Full wwPDB X-ray Structure Validation Report (i)

Mar 3, 2022 – 02:02 AM EST

PDB ID	:	1BIL
Title	:	CRYSTALLOGRAPHIC STUDIES ON THE BINDING MODES OF P2-P3 BUTANEDIAMIDE RENIN INHIBITORS
Authors	:	Tong, L.
Deposited on	:	1995-09-27
Resolution	:	2.40 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

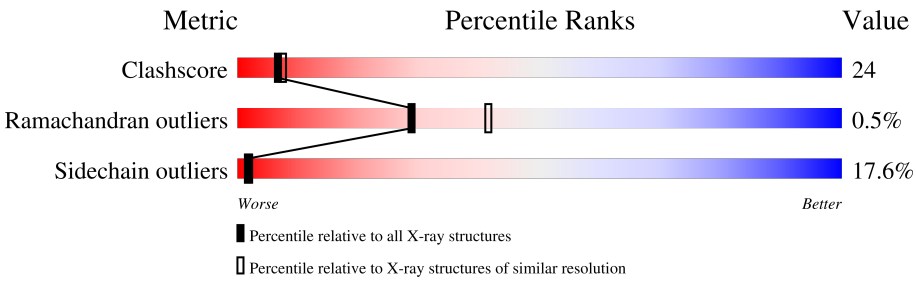
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.27

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	337	55% 30% 13% ··
1	B	337	51% 35% 10% · ·

2 Entry composition (i)

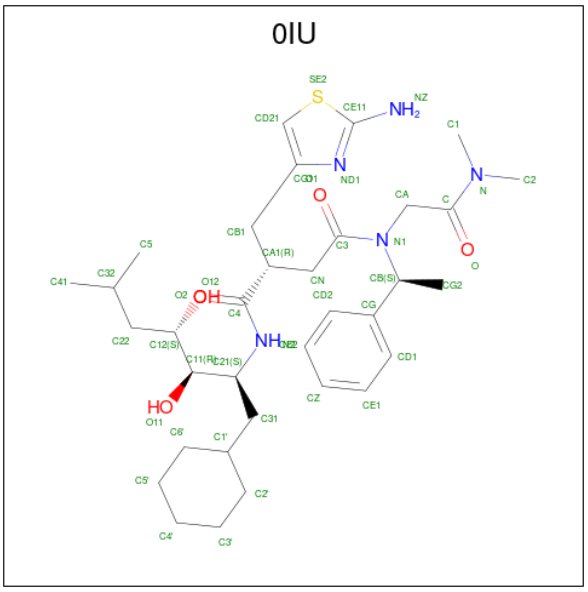
There are 3 unique types of molecules in this entry. The entry contains 5350 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
			Total	C	N	O	S			
1	A	334	2575	1644	416	501	14	0	0	0
1	B	333	2550	1630	408	498	14	0	0	0

- Molecule 2 is (2S)-2-[(2-amino-1,3-thiazol-4-yl)methyl]-N 1 -[(1S,2R,3R)-1-(cyclohexylmethyl)-2,3-dihydroxy-5-methylhexyl]-N 4 -[2-(d imethylamino)-2-oxoethyl]-N 4 -[(1S)-1-phenylethyl]butanediamide (three-letter code: 0IU) (formula: $C_{34}H_{53}N_5O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	45	34	5	5	1	0	0
2	B	1	45	34	5	5	1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	69	69	69	0	0
			Total	O		
3	B	66	66	66	0	0

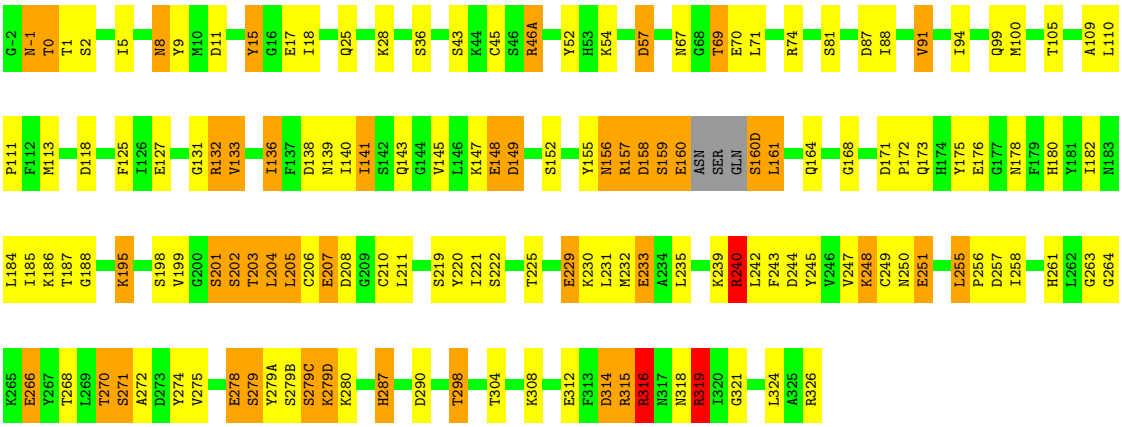
3 Residue-property plots (i)

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. 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.

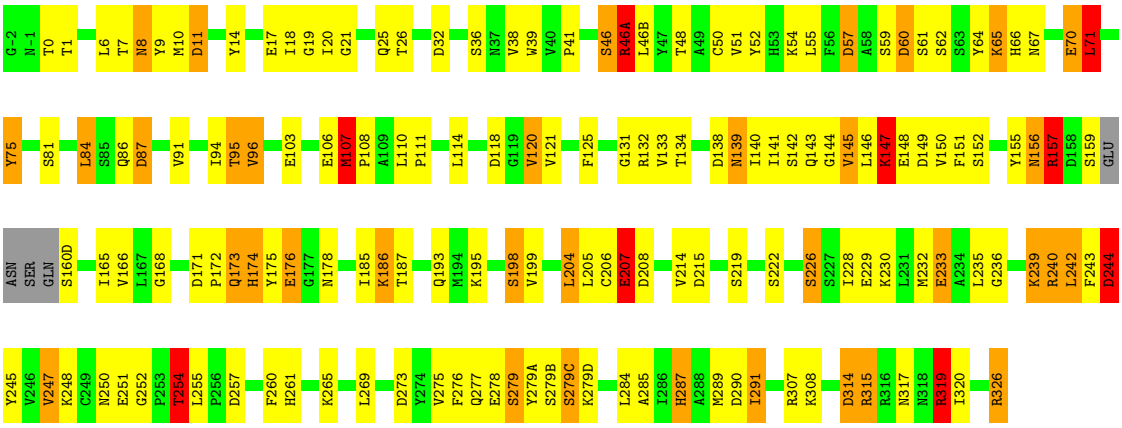
• Molecule 1: Renin

Chain A: 55% 30% 13% ••



• Molecule 1: Renin

Chain B: 51% 35% 10% • •



4 Data and refinement statistics (i)

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	143.10Å 143.10Å 143.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.170 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5350	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 0IU

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	1.13	13/2634 (0.5%)	1.46	33/3570 (0.9%)
1	B	1.12	11/2609 (0.4%)	1.50	41/3540 (1.2%)
All	All	1.12	24/5243 (0.5%)	1.48	74/7110 (1.0%)

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	278	GLU	CD-OE2	11.27	1.38	1.25
1	B	233	GLU	CD-OE1	10.21	1.36	1.25
1	A	229	GLU	CD-OE2	9.36	1.35	1.25
1	A	266	GLU	CD-OE2	9.32	1.35	1.25
1	B	229	GLU	CD-OE1	8.77	1.35	1.25
1	A	70	GLU	CD-OE1	8.54	1.35	1.25
1	A	278	GLU	CD-OE2	8.52	1.35	1.25
1	A	251	GLU	CD-OE1	8.21	1.34	1.25
1	B	17	GLU	CD-OE2	8.11	1.34	1.25
1	B	176	GLU	CD-OE2	8.11	1.34	1.25
1	A	160	GLU	CD-OE1	7.94	1.34	1.25
1	B	106	GLU	CD-OE2	7.86	1.34	1.25
1	A	233	GLU	CD-OE2	7.83	1.34	1.25
1	B	148	GLU	CD-OE1	7.54	1.33	1.25
1	A	207	GLU	CD-OE1	7.28	1.33	1.25
1	A	148	GLU	CD-OE2	6.83	1.33	1.25
1	B	70	GLU	CD-OE2	6.78	1.33	1.25
1	A	17	GLU	CD-OE2	6.41	1.32	1.25
1	B	103	GLU	CD-OE2	6.40	1.32	1.25
1	A	176	GLU	CD-OE1	5.61	1.31	1.25
1	B	157	ARG	NE-CZ	5.43	1.40	1.33
1	A	312	GLU	CD-OE1	-5.33	1.19	1.25
1	A	127	GLU	CD-OE1	-5.09	1.20	1.25
1	B	207	GLU	CD-OE1	5.08	1.31	1.25

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	319	ARG	NE-CZ-NH1	12.87	126.73	120.30
1	A	319	ARG	NE-CZ-NH2	-12.32	114.14	120.30
1	B	319	ARG	NE-CZ-NH1	10.25	125.42	120.30
1	A	158	ASP	CB-CG-OD2	-9.72	109.55	118.30
1	A	74	ARG	NE-CZ-NH1	9.63	125.11	120.30
1	B	149	ASP	CB-CG-OD2	-9.60	109.66	118.30
1	B	11	ASP	CB-CG-OD1	8.45	125.90	118.30
1	B	240	ARG	NE-CZ-NH2	-8.34	116.13	120.30
1	A	208	ASP	CB-CG-OD1	8.25	125.72	118.30
1	B	208	ASP	CB-CG-OD1	-8.14	110.97	118.30
1	B	257	ASP	CB-CG-OD1	7.76	125.28	118.30
1	B	60	ASP	CB-CG-OD2	-7.52	111.53	118.30
1	B	290	ASP	CB-CG-OD1	7.46	125.01	118.30
1	A	57	ASP	CB-CG-OD2	7.23	124.81	118.30
1	A	158	ASP	CB-CG-OD1	7.16	124.75	118.30
1	B	46(A)	ARG	NE-CZ-NH2	-7.16	116.72	120.30
1	A	118	ASP	CB-CG-OD1	7.10	124.69	118.30
1	B	147	LYS	N-CA-CB	7.05	123.30	110.60
1	A	257	ASP	CB-CG-OD1	7.01	124.61	118.30
1	B	314	ASP	CB-CG-OD2	-6.91	112.08	118.30
1	B	279(A)	TYR	CB-CG-CD1	-6.89	116.87	121.00
1	B	87	ASP	CB-CG-OD1	6.81	124.43	118.30
1	B	290	ASP	CB-CG-OD2	-6.76	112.21	118.30
1	B	257	ASP	CB-CG-OD2	-6.73	112.24	118.30
1	A	220	TYR	CB-CG-CD2	-6.68	117.00	121.00
1	B	307	ARG	NE-CZ-NH1	6.64	123.62	120.30
1	A	149	ASP	CB-CG-OD1	-6.59	112.37	118.30
1	B	32	ASP	CB-CG-OD2	-6.57	112.38	118.30
1	B	245	TYR	CB-CG-CD2	-6.54	117.08	121.00
1	A	257	ASP	CB-CG-OD2	-6.49	112.46	118.30
1	B	11	ASP	CB-CG-OD2	-6.44	112.50	118.30
1	B	289	MET	CG-SD-CE	-6.44	89.90	100.20
1	A	118	ASP	CB-CG-OD2	-6.20	112.72	118.30
1	B	244	ASP	CB-CG-OD2	6.19	123.87	118.30
1	A	171	ASP	CB-CG-OD1	6.14	123.83	118.30
1	B	174	HIS	CA-CB-CG	-6.09	103.25	113.60
1	B	87	ASP	CB-CG-OD2	-6.09	112.82	118.30
1	B	118	ASP	CB-CG-OD2	-6.07	112.84	118.30
1	B	240	ARG	NE-CZ-NH1	5.97	123.28	120.30
1	A	314	ASP	CB-CG-OD2	-5.95	112.94	118.30
1	A	155	TYR	CB-CG-CD1	5.85	124.51	121.00
1	B	57	ASP	CB-CG-OD2	-5.81	113.07	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	208	ASP	CB-CG-OD2	-5.80	113.08	118.30
1	B	32	ASP	CB-CG-OD1	5.80	123.52	118.30
1	A	315	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	A	271	SER	N-CA-CB	-5.71	101.94	110.50
1	B	75	TYR	CB-CG-CD2	-5.66	117.61	121.00
1	B	273	ASP	CB-CG-OD2	-5.64	113.22	118.30
1	B	244	ASP	CB-CG-OD1	-5.59	113.27	118.30
1	B	84	LEU	N-CA-CB	-5.57	99.27	110.40
1	A	52	TYR	CB-CG-CD1	5.53	124.32	121.00
1	A	138	ASP	CB-CG-OD2	5.52	123.27	118.30
1	A	316	ARG	NE-CZ-NH1	5.51	123.05	120.30
1	B	71	LEU	N-CA-CB	5.51	121.41	110.40
1	B	208	ASP	CB-CG-OD2	5.43	123.19	118.30
1	B	198	SER	CB-CA-C	-5.37	99.89	110.10
1	A	314	ASP	CB-CG-OD1	5.37	123.13	118.30
1	A	52	TYR	CB-CG-CD2	-5.37	117.78	121.00
1	B	60	ASP	CB-CG-OD1	5.36	123.12	118.30
1	A	171	ASP	CB-CG-OD2	-5.33	113.50	118.30
1	A	87	ASP	CB-CG-OD1	5.28	123.05	118.30
1	A	240	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	50	CYS	CA-CB-SG	-5.26	104.52	114.00
1	A	87	ASP	CB-CG-OD2	-5.26	113.56	118.30
1	A	46(A)	ARG	NE-CZ-NH2	-5.23	117.68	120.30
1	A	220	TYR	CB-CG-CD1	5.20	124.12	121.00
1	A	290	ASP	CB-CG-OD1	5.16	122.94	118.30
1	B	155	TYR	CB-CG-CD2	-5.14	117.91	121.00
1	B	215	ASP	CB-CG-OD2	-5.14	113.68	118.30
1	A	15	TYR	CB-CG-CD2	-5.14	117.92	121.00
1	A	225	THR	CA-CB-CG2	-5.13	105.22	112.40
1	B	254	THR	N-CA-CB	-5.05	100.70	110.30
1	B	118	ASP	CB-CG-OD1	5.05	122.85	118.30
1	B	107	MET	N-CA-CB	5.02	119.64	110.60

There are no chirality outliers.

There are no planarity outliers.

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	2575	0	2507	105	0
1	B	2550	0	2468	117	0
2	A	45	0	53	10	0
2	B	45	0	53	15	0
3	A	69	0	0	4	0
3	B	66	0	0	7	0
All	All	5350	0	5081	242	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 24.

All (242) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:391:0IU:CD21	2:B:391:0IU:HD21	0.97	1.14
2:A:391:0IU:HD21	2:A:391:0IU:CD21	0.97	1.07
2:A:391:0IU:CG1	2:A:391:0IU:SE2	2.55	1.05
2:A:391:0IU:SE2	2:A:391:0IU:CE11	1.74	1.02
2:A:391:0IU:CD21	2:A:391:0IU:SE2	1.77	1.01
1:B:150:VAL:HG12	1:B:314:ASP:HA	1.38	1.01
1:A:-1:ASN:HD21	1:A:147:LYS:HE2	1.26	1.00
2:B:391:0IU:CD21	2:B:391:0IU:SE2	1.78	1.00
2:B:391:0IU:SE2	2:B:391:0IU:CG1	2.59	1.00
2:B:391:0IU:SE2	2:B:391:0IU:CE11	1.78	0.99
1:B:8:ASN:HD21	1:B:11:ASP:H	1.11	0.97
1:A:195:LYS:HZ3	1:A:263:GLY:H	1.14	0.95
1:B:176:GLU:HG3	1:B:326:ARG:HD3	1.47	0.94
1:A:8:ASN:HD21	1:A:11:ASP:H	1.16	0.93
1:B:6:LEU:HD21	1:B:165:ILE:HD12	1.53	0.91
2:A:391:0IU:SE2	2:A:391:0IU:ND1	2.56	0.89
2:A:391:0IU:HD21	2:A:391:0IU:CG1	2.03	0.89
1:B:252:GLY:HA3	1:B:277:GLN:HE22	1.38	0.88
2:B:391:0IU:HD21	2:B:391:0IU:CG1	2.03	0.87
1:A:195:LYS:NZ	1:A:263:GLY:H	1.72	0.86
2:B:391:0IU:SE2	2:B:391:0IU:ND1	2.59	0.85
1:A:156:ASN:ND2	1:A:157:ARG:H	1.75	0.85
1:B:48:THR:HB	1:B:52:TYR:HE2	1.42	0.85
1:B:291:ILE:HD12	2:B:391:0IU:H41	1.60	0.83
1:B:242:LEU:HB2	1:B:243:PHE:CD1	2.13	0.83
2:B:391:0IU:ND1	2:B:391:0IU:NZ	2.22	0.82
1:A:195:LYS:HE2	1:A:264:GLY:H	1.43	0.82
1:A:156:ASN:HD22	1:A:157:ARG:H	1.27	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:391:0IU:SE2	2:A:391:0IU:NZ	2.63	0.81
1:B:8:ASN:ND2	1:B:11:ASP:H	1.78	0.80
1:B:8:ASN:HD21	1:B:11:ASP:N	1.78	0.80
1:A:231:LEU:HD23	1:A:232:MET:HE3	1.67	0.76
1:B:314:ASP:OD2	1:B:319:ARG:HD3	1.85	0.76
2:A:391:0IU:ND1	2:A:391:0IU:NZ	2.29	0.76
2:B:391:0IU:SE2	2:B:391:0IU:NZ	2.68	0.75
1:A:8:ASN:HD21	1:A:11:ASP:N	1.84	0.74
1:A:231:LEU:HD23	1:A:232:MET:CE	2.16	0.74
1:B:242:LEU:HD21	3:B:928:HOH:O	1.87	0.74
1:A:279(D):LYS:O	1:A:280:LYS:HD3	1.88	0.74
1:B:67:ASN:HB3	1:B:84:LEU:O	1.87	0.74
1:A:195:LYS:HD2	1:A:261:HIS:O	1.88	0.73
1:A:8:ASN:ND2	1:A:11:ASP:H	1.86	0.71
1:A:156:ASN:ND2	1:A:157:ARG:N	2.39	0.71
1:A:25:GLN:HE22	1:A:57:ASP:H	1.35	0.71
1:A:279(B):SER:HB3	1:A:280:LYS:HG2	1.72	0.70
1:B:251:GLU:O	1:B:254:THR:HB	1.91	0.70
1:B:156:ASN:HD22	1:B:157:ARG:H	1.39	0.70
1:A:195:LYS:NZ	1:A:263:GLY:N	2.39	0.69
1:B:96:VAL:HG21	1:B:139:ASN:HB3	1.73	0.69
1:A:195:LYS:HZ3	1:A:263:GLY:N	1.87	0.68
1:B:275:VAL:HG22	1:B:284:LEU:HD22	1.75	0.68
1:B:21:GLY:HA2	1:B:87:ASP:OD1	1.94	0.68
1:A:239:LYS:HD2	1:A:240:ARG:O	1.94	0.68
1:B:317:ASN:HB2	1:B:319:ARG:HD2	1.75	0.67
1:B:46:SER:OG	1:B:46(B):LEU:HB2	1.94	0.67
1:B:242:LEU:HB2	1:B:243:PHE:HD1	1.58	0.67
2:B:391:0IU:HD21	2:B:391:0IU:SE2	2.46	0.66
2:A:391:0IU:HD21	2:A:391:0IU:SE2	2.45	0.66
1:A:202:SER:O	1:A:204:LEU:N	2.29	0.65
1:A:239:LYS:O	1:A:239:LYS:HG3	1.97	0.65
1:B:140:ILE:O	1:B:143:GLN:HG2	1.96	0.65
1:B:84:LEU:HD11	1:B:133:VAL:HG11	1.79	0.65
1:B:240:ARG:O	1:B:243:PHE:N	2.30	0.65
1:B:46(A):ARG:HB3	1:B:51:VAL:CG2	2.26	0.65
1:A:205:LEU:CD2	1:A:230:LYS:HE2	2.27	0.64
1:A:195:LYS:HD2	1:A:195:LYS:H	1.64	0.63
1:A:240:ARG:NH1	1:A:244:ASP:OD2	2.32	0.62
1:A:136:ILE:O	1:A:140:ILE:HG13	1.99	0.62
1:B:18:ILE:HG22	1:B:91:VAL:HG22	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:PRO:HB2	1:B:55:LEU:HD23	1.80	0.62
1:A:240:ARG:HB2	1:A:244:ASP:OD1	1.99	0.62
1:B:205:LEU:CD2	1:B:230:LYS:HD2	2.30	0.62
1:B:185:ILE:HD11	1:B:193:GLN:HB2	1.81	0.60
1:A:94:ILE:HG21	1:A:140:ILE:HG23	1.82	0.60
1:B:174:HIS:HE1	3:B:893:HOH:O	1.84	0.60
1:A:-1:ASN:ND2	1:A:147:LYS:HE2	2.08	0.60
1:A:187:THR:HG22	1:A:188:GLY:N	2.17	0.59
1:A:5:ILE:HG23	1:A:161:LEU:HD11	1.84	0.59
1:B:46(A):ARG:HB3	1:B:51:VAL:HG22	1.83	0.59
1:B:171:ASP:OD1	1:B:173:GLN:HB3	2.03	0.59
1:B:198:SER:OG	1:B:261:HIS:HE1	1.85	0.59
1:B:172:PRO:HA	1:B:175:TYR:CE1	2.38	0.59
1:B:111:PRO:O	1:B:114:LEU:HG	2.03	0.58
1:B:8:ASN:HD22	1:B:9:TYR:N	2.02	0.58
1:B:205:LEU:HD21	1:B:230:LYS:HD2	1.86	0.58
1:B:228:ILE:O	1:B:232:MET:HG2	2.04	0.57
1:A:221:ILE:HG13	1:A:304:THR:HB	1.87	0.57
1:B:25:GLN:HE22	1:B:57:ASP:H	1.50	0.57
1:A:156:ASN:HD22	1:A:157:ARG:N	1.98	0.57
1:B:226:SER:O	1:B:230:LYS:HE3	2.04	0.57
1:A:205:LEU:HD23	1:A:230:LYS:HE2	1.85	0.57
1:B:247:VAL:HG21	1:B:255:LEU:HD11	1.85	0.57
1:A:182:ILE:N	1:A:182:ILE:HD12	2.21	0.56
1:B:174:HIS:HD2	1:B:326:ARG:OXT	1.88	0.56
1:B:199:VAL:CG2	1:B:204:LEU:HD22	2.36	0.56
1:A:205:LEU:HD23	1:A:230:LYS:CE	2.36	0.56
1:A:279(B):SER:HB3	1:A:280:LYS:CG	2.36	0.56
1:A:28:LYS:HB2	1:A:28:LYS:NZ	2.22	0.55
1:A:210:CYS:HB2	3:A:858:HOH:O	2.08	0.54
1:B:57:ASP:HB3	1:B:60:ASP:OD1	2.07	0.54
1:B:240:ARG:O	1:B:242:LEU:C	2.46	0.54
1:A:184:LEU:HD12	1:A:318:ASN:ND2	2.22	0.54
1:B:26:THR:O	1:B:54:LYS:HE3	2.07	0.54
1:A:268:THR:O	1:A:308:LYS:NZ	2.40	0.54
1:A:43:SER:O	1:A:46(A):ARG:NH2	2.40	0.54
1:B:317:ASN:CB	1:B:319:ARG:HD2	2.38	0.53
1:A:148:GLU:HB2	1:A:168:GLY:O	2.08	0.53
1:B:243:PHE:CD1	1:B:243:PHE:N	2.77	0.53
1:B:252:GLY:HA3	1:B:277:GLN:NE2	2.18	0.53
1:B:8:ASN:HD21	1:B:11:ASP:CA	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:SER:OG	2:B:391:0IU:HA2	2.09	0.53
1:A:178:ASN:HB2	3:A:846:HOH:O	2.10	0.52
1:A:314:ASP:OD2	1:A:319:ARG:HD2	2.09	0.52
1:B:9:TYR:CZ	1:B:10:MET:HE3	2.45	0.52
1:B:176:GLU:HG3	1:B:326:ARG:CD	2.31	0.52
1:B:242:LEU:HD23	1:B:242:LEU:N	2.25	0.52
1:B:131:GLY:O	1:B:132:ARG:C	2.48	0.52
1:B:279:SER:HB2	1:B:279(B):SER:H	1.75	0.52
1:B:19:GLY:HA2	1:B:25:GLN:O	2.09	0.52
1:A:184:LEU:HD12	1:A:318:ASN:HD22	1.74	0.52
1:A:132:ARG:HB2	1:A:132:ARG:HH11	1.74	0.51
1:A:211:LEU:HB2	1:A:298:THR:HG23	1.92	0.51
1:A:205:LEU:HD21	1:A:230:LYS:HE2	1.92	0.51
1:B:248:LYS:NZ	1:B:279(C):SER:O	2.38	0.51
1:A:270:THR:CG2	1:A:272:ALA:HB3	2.41	0.50
1:B:84:LEU:HD11	1:B:133:VAL:CG1	2.41	0.50
1:A:211:LEU:CB	1:A:298:THR:HG23	2.42	0.50
1:A:67:ASN:OD1	1:A:69:THR:HB	2.12	0.50
1:B:240:ARG:HD2	3:B:859:HOH:O	2.12	0.50
1:B:7:THR:O	1:B:14:TYR:HA	2.12	0.50
1:B:156:ASN:ND2	1:B:157:ARG:H	2.08	0.49
1:A:205:LEU:HD23	1:A:230:LYS:NZ	2.27	0.49
1:A:249:CYS:HB3	1:A:279:SER:O	2.11	0.49
1:A:45:CYS:HB2	1:A:105:THR:HA	1.93	0.49
1:B:206:CYS:HB2	3:B:857:HOH:O	2.11	0.49
1:A:279(B):SER:CB	1:A:280:LYS:HG2	2.41	0.49
1:B:18:ILE:CG2	1:B:91:VAL:HG22	2.42	0.49
1:A:133:VAL:HG22	3:A:812:HOH:O	2.11	0.49
1:B:141:ILE:O	1:B:142:SER:C	2.51	0.49
1:B:173:GLN:HG3	3:B:905:HOH:O	2.12	0.49
1:B:276:PHE:CE2	1:B:285:ALA:HB2	2.48	0.49
1:B:6:LEU:HD21	1:B:165:ILE:CD1	2.34	0.49
1:B:222:SER:HA	1:B:287:HIS:O	2.13	0.49
1:B:110:LEU:HA	1:B:111:PRO:HA	1.57	0.48
1:A:99:GLN:NE2	1:A:139:ASN:OD1	2.46	0.48
1:B:64:TYR:CG	1:B:65:LYS:N	2.80	0.48
1:B:134:THR:HG22	1:B:138:ASP:HB2	1.94	0.48
1:A:8:ASN:HD22	1:A:9:TYR:N	2.11	0.48
1:A:125:PHE:CE2	1:A:187:THR:HG23	2.47	0.48
1:A:231:LEU:HD23	1:A:232:MET:HE2	1.95	0.48
1:A:248:LYS:O	1:A:251:GLU:HB2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:SER:O	1:B:165:ILE:HA	2.14	0.48
1:A:270:THR:HG22	1:A:272:ALA:HB3	1.94	0.47
1:B:291:ILE:CD1	2:B:391:0IU:H41	2.38	0.47
1:B:95:THR:O	1:B:95:THR:HG22	2.14	0.47
1:B:8:ASN:HD22	1:B:8:ASN:C	2.17	0.47
1:A:28:LYS:NZ	1:A:28:LYS:CB	2.78	0.47
1:A:18:ILE:HG22	1:A:91:VAL:HG13	1.96	0.47
1:A:131:GLY:O	1:A:132:ARG:HB2	2.14	0.47
2:B:391:0IU:HA3	2:B:391:0IU:HG22	1.71	0.47
1:B:0:THR:CB	1:B:145:VAL:HG22	2.45	0.47
1:B:199:VAL:HG23	1:B:204:LEU:HD22	1.97	0.46
2:B:391:0IU:HN3	2:B:391:0IU:HB	1.61	0.46
1:A:315:ARG:O	1:A:316:ARG:C	2.52	0.46
1:B:10:MET:O	1:B:11:ASP:HB2	2.15	0.46
1:B:232:MET:O	1:B:236:GLY:N	2.48	0.46
1:A:232:MET:HG3	1:A:245:TYR:CE1	2.51	0.46
1:A:250:ASN:OD1	1:A:279(C):SER:HA	2.16	0.45
1:B:243:PHE:HD1	1:B:243:PHE:H	1.61	0.45
1:A:125:PHE:CE2	1:A:187:THR:CG2	2.99	0.45
1:A:164:GLN:NE2	3:A:923:HOH:O	2.49	0.45
1:B:250:ASN:OD1	1:B:279(C):SER:HA	2.16	0.45
1:B:199:VAL:HG21	1:B:204:LEU:HD22	1.99	0.45
1:B:48:THR:HB	1:B:52:TYR:CE2	2.34	0.45
1:B:173:GLN:HG3	1:B:173:GLN:O	2.15	0.45
1:B:242:LEU:HD23	1:B:242:LEU:H	1.82	0.45
1:A:198:SER:O	1:A:258:ILE:HA	2.16	0.45
1:A:202:SER:O	1:A:204:LEU:HD23	2.17	0.45
1:B:244:ASP:HB3	3:B:884:HOH:O	2.16	0.45
1:A:160:GLU:OE2	1:B:159:SER:HB3	2.17	0.45
1:A:235:LEU:HD22	1:A:256:PRO:HD2	1.98	0.45
1:B:242:LEU:CB	1:B:243:PHE:CD1	2.96	0.45
1:B:214:VAL:HG21	1:B:320:ILE:HD13	1.99	0.44
1:A:111:PRO:HB2	2:A:391:0IU:CD1	2.46	0.44
1:B:285:ALA:HB3	3:B:915:HOH:O	2.17	0.44
1:B:239:LYS:NZ	1:B:244:ASP:O	2.48	0.44
1:B:66:HIS:CG	1:B:67:ASN:N	2.86	0.44
1:B:71:LEU:O	1:B:81:SER:HA	2.18	0.44
1:B:150:VAL:HG12	1:B:314:ASP:CA	2.28	0.44
1:A:8:ASN:HD22	1:A:8:ASN:C	2.20	0.44
1:A:222:SER:HA	1:A:287:HIS:O	2.18	0.44
1:B:39:TRP:HE1	1:B:75:TYR:HH	1.66	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:LYS:O	1:B:186:LYS:HG2	2.18	0.44
1:A:187:THR:CG2	1:A:188:GLY:N	2.81	0.43
1:A:250:ASN:HB3	1:A:279(A):TYR:O	2.18	0.43
1:B:206:CYS:SG	1:B:206:CYS:O	2.77	0.43
1:A:203:THR:HG22	1:A:203:THR:O	2.18	0.43
1:B:147:LYS:NZ	1:B:168:GLY:O	2.45	0.43
1:A:109:ALA:O	1:A:113:MET:HB2	2.19	0.43
1:A:187:THR:HG22	1:A:188:GLY:H	1.83	0.43
1:A:242:LEU:HD23	1:A:243:PHE:CD2	2.53	0.43
1:A:304:THR:O	1:A:304:THR:HG22	2.18	0.43
1:A:159:SER:HG	1:A:160(D):SER:N	2.18	0.42
1:A:185:ILE:O	1:A:185:ILE:HG22	2.19	0.42
1:A:195:LYS:HD3	1:A:261:HIS:CE1	2.54	0.42
1:B:8:ASN:HD21	1:B:11:ASP:HA	1.83	0.42
1:B:151:PHE:HA	1:B:166:VAL:O	2.18	0.42
1:A:242:LEU:HD23	1:A:243:PHE:CE2	2.54	0.42
1:B:133:VAL:HG23	1:B:133:VAL:O	2.19	0.42
1:B:156:ASN:HD22	1:B:157:ARG:N	2.11	0.42
1:A:172:PRO:HA	1:A:175:TYR:CE1	2.55	0.42
1:A:235:LEU:CD2	1:A:256:PRO:HD2	2.49	0.42
1:B:65:LYS:HD3	1:B:65:LYS:HA	1.75	0.42
1:A:8:ASN:HD21	1:A:11:ASP:CA	2.32	0.42
1:B:143:GLN:HG3	1:B:145:VAL:HG12	2.02	0.42
1:B:48:THR:O	1:B:52:TYR:CD2	2.72	0.42
1:B:61:SER:HB3	1:B:64:TYR:HB2	2.01	0.42
1:A:206:CYS:O	1:A:206:CYS:SG	2.78	0.42
1:B:39:TRP:NE1	1:B:120:VAL:HG13	2.35	0.42
1:A:187:THR:HG23	1:A:318:ASN:HD21	1.85	0.42
1:A:255:LEU:HB3	1:A:274:TYR:OH	2.19	0.42
1:B:38:VAL:HG13	1:B:121:VAL:HG22	2.01	0.41
1:B:107:MET:HA	1:B:108:PRO:HD2	1.83	0.41
1:A:204:LEU:HB3	1:A:205:LEU:HG	2.02	0.41
2:B:391:0IU:CA	2:B:391:0IU:HD1	2.51	0.41
1:B:1:THR:OG1	1:B:147:LYS:HD2	2.21	0.41
1:A:0:THR:OG1	1:A:145:VAL:O	2.39	0.41
1:A:28:LYS:HB2	1:A:28:LYS:HZ2	1.86	0.41
1:A:180:HIS:O	1:A:321:GLY:HA2	2.21	0.41
1:B:144:GLY:O	1:B:146:LEU:N	2.54	0.41
1:B:239:LYS:NZ	1:B:242:LEU:O	2.45	0.41
1:B:260:PHE:HE2	1:B:269:LEU:HD22	1.86	0.41
1:A:8:ASN:HD21	1:A:11:ASP:HA	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:THR:OG1	1:A:318:ASN:ND2	2.55	0.40
1:A:199:VAL:O	1:A:201:SER:N	2.54	0.40
1:A:271:SER:O	1:A:275:VAL:HG23	2.21	0.40
1:A:15:TYR:CE2	1:A:28:LYS:HD2	2.56	0.40
1:B:8:ASN:ND2	1:B:8:ASN:C	2.74	0.40
1:B:125:PHE:CE1	1:B:315:ARG:HD3	2.56	0.40
1:B:207:GLU:H	1:B:207:GLU:HG2	1.28	0.40
1:A:5:ILE:HG23	1:A:161:LEU:CD1	2.50	0.40
1:A:141:ILE:N	1:A:141:ILE:HD13	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	330/337 (98%)	310 (94%)	19 (6%)	1 (0%)	41	55
1	B	329/337 (98%)	313 (95%)	14 (4%)	2 (1%)	25	36
All	All	659/674 (98%)	623 (94%)	33 (5%)	3 (0%)	29	41

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	203	THR
1	B	242	LEU
1	B	145	VAL

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%)	230 (81%)	54 (19%)	1 1
1	B	280/287 (98%)	235 (84%)	45 (16%)	2 3
All	All	564/574 (98%)	465 (82%)	99 (18%)	2 2

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	ASN
1	A	0	THR
1	A	1	THR
1	A	2	SER
1	A	8	ASN
1	A	36	SER
1	A	54	LYS
1	A	69	THR
1	A	71	LEU
1	A	81	SER
1	A	88	ILE
1	A	91	VAL
1	A	100	MET
1	A	110	LEU
1	A	132	ARG
1	A	133	VAL
1	A	136	ILE
1	A	141	ILE
1	A	143	GLN
1	A	149	ASP
1	A	152	SER
1	A	156	ASN
1	A	157	ARG
1	A	158	ASP
1	A	159	SER
1	A	160(D)	SER
1	A	161	LEU
1	A	173	GLN
1	A	186	LYS
1	A	195	LYS
1	A	201	SER
1	A	202	SER
1	A	204	LEU

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Mol	Chain	Res	Type
1	A	205	LEU
1	A	207	GLU
1	A	219	SER
1	A	229	GLU
1	A	233	GLU
1	A	240	ARG
1	A	247	VAL
1	A	248	LYS
1	A	255	LEU
1	A	266	GLU
1	A	270	THR
1	A	278	GLU
1	A	279	SER
1	A	279(C)	SER
1	A	279(D)	LYS
1	A	287	HIS
1	A	298	THR
1	A	316	ARG
1	A	319	ARG
1	A	324	LEU
1	A	326	ARG
1	B	8	ASN
1	B	20	ILE
1	B	36	SER
1	B	46	SER
1	B	46(A)	ARG
1	B	59	SER
1	B	62	SER
1	B	65	LYS
1	B	70	GLU
1	B	71	LEU
1	B	86	GLN
1	B	94	ILE
1	B	95	THR
1	B	96	VAL
1	B	107	MET
1	B	120	VAL
1	B	139	ASN
1	B	147	LYS
1	B	156	ASN
1	B	157	ARG
1	B	160(D)	SER

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Mol	Chain	Res	Type
1	B	173	GLN
1	B	178	ASN
1	B	186	LYS
1	B	187	THR
1	B	195	LYS
1	B	204	LEU
1	B	207	GLU
1	B	226	SER
1	B	233	GLU
1	B	235	LEU
1	B	239	LYS
1	B	244	ASP
1	B	247	VAL
1	B	254	THR
1	B	265	LYS
1	B	279	SER
1	B	279(C)	SER
1	B	279(D)	LYS
1	B	287	HIS
1	B	291	ILE
1	B	308	LYS
1	B	315	ARG
1	B	319	ARG
1	B	326	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-1	ASN
1	A	8	ASN
1	A	25	GLN
1	A	86	GLN
1	A	99	GLN
1	A	139	ASN
1	A	156	ASN
1	A	174	HIS
1	A	277	GLN
1	A	317	ASN
1	A	318	ASN
1	B	8	ASN
1	B	25	GLN
1	B	156	ASN

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Mol	Chain	Res	Type
1	B	174	HIS
1	B	180	HIS
1	B	183	ASN
1	B	261	HIS
1	B	277	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

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	$\#\|Z\| > 2$
2	0IU	B	391	-	43,47,47	1.70	10 (23%)	48,64,64	2.03	13 (27%)
2	0IU	A	391	-	43,47,47	1.76	8 (18%)	48,64,64	1.93	10 (20%)

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
2	0IU	B	391	-	-	4/51/60/60	0/3/3/3
2	0IU	A	391	-	-	5/51/60/60	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	391	0IU	CD2-CG	-5.24	1.30	1.39
2	B	391	0IU	CD21-SE2	5.14	1.78	1.70
2	A	391	0IU	CD21-SE2	4.21	1.77	1.70
2	A	391	0IU	CD1-CG	-4.01	1.32	1.39
2	B	391	0IU	CB-N1	-3.48	1.43	1.48
2	B	391	0IU	CD1-CG	-3.39	1.33	1.39
2	B	391	0IU	CD2-CG	-3.14	1.34	1.39
2	A	391	0IU	C31-C21	2.98	1.57	1.52
2	B	391	0IU	CA-N1	2.54	1.48	1.45
2	B	391	0IU	C4-N2	2.48	1.39	1.34
2	A	391	0IU	CE2-CD2	-2.39	1.33	1.38
2	B	391	0IU	C-N	-2.36	1.31	1.34
2	B	391	0IU	C31-C21	2.30	1.56	1.52
2	A	391	0IU	CZ-CE1	-2.28	1.32	1.38
2	A	391	0IU	CE1-CD1	-2.18	1.34	1.38
2	B	391	0IU	CG2-CB	2.16	1.57	1.52
2	B	391	0IU	CE11-NZ	-2.14	1.28	1.35
2	A	391	0IU	CZ-CE2	-2.02	1.32	1.38

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	391	0IU	CG2-CB-CG	-7.42	94.52	113.24
2	A	391	0IU	CN-CA1-C4	-5.70	101.52	109.77
2	A	391	0IU	CB1-CG1-CD21	-5.60	120.01	130.18
2	A	391	0IU	CG2-CB-N1	-5.18	104.98	110.38
2	B	391	0IU	CA-C-N	5.16	120.49	116.36
2	B	391	0IU	CB1-CG1-CD21	-4.10	122.74	130.18
2	A	391	0IU	C32-C22-C12	-3.80	109.54	114.82
2	B	391	0IU	CN-CA1-C4	-3.56	104.61	109.77
2	A	391	0IU	O2-C4-CA1	-3.41	117.69	122.12
2	A	391	0IU	CA-C-N	3.34	119.03	116.36
2	B	391	0IU	C4'-C3'-C2'	-3.20	104.88	111.42
2	B	391	0IU	C-CA-N1	-3.05	108.45	112.06
2	B	391	0IU	O-C-N	-2.80	117.30	121.83
2	B	391	0IU	C32-C22-C12	-2.75	111.00	114.82
2	B	391	0IU	CG2-CB-N1	-2.40	107.89	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	391	0IU	C4'-C5'-C6'	-2.29	106.74	111.42
2	A	391	0IU	C4'-C5'-C6'	-2.28	106.77	111.42
2	A	391	0IU	O1-C3-N1	-2.27	116.62	121.54
2	A	391	0IU	C31-C1'-C6'	-2.21	106.94	111.73
2	B	391	0IU	C2-N-C1	2.15	122.44	115.77
2	B	391	0IU	O2-C4-CA1	-2.14	119.33	122.12
2	B	391	0IU	CG-CB-N1	2.09	113.08	110.40
2	A	391	0IU	CA-N1-CB	2.00	120.70	117.90

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	391	0IU	C4-CA1-CB1-CG1
2	A	391	0IU	O11-C11-C21-C31
2	A	391	0IU	C4-CA1-CB1-CG1
2	B	391	0IU	CG2-CB-N1-CA
2	B	391	0IU	CG2-CB-N1-C3
2	A	391	0IU	O11-C11-C21-N2
2	A	391	0IU	C12-C11-C21-N2
2	B	391	0IU	N1-CB-CG-CD2
2	A	391	0IU	C12-C11-C21-C31

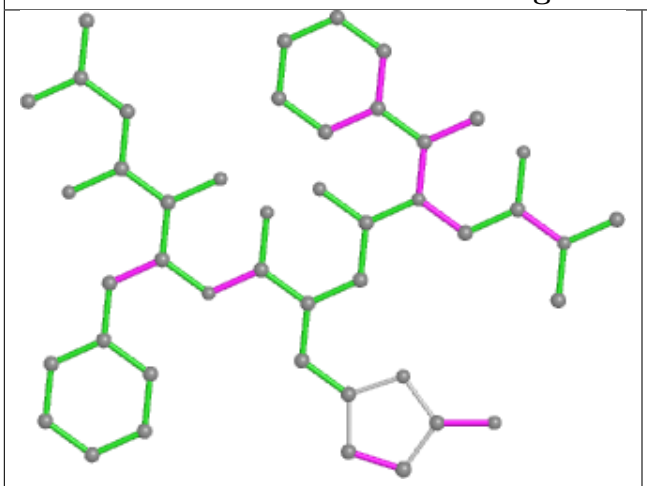
There are no ring outliers.

2 monomers are involved in 25 short contacts:

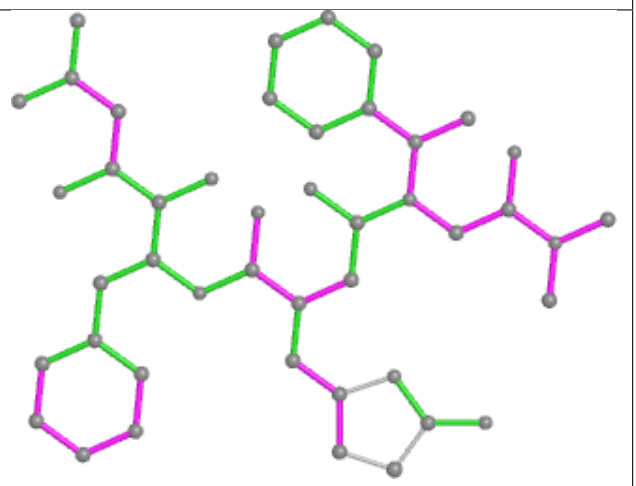
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	391	0IU	15	0
2	A	391	0IU	10	0

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 then 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.

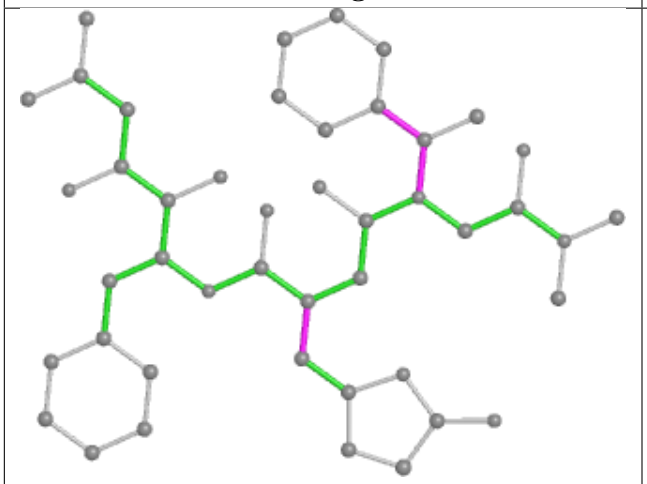
Ligand 0IU B 391



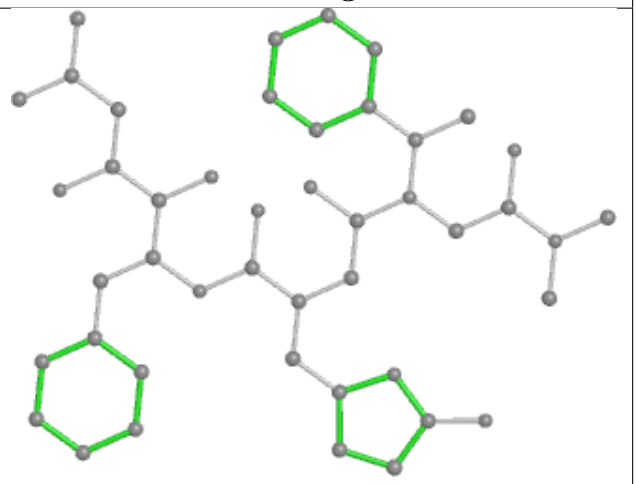
Bond lengths



Bond angles

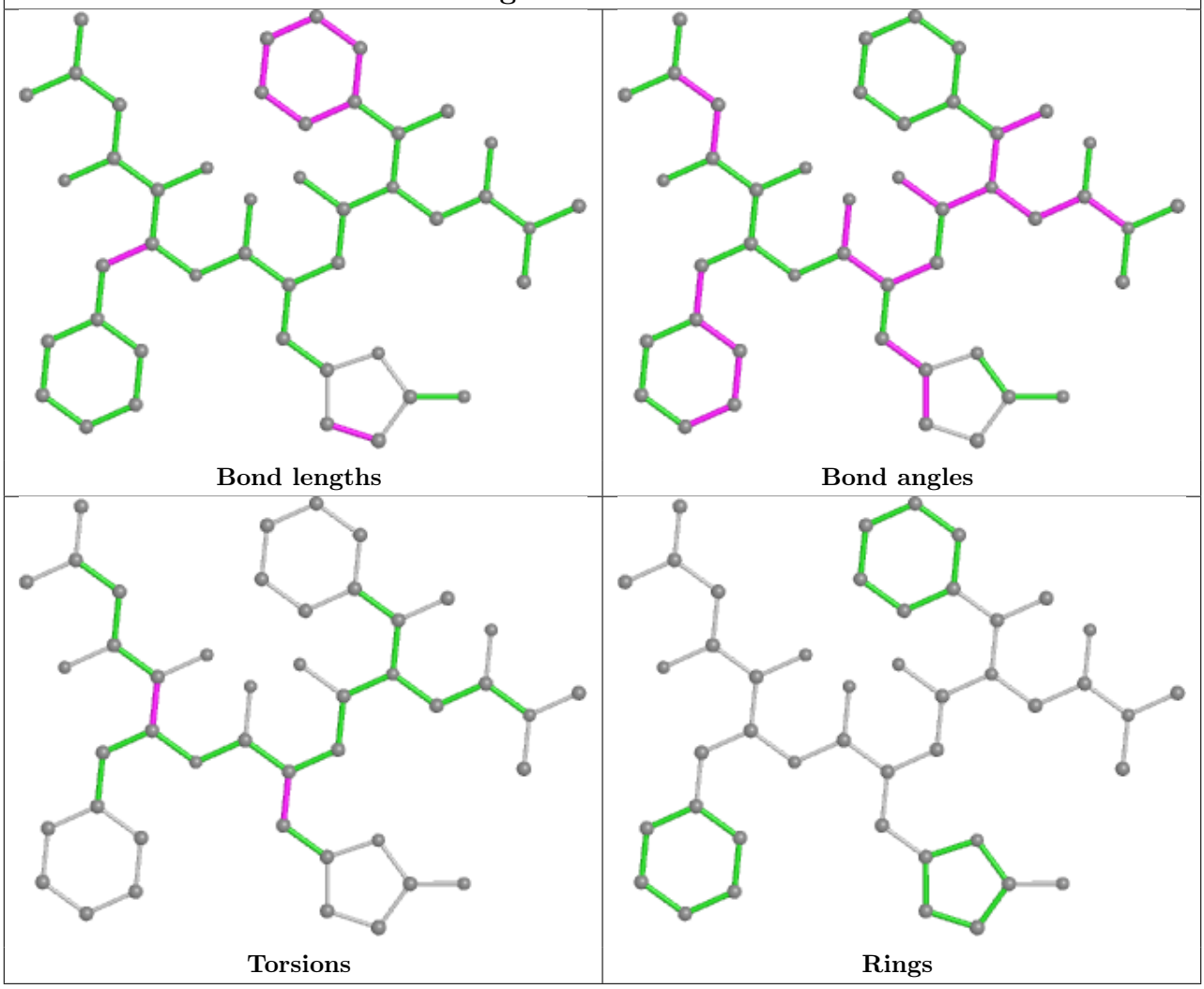


Torsions



Rings

Ligand 0IU A 391



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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.