropylmethyl)-4,5-dihydroxy-6-phenylhexanamide (CCD ID: 03D) (formula: $C_{30}H_{49}NO_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			36	30	1	5		
3	B	1	Total	C	N	O	0	0
			36	30	1	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	206	Total	O	0	0
			206	206		
4	B	160	Total	O	0	0
			160	160		

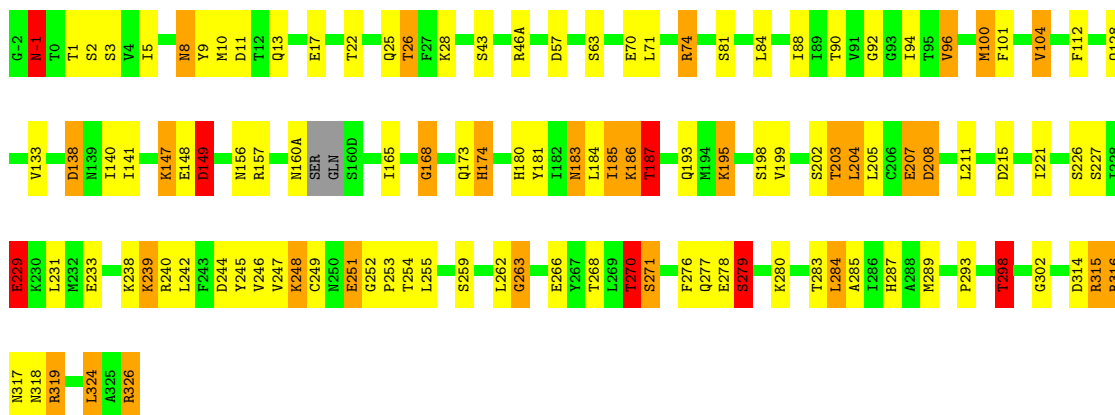
3 Residue-property plots

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

Note EDS was not executed.

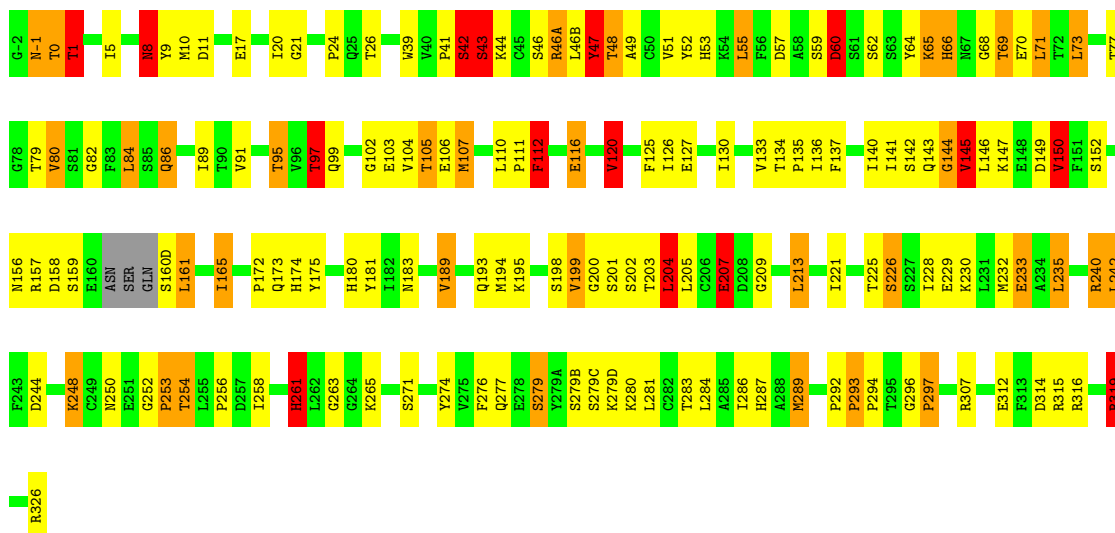
- Molecule 1: RENIN

Chain A: 



- Molecule 1: RENIN

Chain B: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	141.10Å 141.10Å 141.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.215 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5593	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, 03D

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.41	7/2635 (0.3%)	1.72	37/3572 (1.0%)
1	B	1.43	9/2610 (0.3%)	1.80	57/3542 (1.6%)
All	All	1.42	16/5245 (0.3%)	1.76	94/7114 (1.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	1	0

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	159	SER	C-N	-7.81	1.22	1.33
1	B	102	GLY	CA-C	-6.85	1.47	1.52
1	A	251	GLU	CD-OE1	6.83	1.38	1.25
1	A	74	ARG	CZ-NH2	6.72	1.42	1.33
1	B	289	MET	N-CA	-5.95	1.39	1.46
1	B	142	SER	CA-C	-5.74	1.44	1.52
1	A	233	GLU	CD-OE1	5.65	1.36	1.25
1	B	47	TYR	CA-C	-5.50	1.45	1.52
1	A	229	GLU	CD-OE1	5.41	1.35	1.25
1	B	17	GLU	CD-OE2	5.39	1.35	1.25
1	B	165	ILE	CA-CB	-5.30	1.48	1.54
1	A	185	ILE	N-CA	-5.27	1.40	1.46
1	B	207	GLU	CD-OE1	5.23	1.35	1.25
1	A	22	THR	C-O	-5.18	1.21	1.25
1	B	261	HIS	CG-CD2	5.08	1.41	1.35
1	A	13	GLN	C-O	-5.05	1.18	1.24

All (94) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	208	ASP	N-CA-CB	10.20	124.30	114.10
1	B	107	MET	O-C-N	9.76	127.12	121.27
1	A	263	GLY	CA-C-N	-9.68	102.44	121.41
1	A	263	GLY	C-N-CA	-9.68	102.44	121.41
1	B	292	PRO	N-CA-CB	9.54	108.18	103.22
1	A	293	PRO	N-CA-CB	9.19	108.34	103.19
1	A	180	HIS	CA-CB-CG	-9.02	104.78	113.80
1	A	183	ASN	CA-CB-CG	-8.56	104.04	112.60
1	A	293	PRO	N-CA-C	-8.50	102.31	110.47
1	A	270	THR	N-CA-CB	-8.27	97.64	110.46
1	B	107	MET	CA-C-N	8.07	128.54	119.83
1	B	107	MET	C-N-CA	8.07	128.54	119.83
1	A	26	THR	N-CA-CB	-7.75	97.67	110.23
1	B	174	HIS	CA-CB-CG	-7.56	106.24	113.80
1	B	319	ARG	NE-CZ-NH2	-7.41	112.53	119.20
1	B	144	GLY	N-CA-C	-7.39	98.76	111.47
1	B	319	ARG	NE-CZ-NH1	7.37	128.87	121.50
1	B	279(C)	SER	N-CA-C	-7.36	104.28	113.18
1	B	199	VAL	CB-CA-C	7.27	120.94	110.98
1	B	202	SER	N-CA-C	7.26	121.03	108.76
1	B	43	SER	N-CA-C	-7.13	102.33	111.02
1	A	293	PRO	CB-CA-C	-7.08	104.74	111.39
1	B	60	ASP	CA-C-N	-7.03	113.21	123.05
1	B	60	ASP	C-N-CA	-7.03	113.21	123.05
1	B	293	PRO	N-CA-C	-6.55	102.70	110.70
1	A	10	MET	CB-CA-C	-6.51	101.43	111.66
1	B	95	THR	OG1-CB-CG2	6.44	122.19	109.30
1	B	71	LEU	CB-CA-C	-6.39	96.40	109.38
1	A	185	ILE	CB-CA-C	-6.26	103.87	111.88
1	A	207	GLU	CA-C-N	-6.22	117.10	126.86
1	A	207	GLU	C-N-CA	-6.22	117.10	126.86
1	B	47	TYR	N-CA-CB	6.21	119.33	109.51
1	B	69	THR	CB-CA-C	6.21	117.44	109.80
1	B	149	ASP	N-CA-CB	-6.19	102.98	110.53
1	A	104	VAL	CB-CA-C	6.17	119.17	111.15
1	A	298	THR	CA-CB-OG1	6.12	118.78	109.60
1	A	203	THR	CA-C-N	-6.07	112.61	122.26
1	A	203	THR	C-N-CA	-6.07	112.61	122.26
1	A	70	GLU	CA-CB-CG	-6.04	102.02	114.10
1	B	95	THR	CA-CB-OG1	5.97	118.56	109.60
1	B	60	ASP	CA-CB-CG	-5.94	106.66	112.60
1	A	-1	ASN	CA-CB-CG	5.93	118.53	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	187	THR	CA-CB-OG1	-5.89	100.77	109.60
1	A	81	SER	CA-CB-OG	-5.86	99.38	111.10
1	A	149	ASP	CB-CA-C	-5.85	103.95	111.50
1	B	26	THR	CA-CB-OG1	-5.84	100.85	109.60
1	B	8	ASN	CA-CB-CG	5.82	118.42	112.60
1	A	101	PHE	CA-C-N	-5.81	116.73	122.36
1	A	101	PHE	C-N-CA	-5.81	116.73	122.36
1	B	279	SER	N-CA-CB	-5.79	102.41	111.56
1	A	112	PHE	N-CA-C	5.72	119.52	112.54
1	B	221	ILE	N-CA-C	-5.70	100.37	108.58
1	B	201	SER	CA-C-N	-5.69	113.97	122.74
1	B	201	SER	C-N-CA	-5.69	113.97	122.74
1	B	112	PHE	CA-CB-CG	-5.67	108.12	113.80
1	B	137	PHE	O-C-N	5.66	128.12	122.12
1	A	149	ASP	N-CA-CB	-5.64	101.79	110.25
1	A	203	THR	N-CA-CB	5.60	119.95	110.49
1	B	44	LYS	CA-C-N	-5.49	113.57	121.42
1	B	44	LYS	C-N-CA	-5.49	113.57	121.42
1	B	319	ARG	CD-NE-CZ	5.47	132.05	124.40
1	B	292	PRO	CA-C-N	-5.42	114.79	120.38
1	B	292	PRO	C-N-CA	-5.42	114.79	120.38
1	B	1	THR	N-CA-CB	-5.41	101.47	111.53
1	B	242	LEU	N-CA-CB	5.38	119.58	110.49
1	A	221	ILE	N-CA-C	-5.37	100.68	108.36
1	B	276	PHE	CA-CB-CG	-5.37	108.43	113.80
1	A	293	PRO	O-C-N	5.32	123.76	121.31
1	B	204	LEU	N-CA-CB	-5.32	102.33	110.26
1	A	185	ILE	N-CA-CB	-5.31	104.62	110.51
1	B	189	VAL	CA-CB-CG1	5.29	119.39	110.40
1	B	42	SER	N-CA-C	5.28	118.71	110.32
1	B	145	VAL	N-CA-C	5.26	120.27	109.34
1	A	174	HIS	N-CA-C	5.24	120.32	113.30
1	A	268	THR	N-CA-C	5.19	117.70	109.24
1	B	-1	ASN	CA-CB-CG	5.15	117.75	112.60
1	A	263	GLY	O-C-N	-5.14	116.02	122.70
1	B	116	GLU	CB-CG-CD	-5.13	103.88	112.60
1	B	293	PRO	N-CA-CB	5.13	108.05	103.08
1	A	168	GLY	CA-C-N	5.12	126.75	120.51
1	A	168	GLY	C-N-CA	5.12	126.75	120.51
1	B	297	PRO	N-CA-CB	5.11	108.22	102.60
1	B	97	THR	N-CA-CB	5.10	117.86	109.95
1	B	120	VAL	CA-CB-CG1	5.09	119.06	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	150	VAL	CA-CB-CG1	5.09	119.05	110.40
1	B	53	HIS	CA-CB-CG	-5.08	108.72	113.80
1	A	279	SER	N-CA-CB	5.07	119.06	110.49
1	B	26	THR	N-CA-C	5.06	117.87	109.46
1	B	95	THR	CA-CB-CG2	5.06	119.11	110.50
1	B	145	VAL	CA-CB-CG1	5.06	119.00	110.40
1	B	66	HIS	N-CA-C	5.05	117.56	110.23
1	B	0	THR	CA-C-N	5.02	130.58	122.29
1	B	0	THR	C-N-CA	5.02	130.58	122.29
1	B	281	LEU	CA-C-O	-5.01	114.97	120.43

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	95	THR	CB

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	A	2576	0	2506	98	0
1	B	2551	0	2467	135	0
2	A	14	0	13	0	0
2	B	14	0	13	0	0
3	A	36	0	49	0	0
3	B	36	0	49	6	0
4	A	206	0	0	14	1
4	B	160	0	0	10	1
All	All	5593	0	5097	233	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 23.

All (233) 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)
1:B:46(A):ARG:HG3	1:B:46(A):ARG:HH11	1.26	0.97
1:A:156:ASN:HD22	1:A:157:ARG:H	1.18	0.92
1:A:25:GLN:HE22	1:A:57:ASP:H	1.23	0.86
1:A:185:ILE:HG22	1:A:186:LYS:HG2	1.57	0.86
1:B:49:ALA:CB	1:B:107:MET:HG2	2.10	0.80
1:A:226:SER:HA	1:A:229:GLU:HG2	1.64	0.79
1:B:20:ILE:HG12	1:B:89:ILE:HG12	1.63	0.78
1:B:253:PRO:HB2	4:B:1160:HOH:O	1.84	0.77
1:B:47:TYR:OH	1:B:106:GLU:HG3	1.86	0.76
1:B:289:MET:HG2	3:B:391:03D:H16A	1.67	0.76
1:B:314:ASP:OD2	1:B:319:ARG:HD2	1.87	0.75
1:B:46(A):ARG:HG3	1:B:46(A):ARG:NH1	1.94	0.74
1:A:211:LEU:HB2	1:A:298:THR:HG23	1.68	0.74
1:B:41:PRO:HB2	1:B:55:LEU:HD23	1.70	0.74
1:B:71:LEU:HD11	1:B:84:LEU:HD13	1.69	0.73
1:B:8:ASN:HD21	1:B:11:ASP:H	1.36	0.73
1:A:289:MET:HB3	4:A:1094:HOH:O	1.89	0.72
1:A:156:ASN:HD22	1:A:157:ARG:N	1.87	0.71
1:A:8:ASN:HD21	1:A:11:ASP:H	1.37	0.71
1:A:283:THR:HG22	1:A:284:LEU:N	2.07	0.70
1:B:261:HIS:HB3	4:B:872:HOH:O	1.90	0.70
1:A:202:SER:OG	1:A:204:LEU:HD23	1.91	0.70
1:B:279(B):SER:HB2	1:B:279(D):LYS:H	1.57	0.69
1:B:64:TYR:OH	1:B:66:HIS:HA	1.93	0.68
1:B:51:VAL:HG12	1:B:52:TYR:CD1	2.29	0.68
1:B:51:VAL:HG12	1:B:52:TYR:CE1	2.29	0.67
1:A:252:GLY:HA3	1:A:277:GLN:HE22	1.58	0.67
1:B:160(D):SER:N	4:B:937:HOH:O	2.27	0.67
1:B:42:SER:O	1:B:55:LEU:HD22	1.94	0.67
3:B:391:03D:H3	3:B:391:03D:H25A	1.77	0.67
1:A:156:ASN:ND2	1:A:157:ARG:H	1.90	0.66
1:B:193:GLN:HG2	1:B:195:LYS:NZ	2.10	0.66
1:A:195:LYS:NZ	1:A:263:GLY:H	1.93	0.66
1:B:49:ALA:HB3	1:B:107:MET:HG2	1.78	0.65
1:A:326:ARG:NH1	4:A:1110:HOH:O	2.29	0.65
1:B:77:THR:HB	1:B:111:PRO:HG3	1.79	0.65
1:A:187:THR:OG1	1:A:318:ASN:ND2	2.29	0.65
1:B:225:THR:O	1:B:229:GLU:HG3	1.97	0.64
1:A:270:THR:HG21	4:B:956:HOH:O	1.96	0.64
1:A:315:ARG:NH2	4:A:982:HOH:O	2.30	0.64
1:B:64:TYR:CZ	1:B:66:HIS:HA	2.32	0.63
1:B:226:SER:O	1:B:230:LYS:HG3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:ASN:ND2	1:B:11:ASP:H	1.95	0.63
1:B:193:GLN:HG2	1:B:195:LYS:HZ3	1.63	0.63
1:B:235:LEU:HD12	1:B:256:PRO:HD2	1.80	0.63
1:B:283:THR:C	1:B:284:LEU:HD23	2.24	0.62
1:B:68:GLY:O	1:B:69:THR:C	2.39	0.62
1:A:244:ASP:OD2	1:A:283:THR:HG23	2.00	0.62
1:A:147:LYS:HD2	1:A:148:GLU:HG3	1.81	0.62
1:A:160(A):ASN:N	4:A:898:HOH:O	2.32	0.62
1:A:226:SER:O	1:A:229:GLU:HG3	1.99	0.62
1:A:174:HIS:HE1	4:A:1005:HOH:O	1.83	0.62
1:A:211:LEU:CB	1:A:298:THR:HG23	2.30	0.62
1:B:213:LEU:HD11	3:B:391:03D:H35A	1.82	0.61
1:B:49:ALA:HB1	1:B:107:MET:HG2	1.83	0.61
1:A:271:SER:N	4:A:892:HOH:O	2.33	0.60
1:A:133:VAL:HG23	4:A:837:HOH:O	2.01	0.60
1:B:8:ASN:HD21	1:B:11:ASP:N	2.00	0.59
1:A:8:ASN:ND2	1:A:11:ASP:H	2.00	0.59
1:A:141:ILE:HD12	1:A:149:ASP:OD1	2.02	0.59
1:B:173:GLN:HA	1:B:326:ARG:HH21	1.67	0.58
1:A:8:ASN:C	1:A:8:ASN:HD22	2.12	0.58
1:A:195:LYS:HZ3	1:A:263:GLY:H	1.51	0.58
1:A:8:ASN:HD21	1:A:11:ASP:N	2.01	0.57
1:A:314:ASP:OD2	1:A:319:ARG:HD2	2.04	0.57
1:B:8:ASN:C	1:B:8:ASN:HD22	2.12	0.57
1:B:82:GLY:HA3	1:B:103:GLU:O	2.04	0.57
1:B:140:ILE:O	1:B:143:GLN:HB3	2.05	0.57
1:A:156:ASN:ND2	1:A:157:ARG:N	2.50	0.57
1:A:270:THR:HG23	4:B:875:HOH:O	2.03	0.57
1:A:287:HIS:HE1	4:A:1096:HOH:O	1.87	0.57
1:A:185:ILE:CG2	1:A:186:LYS:HG2	2.34	0.56
1:B:111:PRO:HD2	1:B:112:PHE:CE2	2.39	0.56
1:B:183:ASN:HD22	1:B:319:ARG:HB3	1.70	0.56
1:A:283:THR:HG22	1:A:284:LEU:H	1.70	0.56
1:A:138:ASP:OD1	1:A:315:ARG:NH2	2.39	0.56
1:A:-1:ASN:O	1:A:-1:ASN:ND2	2.38	0.56
1:B:42:SER:C	1:B:55:LEU:HB3	2.31	0.56
1:B:316:ARG:HB3	4:B:953:HOH:O	2.05	0.55
1:A:199:VAL:CG2	1:A:231:LEU:HD12	2.37	0.55
1:A:240:ARG:HD2	1:A:246:VAL:HG13	1.87	0.55
1:A:71:LEU:HD23	1:A:71:LEU:N	2.21	0.55
1:A:252:GLY:HA3	1:A:277:GLN:NE2	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:GLY:C	1:B:254:THR:H	2.16	0.54
1:A:46(A):ARG:HD2	4:A:874:HOH:O	2.07	0.54
1:B:41:PRO:CB	1:B:55:LEU:HD23	2.36	0.54
1:A:277:GLN:HG2	4:A:969:HOH:O	2.07	0.54
1:A:185:ILE:HG22	1:A:186:LYS:CG	2.34	0.54
1:B:125:PHE:CE2	1:B:315:ARG:HD3	2.43	0.54
1:B:48:THR:HG22	4:B:1069:HOH:O	2.08	0.53
1:B:11:ASP:OD1	1:B:158:ASP:HB2	2.08	0.53
1:B:204:LEU:HD13	1:B:205:LEU:HG	1.91	0.53
1:B:51:VAL:CG1	1:B:52:TYR:CE1	2.92	0.53
1:B:161:LEU:HD23	1:B:161:LEU:O	2.08	0.53
1:B:60:ASP:OD1	1:B:60:ASP:N	2.38	0.52
1:B:195:LYS:HG3	4:B:1065:HOH:O	2.09	0.52
1:B:42:SER:O	1:B:55:LEU:HB3	2.08	0.52
1:A:185:ILE:HG22	1:A:186:LYS:N	2.22	0.52
1:B:86:GLN:OE1	1:B:97:THR:HG22	2.10	0.52
1:A:84:LEU:HB3	1:A:100:MET:HE1	1.91	0.52
1:A:315:ARG:O	1:A:316:ARG:C	2.49	0.51
1:A:226:SER:HA	1:A:229:GLU:CG	2.39	0.51
1:B:73:LEU:O	1:B:79:THR:HA	2.10	0.51
1:A:240:ARG:HD3	1:A:244:ASP:OD1	2.11	0.50
1:B:199:VAL:HG13	1:B:256:PRO:HG2	1.94	0.50
1:A:63:SER:CB	1:A:88:ILE:HD12	2.42	0.50
1:A:43:SER:O	1:A:46(A):ARG:NH2	2.44	0.50
1:B:150:VAL:HG13	1:B:312:GLU:HG3	1.92	0.49
1:B:200:GLY:HA2	4:B:1164:HOH:O	2.11	0.49
1:B:144:GLY:O	1:B:145:VAL:HB	2.11	0.49
1:B:289:MET:HG2	3:B:391:03D:C16	2.40	0.49
1:B:161:LEU:HD23	1:B:161:LEU:N	2.27	0.49
1:B:205:LEU:HD21	1:B:230:LYS:HB2	1.93	0.49
1:A:147:LYS:NZ	1:A:148:GLU:OE2	2.42	0.49
1:A:84:LEU:HD13	1:A:100:MET:CE	2.43	0.49
1:A:8:ASN:HD21	1:A:11:ASP:HA	1.78	0.49
1:B:293:PRO:HA	1:B:294:PRO:C	2.36	0.49
1:A:253:PRO:HD3	1:A:277:GLN:NE2	2.27	0.49
1:B:130:ILE:HD12	1:B:130:ILE:N	2.28	0.48
1:B:271:SER:HA	1:B:274:TYR:CE2	2.47	0.48
1:B:203:THR:HG22	1:B:207:GLU:OE1	2.13	0.48
1:A:90:THR:HA	1:A:94:ILE:O	2.13	0.48
1:B:39:TRP:NE1	1:B:120:VAL:HG13	2.29	0.48
1:B:296:GLY:HA2	1:B:297:PRO:C	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:HIS:HE1	1:B:263:GLY:O	1.96	0.47
1:B:199:VAL:CG1	1:B:256:PRO:HG2	2.44	0.47
1:B:199:VAL:O	1:B:200:GLY:C	2.52	0.47
1:A:248:LYS:HE3	1:A:251:GLU:HG3	1.95	0.47
1:A:2:SER:OG	1:A:92:GLY:O	2.29	0.47
1:A:63:SER:HB2	1:A:88:ILE:HD12	1.97	0.47
1:B:156:ASN:HD22	1:B:157:ARG:H	1.63	0.47
1:A:133:VAL:HG23	1:A:133:VAL:O	2.15	0.47
1:B:198:SER:OG	1:B:261:HIS:HE1	1.98	0.46
1:B:252:GLY:HA3	1:B:277:GLN:HE22	1.80	0.46
1:B:-1:ASN:HA	1:B:147:LYS:HG3	1.97	0.46
1:A:8:ASN:ND2	1:A:8:ASN:C	2.73	0.46
1:B:10:MET:O	1:B:11:ASP:HB2	2.16	0.46
1:A:184:LEU:O	1:A:185:ILE:C	2.56	0.46
1:B:99:GLN:NE2	1:B:136:ILE:HA	2.30	0.46
1:A:248:LYS:O	1:A:251:GLU:HB2	2.16	0.46
1:B:107:MET:HG2	1:B:107:MET:O	2.16	0.46
1:B:130:ILE:N	1:B:130:ILE:CD1	2.79	0.46
1:B:233:GLU:O	1:B:233:GLU:HG3	2.14	0.46
1:A:148:GLU:O	1:A:168:GLY:HA2	2.16	0.45
1:A:3:SER:HA	1:A:165:ILE:O	2.17	0.45
1:B:161:LEU:N	1:B:161:LEU:CD2	2.79	0.45
1:B:261:HIS:ND1	1:B:261:HIS:N	2.64	0.45
1:A:186:LYS:HB3	1:A:186:LYS:HE3	1.65	0.45
1:A:316:ARG:CZ	1:A:317:ASN:ND2	2.79	0.45
1:B:204:LEU:HB3	1:B:205:LEU:HG	1.99	0.45
1:B:51:VAL:C	1:B:52:TYR:CD1	2.95	0.45
1:B:161:LEU:HD23	1:B:161:LEU:C	2.42	0.45
1:B:97:THR:HG21	4:B:1016:HOH:O	2.17	0.45
1:A:183:ASN:OD1	1:A:183:ASN:N	2.49	0.45
1:A:278:GLU:O	1:A:279:SER:HB3	2.16	0.45
1:A:283:THR:CG2	1:A:284:LEU:N	2.78	0.45
1:B:279(B):SER:OG	1:B:280:LYS:HB2	2.17	0.45
1:B:284:LEU:HD23	1:B:284:LEU:N	2.29	0.45
1:B:-1:ASN:C	1:B:147:LYS:HG3	2.42	0.44
1:A:276:PHE:CE2	1:A:285:ALA:HB2	2.52	0.44
1:B:48:THR:O	1:B:52:TYR:N	2.49	0.44
1:B:111:PRO:HB2	3:B:391:03D:H8	1.99	0.44
1:B:181:TYR:CD2	1:B:319:ARG:HD3	2.51	0.44
1:A:8:ASN:HD21	1:A:11:ASP:CA	2.30	0.44
1:A:128:GLN:HG3	4:A:902:HOH:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:GLN:O	1:A:195:LYS:NZ	2.48	0.44
1:B:152:SER:O	1:B:165:ILE:HA	2.17	0.44
1:B:157:ARG:HG2	1:B:307:ARG:HD2	1.98	0.44
1:B:21:GLY:O	1:B:24:PRO:HA	2.17	0.44
1:B:43:SER:HB3	1:B:57:ASP:HA	2.00	0.44
1:A:253:PRO:HD3	1:A:277:GLN:HE22	1.82	0.44
1:A:280:LYS:HD3	1:A:280:LYS:HA	1.52	0.44
1:A:84:LEU:HD13	1:A:100:MET:HE1	2.00	0.43
1:A:100:MET:HE3	1:A:100:MET:HB3	1.55	0.43
1:A:181:TYR:HD1	4:A:859:HOH:O	2.00	0.43
1:A:326:ARG:HH12	1:B:1:THR:HG22	1.84	0.43
1:B:46(A):ARG:HB3	1:B:51:VAL:HG22	2.00	0.43
1:B:228:ILE:O	1:B:232:MET:HG2	2.18	0.43
1:B:277:GLN:C	1:B:279:SER:H	2.25	0.43
1:A:74:ARG:HD3	4:A:1149:HOH:O	2.18	0.43
1:A:316:ARG:CZ	1:A:317:ASN:HD21	2.31	0.43
1:B:111:PRO:HB2	3:B:391:03D:C8	2.49	0.43
1:B:125:PHE:HB3	1:B:127:GLU:OE1	2.19	0.43
1:A:249:CYS:HB3	1:A:279:SER:O	2.19	0.43
1:B:69:THR:HB	1:B:84:LEU:HD23	2.00	0.43
1:A:262:LEU:O	1:A:263:GLY:C	2.62	0.43
1:B:49:ALA:HB1	1:B:107:MET:CG	2.48	0.43
1:B:52:TYR:CD1	1:B:52:TYR:N	2.86	0.42
1:B:240:ARG:N	1:B:244:ASP:O	2.49	0.42
1:B:51:VAL:C	1:B:52:TYR:HD1	2.27	0.42
1:B:82:GLY:HA2	1:B:105:THR:HB	2.01	0.42
1:B:111:PRO:HD2	1:B:112:PHE:CD2	2.55	0.42
1:A:248:LYS:O	1:A:249:CYS:C	2.62	0.42
1:B:80:VAL:HG13	1:B:104:VAL:CG1	2.50	0.42
1:B:172:PRO:HA	1:B:175:TYR:CE1	2.54	0.42
1:B:198:SER:OG	1:B:261:HIS:CE1	2.73	0.42
1:B:0:THR:O	1:B:146:LEU:HA	2.20	0.42
1:B:286:ILE:C	1:B:287:HIS:HD2	2.27	0.42
1:A:174:HIS:HD2	1:A:326:ARG:OXT	2.02	0.41
1:A:215:ASP:O	1:A:302:GLY:HA2	2.18	0.41
1:A:242:LEU:HD12	1:A:242:LEU:HA	1.80	0.41
1:A:278:GLU:HG3	4:A:992:HOH:O	2.19	0.41
1:B:64:TYR:CG	1:B:65:LYS:N	2.88	0.41
1:B:134:THR:HA	1:B:135:PRO:HD3	1.91	0.41
1:B:8:ASN:ND2	1:B:8:ASN:C	2.76	0.41
1:A:8:ASN:HD22	1:A:9:TYR:N	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:VAL:HG22	1:B:314:ASP:HA	2.02	0.41
1:B:9:TYR:CD2	1:B:116:GLU:HG3	2.56	0.41
1:B:89:ILE:HD12	1:B:99:GLN:HG2	2.03	0.41
1:B:156:ASN:ND2	1:B:157:ARG:H	2.19	0.41
1:B:5:ILE:HD12	1:B:5:ILE:HG23	1.92	0.41
1:B:48:THR:HG22	1:B:52:TYR:CE1	2.56	0.41
1:B:107:MET:HE2	1:B:112:PHE:HB3	2.03	0.41
1:B:125:PHE:CE2	1:B:315:ARG:CD	3.04	0.41
1:B:165:ILE:HG21	1:B:165:ILE:HD13	1.89	0.41
1:B:252:GLY:C	1:B:254:THR:N	2.79	0.41
1:A:141:ILE:HD12	1:A:141:ILE:HG23	1.82	0.41
1:A:324:LEU:HD23	1:A:324:LEU:HA	1.83	0.41
1:B:252:GLY:O	1:B:254:THR:N	2.54	0.41
1:B:277:GLN:C	1:B:279:SER:N	2.78	0.41
1:A:204:LEU:HD12	1:A:205:LEU:HG	2.01	0.40
1:A:199:VAL:HG22	1:A:231:LEU:HD12	2.03	0.40
1:A:202:SER:O	1:A:204:LEU:N	2.54	0.40
1:B:48:THR:O	1:B:51:VAL:N	2.54	0.40
1:B:286:ILE:C	1:B:287:HIS:CD2	2.99	0.40
1:B:252:GLY:N	1:B:253:PRO:HD2	2.37	0.40
1:A:239:LYS:HG3	1:A:245:TYR:CZ	2.57	0.40
1:B:8:ASN:HD21	1:B:11:ASP:HA	1.86	0.40
1:B:198:SER:O	1:B:258:ILE:HA	2.21	0.40
1:A:96:VAL:HG11	1:A:140:ILE:HG13	2.04	0.40
1:B:46(A):ARG:HB3	1:B:51:VAL:CG2	2.52	0.40
1:B:194:MET:O	1:B:209:GLY:HA2	2.22	0.40
1:B:248:LYS:HB3	1:B:250:ASN:OD1	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:927:HOH:O	4:B:917:HOH:O[6_456]	2.11	0.09

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	331/337 (98%)	315 (95%)	14 (4%)	2 (1%)	21	11
1	B	330/337 (98%)	315 (96%)	12 (4%)	3 (1%)	14	5
All	All	661/674 (98%)	630 (95%)	26 (4%)	5 (1%)	16	6

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	279	SER
1	A	203	THR
1	B	242	LEU
1	B	112	PHE
1	B	253	PRO

5.3.2 Protein sidechains [i](#)

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	284/287 (99%)	244 (86%)	40 (14%)	3	1
1	B	280/287 (98%)	236 (84%)	44 (16%)	2	0
All	All	564/574 (98%)	480 (85%)	84 (15%)	3	0

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	ASN
1	A	1	THR
1	A	5	ILE
1	A	8	ASN
1	A	17	GLU
1	A	26	THR
1	A	28	LYS

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Mol	Chain	Res	Type
1	A	96	VAL
1	A	100	MET
1	A	104	VAL
1	A	138	ASP
1	A	147	LYS
1	A	149	ASP
1	A	173	GLN
1	A	186	LYS
1	A	187	THR
1	A	195	LYS
1	A	198	SER
1	A	204	LEU
1	A	207	GLU
1	A	208	ASP
1	A	227	SER
1	A	229	GLU
1	A	238	LYS
1	A	239	LYS
1	A	247	VAL
1	A	248	LYS
1	A	254	THR
1	A	255	LEU
1	A	259	SER
1	A	266	GLU
1	A	270	THR
1	A	271	SER
1	A	284	LEU
1	A	298	THR
1	A	315	ARG
1	A	316	ARG
1	A	319	ARG
1	A	324	LEU
1	A	326	ARG
1	B	1	THR
1	B	8	ASN
1	B	42	SER
1	B	43	SER
1	B	46	SER
1	B	46(A)	ARG
1	B	46(B)	LEU
1	B	47	TYR
1	B	48	THR

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Mol	Chain	Res	Type
1	B	55	LEU
1	B	59	SER
1	B	60	ASP
1	B	62	SER
1	B	65	LYS
1	B	70	GLU
1	B	73	LEU
1	B	80	VAL
1	B	84	LEU
1	B	86	GLN
1	B	91	VAL
1	B	95	THR
1	B	97	THR
1	B	105	THR
1	B	110	LEU
1	B	120	VAL
1	B	126	ILE
1	B	133	VAL
1	B	141	ILE
1	B	145	VAL
1	B	150	VAL
1	B	161	LEU
1	B	189	VAL
1	B	204	LEU
1	B	207	GLU
1	B	213	LEU
1	B	226	SER
1	B	233	GLU
1	B	235	LEU
1	B	240	ARG
1	B	248	LYS
1	B	254	THR
1	B	261	HIS
1	B	265	LYS
1	B	319	ARG

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	-1	ASN
1	A	8	ASN
1	A	25	GLN

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Mol	Chain	Res	Type
1	A	99	GLN
1	A	128	GLN
1	A	139	ASN
1	A	156	ASN
1	A	173	GLN
1	A	174	HIS
1	A	277	GLN
1	A	287	HIS
1	A	317	ASN
1	A	318	ASN
1	B	8	ASN
1	B	25	GLN
1	B	66	HIS
1	B	99	GLN
1	B	128	GLN
1	B	139	ASN
1	B	143	GLN
1	B	156	ASN
1	B	164	GLN
1	B	174	HIS
1	B	180	HIS
1	B	183	ASN
1	B	261	HIS
1	B	277	GLN
1	B	287	HIS
1	B	318	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

4 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
3	03D	B	391	-	38,38,38	1.57	7 (18%)	42,51,51	2.33	6 (14%)
2	NAG	B	367	1	14,14,15	1.63	2 (14%)	17,19,21	1.99	7 (41%)
2	NAG	A	367	1	14,14,15	0.94	1 (7%)	17,19,21	1.79	5 (29%)
3	03D	A	391	-	38,38,38	1.64	4 (10%)	42,51,51	1.83	7 (16%)

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
3	03D	B	391	-	-	2/40/50/50	0/3/3/3
2	NAG	B	367	1	-	3/6/23/26	0/1/1/1
2	NAG	A	367	1	-	0/6/23/26	0/1/1/1
3	03D	A	391	-	-	0/40/50/50	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	391	03D	C20-C21	4.63	1.61	1.53
3	A	391	03D	C22-C21	4.05	1.58	1.53
3	B	391	03D	C18-C15	3.71	1.59	1.53
2	B	367	NAG	C1-C2	3.70	1.57	1.52
3	A	391	03D	C3-C4	3.67	1.60	1.51
2	B	367	NAG	C8-C7	2.92	1.56	1.50
3	B	391	03D	C22-C21	2.69	1.56	1.53
3	B	391	03D	C3-C4	2.44	1.57	1.51
2	A	367	NAG	C8-C7	2.37	1.55	1.50
3	A	391	03D	C18-C13	2.22	1.58	1.53
3	B	391	03D	C9-C4	-2.19	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	391	03D	C27-C28	2.12	1.58	1.53
3	B	391	03D	C28-C23	2.03	1.58	1.52
3	B	391	03D	C6-C5	-2.03	1.35	1.38

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	391	03D	C18-C15-C16	-11.52	103.69	119.71
3	A	391	03D	C18-C15-C16	-7.29	109.56	119.71
3	B	391	03D	C18-C15-C17	-5.85	111.57	119.71
3	A	391	03D	C18-C15-C17	-5.13	112.58	119.71
2	A	367	NAG	C6-C5-C4	-3.56	104.29	113.02
2	B	367	NAG	O4-C4-C3	3.49	118.60	110.38
3	B	391	03D	C4-C3-C2	-3.27	107.44	113.39
2	A	367	NAG	C8-C7-N2	3.20	121.42	116.12
2	B	367	NAG	C4-C3-C2	-3.13	106.44	111.02
2	A	367	NAG	O7-C7-C8	-3.10	116.53	122.05
3	B	391	03D	C33-C32-C31	-3.09	110.53	114.82
2	B	367	NAG	C1-C2-N2	3.07	115.27	110.43
2	B	367	NAG	C6-C5-C4	-3.01	105.62	113.02
3	A	391	03D	C33-C32-C31	-3.01	110.64	114.82
3	B	391	03D	O19-C12-C13	-2.96	117.61	122.19
2	A	367	NAG	C4-C3-C2	-2.96	106.68	111.02
2	B	367	NAG	C1-O5-C5	-2.92	108.27	112.19
3	B	391	03D	C22-C21-C20	-2.86	107.60	112.18
3	A	391	03D	O36-C31-C32	2.82	114.79	109.15
3	A	391	03D	C22-C21-C20	-2.75	107.78	112.18
2	A	367	NAG	O5-C5-C4	-2.41	104.95	110.83
3	A	391	03D	O10-C1-C2	-2.35	104.63	109.57
2	B	367	NAG	C8-C7-N2	2.33	119.98	116.12
2	B	367	NAG	O5-C5-C4	-2.33	105.16	110.83
3	A	391	03D	C4-C3-C2	-2.16	109.45	113.39

There are no chirality outliers.

All (5) torsion outliers are listed below:

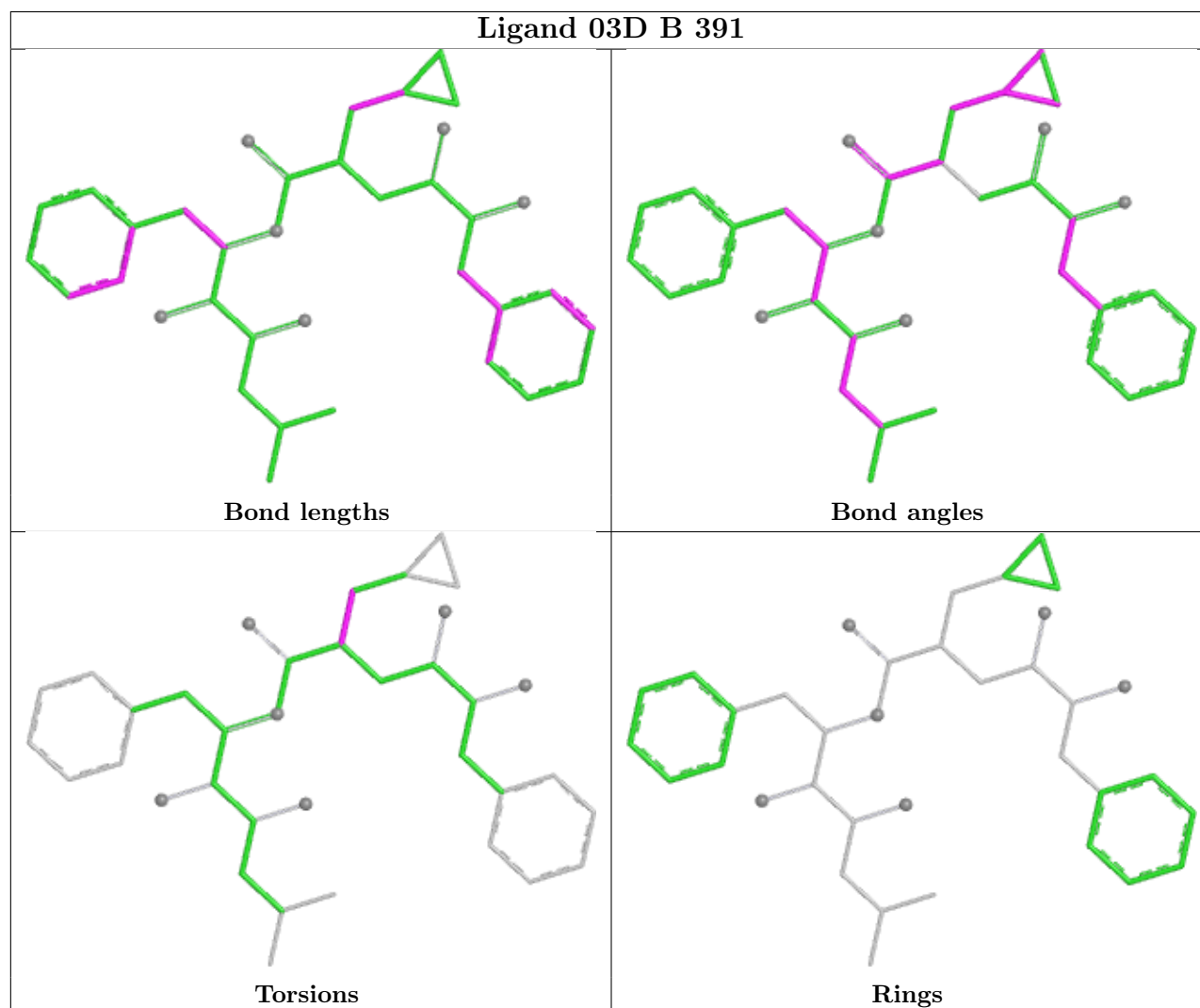
Mol	Chain	Res	Type	Atoms
3	B	391	03D	C12-C13-C18-C15
2	B	367	NAG	O7-C7-N2-C2
2	B	367	NAG	C8-C7-N2-C2
2	B	367	NAG	O5-C5-C6-O6
3	B	391	03D	C14-C13-C18-C15

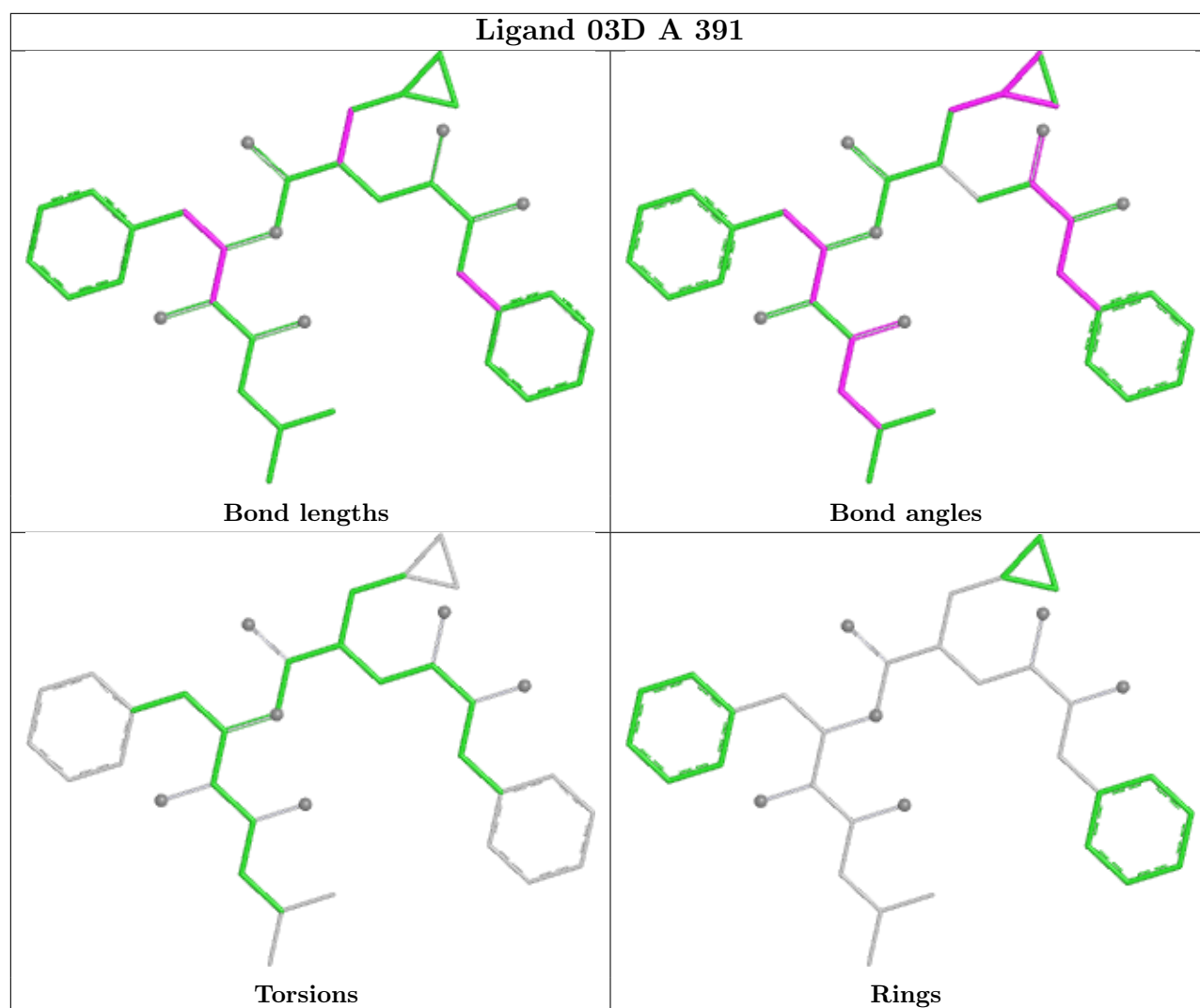
There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	391	03D	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.