



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:54 AM UTC

PDB ID : 1NA1 / pdb_00001na1
Title : The structure of HRV14 when complexed with Pleconaril
Authors : Zhang, Y.; Simpson, A.A.; Bator, C.M.; Chakravarty, S.; Pevear, D.C.;
Skochko, G.A.; Tull, T.M.; Diana, G.; Rossmann, M.G.
Deposited on : 2002-11-26
Resolution : 3.30 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

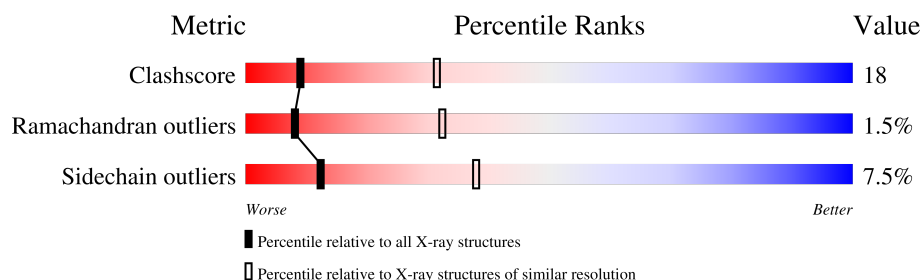
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)

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	289	
2	B	262	
3	C	236	
4	D	68	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	W11	A	7001	-	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6537 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 Coat protein VP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2170	1373	375	414	8			

- Molecule 2 is a protein called Coat protein VP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	255	Total	C	N	O	S	0	0	0
			1952	1238	330	372	12			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	170	LEU	ILE	SEE REMARK 999	UNP P03303

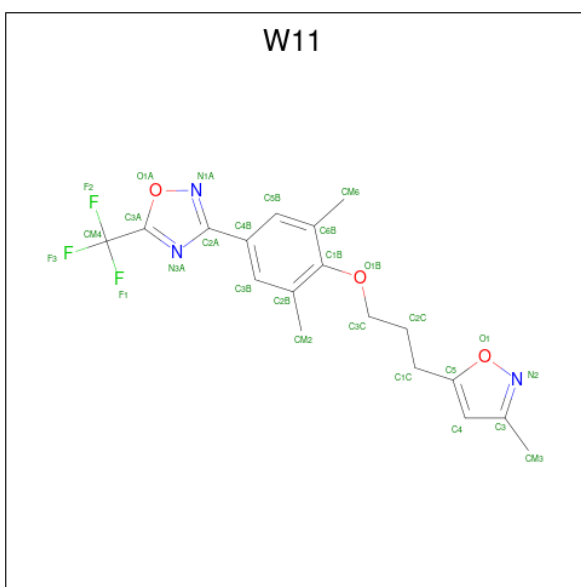
- Molecule 3 is a protein called Coat protein VP3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	236	Total	C	N	O	S	0	0	0
			1849	1184	305	353	7			

- Molecule 4 is a protein called Coat protein VP4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	40	Total	C	N	O	S	0	0	0
			297	186	47	62	2			

- Molecule 5 is 3-{3,5-DIMETHYL-4-[3-(3-METHYL-ISOXAZOL-5-YL)-PROPOXY]-PHENYL}-5-TRIFLUOROMETHYL-[1,2,4]OXADIAZOLE (CCD ID: W11) (formula: C₁₈H₁₈F₃N₃O₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	F	N	O	0	0
			27	18	3	3	3		

- Molecule 6 is water.

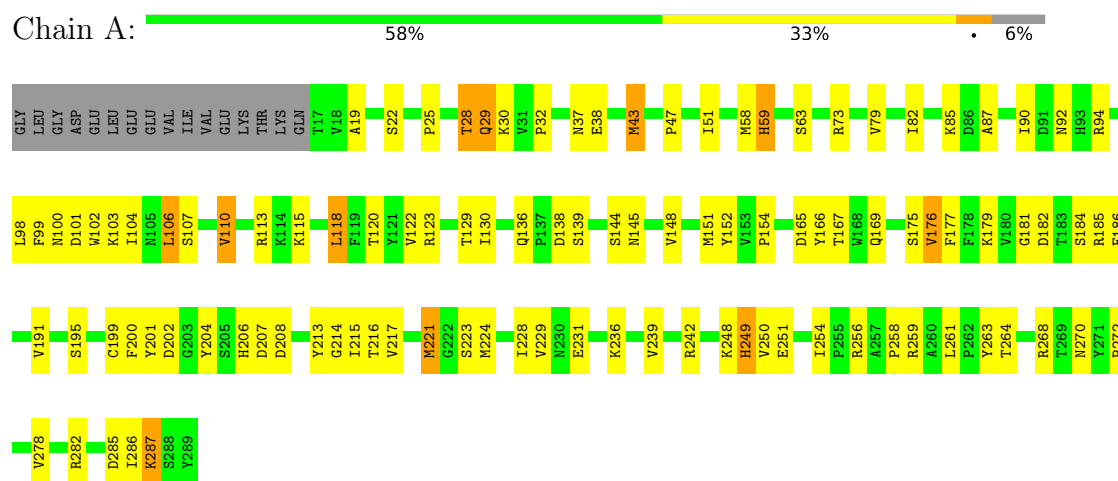
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	87	Total O 87 87	0	0
6	B	69	Total O 69 69	0	0
6	C	80	Total O 80 80	0	0
6	D	6	Total O 6 6	0	0

3 Residue-property plots

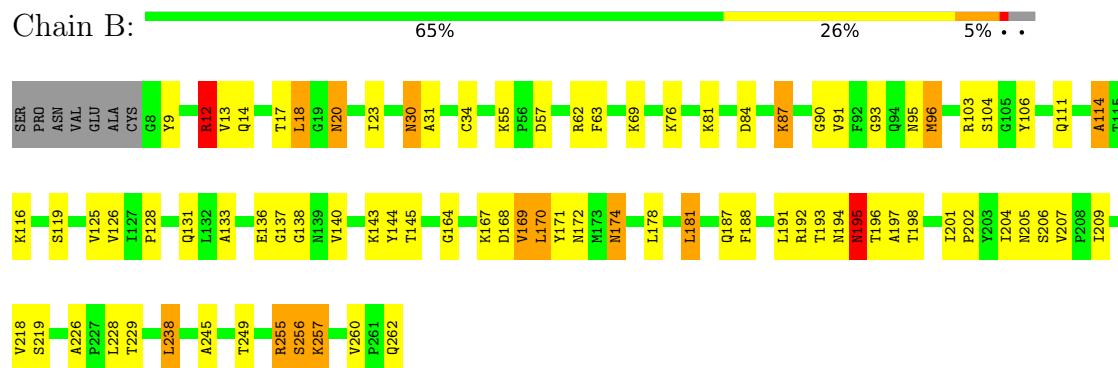
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

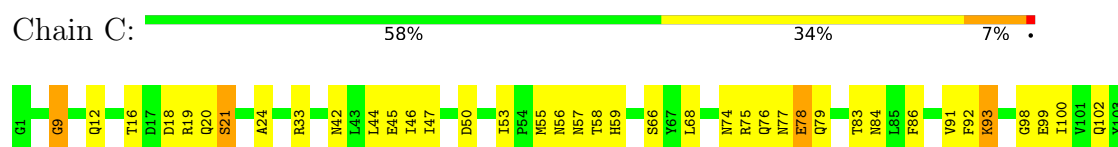
• Molecule 1: Coat protein VP1

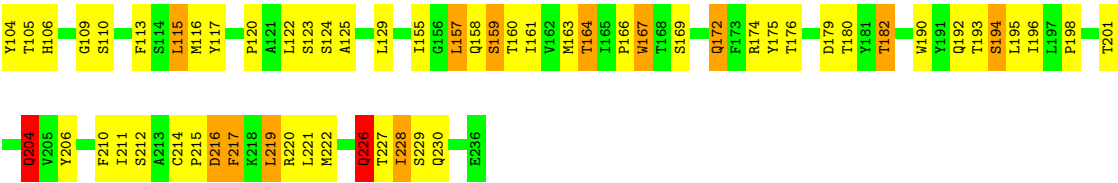


• Molecule 2: Coat protein VP2



• Molecule 3: Coat protein VP3





• Molecule 4: Coat protein VP4



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	437.16 Å 437.16 Å 437.16 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.30	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-3.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.222 , 0.228	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6537	wwPDB-VP
Average B, all atoms (Å ²)	38.0	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: W11

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	0.57	0/2228	1.11	15/3031 (0.5%)
2	B	0.62	1/2001 (0.0%)	1.05	12/2735 (0.4%)
3	C	6.96	1/1898 (0.1%)	1.64	14/2597 (0.5%)
4	D	0.68	1/302 (0.3%)	1.16	2/406 (0.5%)
All	All	3.81	3/6429 (0.0%)	1.28	43/8769 (0.5%)

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	C	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	226	GLN	C-O	302.05	5.25	1.24
2	B	193	THR	C-N	-5.20	1.25	1.33
4	D	47	MET	SD-CE	5.13	1.92	1.79

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	226	GLN	O-C-N	-55.80	50.41	122.39
3	C	226	GLN	CA-C-O	32.34	158.56	119.43
1	A	286	ILE	N-CA-C	-9.74	101.60	112.80
1	A	110	VAL	N-CA-C	9.38	122.77	111.05
1	A	29	GLN	N-CA-C	-9.05	102.25	113.38
1	A	166	TYR	N-CA-C	-8.61	101.63	112.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	217	PHE	N-CA-C	8.15	121.76	110.35
1	A	118	LEU	N-CA-C	-7.72	103.65	113.23
1	A	250	VAL	N-CA-C	7.66	119.62	109.21
3	C	180	THR	N-CA-C	7.52	119.12	111.07
3	C	159	SER	N-CA-C	7.37	119.32	111.28
1	A	98	LEU	N-CA-C	-7.06	103.63	111.82
4	D	41	ALA	N-CA-C	6.28	118.93	111.71
1	A	199	CYS	N-CA-C	-6.23	105.51	113.23
1	A	169	GLN	N-CA-C	-6.22	104.56	111.71
2	B	226	ALA	N-CA-C	-6.15	100.69	109.24
2	B	207	VAL	CA-C-N	6.06	125.54	119.24
2	B	207	VAL	C-N-CA	6.06	125.54	119.24
2	B	256	SER	N-CA-C	6.03	123.65	110.80
3	C	46	ILE	N-CA-C	6.03	118.63	111.09
1	A	229	VAL	N-CA-C	6.01	117.47	110.62
2	B	201	ILE	N-CA-C	5.95	114.51	107.73
3	C	74	ASN	N-CA-C	5.86	119.26	111.30
2	B	194	ASN	N-CA-C	5.84	118.99	110.17
2	B	57	ASP	CB-CA-C	-5.75	109.97	116.63
4	D	46	SER	N-CA-C	-5.67	102.39	111.02
1	A	231	GLU	N-CA-C	-5.64	103.19	110.19
2	B	55	LYS	CA-C-N	5.61	125.28	119.56
2	B	55	LYS	C-N-CA	5.61	125.28	119.56
3	C	9	GLY	N-CA-C	-5.54	107.89	114.48
2	B	164	GLY	CA-C-N	5.50	126.65	120.66
2	B	164	GLY	C-N-CA	5.50	126.65	120.66
2	B	12	ARG	N-CA-C	5.50	117.35	111.36
1	A	181	GLY	N-CA-C	-5.48	108.17	114.69
3	C	160	THR	N-CA-C	5.39	118.60	109.76
3	C	92	PHE	N-CA-C	5.31	117.81	111.71
1	A	28	THR	N-CA-C	5.28	117.30	109.69
1	A	191	VAL	CB-CA-C	-5.20	106.21	112.19
3	C	93	LYS	N-CA-C	5.11	118.29	111.75
3	C	215	PRO	CA-C-N	5.05	127.01	120.44
3	C	215	PRO	C-N-CA	5.05	127.01	120.44
3	C	204	GLN	N-CA-C	5.04	117.61	109.40
1	A	59	HIS	N-CA-C	-5.00	105.84	112.24

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	226	GLN	Mainchain

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	2170	0	2107	93	0
2	B	1952	0	1927	80	0
3	C	1849	0	1833	92	0
4	D	297	0	294	6	0
5	A	27	0	18	4	0
6	A	87	0	0	2	0
6	B	69	0	0	1	0
6	C	80	0	0	3	0
6	D	6	0	0	0	0
All	All	6537	0	6179	229	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 18.

All (229) 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)
5:A:7001:W11:O1A	5:A:7001:W11:N1A	1.57	1.18
1:A:47:PRO:HA	3:C:164:THR:HG21	1.46	0.94
3:C:55:MET:HE3	3:C:91:VAL:HG11	1.51	0.91
1:A:282:ARG:HG3	3:C:57:ASN:HB2	1.53	0.89
3:C:42:ASN:HD22	3:C:44:LEU:H	1.17	0.86
1:A:38:GLU:OE1	3:C:116:MET:HE2	1.76	0.85
1:A:228:ILE:HD11	1:A:239:VAL:HG21	1.60	0.82
2:B:205:ASN:HD22	2:B:206:SER:H	1.28	0.81
2:B:255:ARG:HG2	2:B:256:SER:N	1.96	0.81
2:B:187:GLN:NE2	2:B:198:THR:H	1.82	0.78
1:A:151:MET:HA	1:A:175:SER:HB3	1.67	0.76
2:B:195:ASN:C	2:B:195:ASN:HD22	1.93	0.75
3:C:226:GLN:O	3:C:227:THR:HB	1.87	0.74
2:B:12:ARG:NH2	3:C:157:LEU:HD21	2.01	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:9:GLY:O	3:C:12:GLN:HG2	1.87	0.74
2:B:256:SER:O	2:B:257:LYS:HB3	1.86	0.74
2:B:195:ASN:ND2	2:B:196:THR:HG23	2.03	0.74
1:A:282:ARG:HE	3:C:57:ASN:HD22	1.34	0.74
2:B:205:ASN:ND2	2:B:206:SER:H	1.86	0.73
2:B:84:ASP:OD1	2:B:87:LYS:HE3	1.89	0.73
1:A:201:TYR:H	2:B:131:GLN:HE21	1.36	0.72
1:A:285:ASP:HB3	1:A:287:LYS:H	1.53	0.72
2:B:12:ARG:NH1	2:B:12:ARG:HG3	2.04	0.71
2:B:30:ASN:HD21	4:D:58:ASP:H	1.37	0.71
3:C:164:THR:HG22	3:C:166:PRO:HD3	1.74	0.70
2:B:128:PRO:HA	2:B:218:VAL:HG12	1.72	0.70
2:B:12:ARG:HG3	2:B:12:ARG:HH11	1.57	0.69
1:A:100:ASN:HD22	1:A:101:ASP:H	1.39	0.68
3:C:79:GLN:HB2	3:C:190:TRP:CZ3	2.28	0.68
1:A:201:TYR:H	2:B:131:GLN:NE2	1.91	0.68
2:B:114:ALA:HB2	2:B:238:LEU:HD21	1.76	0.68
1:A:58:MET:HE1	3:C:216:ASP:O	1.92	0.68
1:A:100:ASN:ND2	1:A:101:ASP:H	1.92	0.68
3:C:182:THR:HG23	3:C:182:THR:O	1.95	0.67
1:A:175:SER:HA	5:A:7001:W11:F3	1.82	0.67
2:B:170:LEU:HD13	2:B:170:LEU:O	1.94	0.67
3:C:125:ALA:HB3	3:C:155:ILE:HD12	1.76	0.67
2:B:12:ARG:HH21	3:C:157:LEU:HD21	1.57	0.67
3:C:53:ILE:HD11	3:C:211:ILE:HB	1.77	0.67
2:B:12:ARG:HH21	3:C:157:LEU:CD2	2.09	0.66
2:B:195:ASN:C	2:B:195:ASN:ND2	2.53	0.65
1:A:261:LEU:HD11	2:B:171:TYR:CD1	2.31	0.65
3:C:50:ASP:HB3	3:C:212:SER:HB3	1.79	0.65
1:A:152:TYR:CE1	1:A:154:PRO:HG3	2.32	0.65
1:A:258:PRO:CB	3:C:99:GLU:HG2	2.27	0.65
2:B:81:LYS:HG3	2:B:219:SER:HB3	1.78	0.65
1:A:202:ASP:OD2	2:B:81:LYS:HE3	1.97	0.64
3:C:98:GLY:O	3:C:102:GLN:HG3	1.97	0.64
3:C:20:GLN:HE22	4:D:30:ASN:HA	1.63	0.64
1:A:82:ILE:HD12	1:A:228:ILE:HG13	1.80	0.63
2:B:63:PHE:CD2	2:B:245:ALA:HB2	2.33	0.63
3:C:42:ASN:O	3:C:45:GLU:HB2	1.99	0.63
2:B:228:LEU:HD12	2:B:229:THR:H	1.63	0.63
1:A:270:ASN:HA	2:B:133:ALA:HB1	1.80	0.63
2:B:205:ASN:HD22	2:B:206:SER:N	1.96	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:20:ASN:HD21	2:B:62:ARG:HE	1.46	0.62
1:A:37:ASN:HB2	3:C:116:MET:HE3	1.82	0.61
1:A:28:THR:HB	1:A:30:LYS:H	1.65	0.61
3:C:18:ASP:OD1	4:D:40:SER:HB2	2.00	0.61
1:A:176:VAL:HG22	5:A:7001:W11:F2	1.91	0.61
1:A:47:PRO:CA	3:C:164:THR:HG21	2.27	0.61
1:A:104:ILE:HG13	1:A:223:SER:HA	1.83	0.60
2:B:195:ASN:HD22	2:B:196:THR:N	1.98	0.60
2:B:136:GLU:HB3	2:B:140:VAL:HG21	1.83	0.60
1:A:282:ARG:HE	3:C:57:ASN:ND2	1.99	0.59
2:B:255:ARG:CG	2:B:256:SER:N	2.60	0.59
3:C:42:ASN:ND2	3:C:44:LEU:H	1.95	0.58
1:A:38:GLU:O	2:B:187:GLN:HG3	2.02	0.58
2:B:256:SER:O	2:B:257:LYS:CB	2.51	0.58
3:C:42:ASN:HD22	3:C:44:LEU:N	1.97	0.58
3:C:68:LEU:HB3	3:C:206:TYR:CD2	2.39	0.57
1:A:217:VAL:HG11	6:A:7004:HOH:O	2.04	0.57
1:A:120:THR:HB	1:A:254:ILE:HB	1.86	0.57
3:C:57:ASN:O	3:C:59:HIS:N	2.38	0.57
2:B:187:GLN:HE21	2:B:197:ALA:HA	1.70	0.56
1:A:186:PHE:HD2	3:C:24:ALA:HB2	1.69	0.56
3:C:204:GLN:N	6:C:241:HOH:O	2.38	0.56
2:B:12:ARG:NH2	3:C:157:LEU:CD2	2.69	0.56
2:B:205:ASN:ND2	2:B:206:SER:N	2.53	0.55
4:D:56:VAL:HG11	4:D:60:MET:HE2	1.88	0.55
1:A:107:SER:HB2	1:A:113:ARG:NH1	2.21	0.55
1:A:248:LYS:HE3	4:D:36:ALA:O	2.08	0.54
3:C:75:ARG:HD2	3:C:78:GLU:OE2	2.08	0.54
3:C:77:ASN:HB3	6:C:275:HOH:O	2.07	0.54
3:C:50:ASP:HB3	3:C:212:SER:CB	2.36	0.54
3:C:84:ASN:HD22	3:C:86:PHE:H	1.55	0.54
1:A:115:LYS:O	1:A:118:LEU:HB2	2.07	0.54
1:A:129:THR:OG1	1:A:185:ARG:NH1	2.36	0.54
3:C:115:LEU:HD22	3:C:129:LEU:HD21	1.89	0.53
2:B:116:LYS:HB3	3:C:122:LEU:HG	1.90	0.53
1:A:100:ASN:HD22	1:A:101:ASP:N	2.06	0.53
1:A:259:ARG:HD2	1:A:263:TYR:CZ	2.43	0.53
1:A:282:ARG:HG3	3:C:57:ASN:CB	2.32	0.52
2:B:104:SER:HB2	2:B:249:THR:HG23	1.91	0.52
2:B:20:ASN:C	2:B:20:ASN:HD22	2.18	0.52
1:A:264:THR:CG2	1:A:272:PRO:HD3	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:193:THR:O	3:C:194:SER:HB3	2.09	0.52
1:A:228:ILE:HD11	1:A:239:VAL:CG2	2.36	0.52
1:A:282:ARG:NH2	3:C:55:MET:HE2	2.25	0.52
1:A:251:GLU:HA	6:C:311:HOH:O	2.10	0.51
2:B:170:LEU:HD12	2:B:171:TYR:CE2	2.46	0.51
2:B:17:THR:O	2:B:18:LEU:HD13	2.11	0.51
3:C:182:THR:O	3:C:182:THR:CG2	2.60	0.50
2:B:12:ARG:HH11	2:B:12:ARG:CG	2.22	0.50
3:C:169:SER:HB2	3:C:174:ARG:NH1	2.26	0.50
1:A:206:HIS:NE2	1:A:208:ASP:HB2	2.27	0.50
1:A:259:ARG:CD	1:A:263:TYR:CZ	2.94	0.50
2:B:34:CYS:HB2	2:B:202:PRO:HD2	1.94	0.50
1:A:102:TRP:O	1:A:223:SER:HB2	2.12	0.50
1:A:214:GLY:C	1:A:216:THR:H	2.19	0.50
2:B:30:ASN:HD22	2:B:31:ALA:N	2.10	0.50
1:A:51:ILE:HD13	3:C:166:PRO:HG3	1.94	0.49
1:A:100:ASN:ND2	1:A:101:ASP:N	2.60	0.49
3:C:214:CYS:C	3:C:216:ASP:H	2.20	0.49
3:C:76:GLN:O	3:C:78:GLU:N	2.37	0.49
3:C:55:MET:CE	3:C:91:VAL:HG11	2.35	0.48
1:A:107:SER:HB2	1:A:113:ARG:HD2	1.95	0.48
1:A:185:ARG:HG3	1:A:186:PHE:N	2.29	0.48
1:A:258:PRO:HB2	3:C:99:GLU:HG2	1.94	0.48
3:C:120:PRO:HG2	3:C:123:SER:HB3	1.96	0.48
3:C:110:SER:H	3:C:216:ASP:HB3	1.78	0.48
2:B:172:ASN:HB2	2:B:174:ASN:ND2	2.27	0.48
2:B:13:VAL:O	2:B:14:GLN:HG2	2.13	0.48
3:C:124:SER:O	3:C:125:ALA:HB2	2.13	0.48
3:C:57:ASN:HD21	3:C:91:VAL:CG1	2.26	0.48
1:A:113:ARG:NE	1:A:263:TYR:CE2	2.81	0.48
1:A:110:VAL:HG22	3:C:230:GLN:OE1	2.14	0.48
2:B:103:ARG:HB3	2:B:209:ILE:HG13	1.94	0.48
1:A:228:ILE:CD1	1:A:239:VAL:HG21	2.37	0.48
1:A:259:ARG:HD3	1:A:263:TYR:CE1	2.49	0.48
3:C:93:LYS:HG2	3:C:102:GLN:HE22	1.79	0.48
1:A:186:PHE:CD2	3:C:24:ALA:HB2	2.47	0.47
3:C:228:ILE:HD13	3:C:229:SER:H	1.78	0.47
1:A:107:SER:CB	1:A:113:ARG:HD2	2.44	0.47
2:B:23:ILE:HD12	2:B:63:PHE:CZ	2.49	0.47
3:C:117:TYR:CD2	3:C:155:ILE:HD13	2.49	0.47
1:A:103:LYS:HA	1:A:223:SER:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:33:ARG:HG3	3:C:33:ARG:HH11	1.78	0.47
3:C:129:LEU:N	3:C:129:LEU:HD12	2.30	0.47
1:A:207:ASP:HB3	2:B:260:VAL:HG12	1.95	0.47
3:C:84:ASN:ND2	3:C:86:PHE:H	2.13	0.47
1:A:73:ARG:HD2	3:C:222:MET:HE2	1.97	0.47
2:B:187:GLN:HE21	2:B:198:THR:H	1.62	0.47
3:C:42:ASN:ND2	3:C:44:LEU:HB2	2.30	0.47
2:B:174:ASN:N	2:B:174:ASN:HD22	2.13	0.46
1:A:204:TYR:CZ	2:B:144:TYR:HA	2.50	0.46
1:A:221:MET:HE2	1:A:221:MET:HA	1.98	0.46
3:C:50:ASP:CA	3:C:212:SER:HB3	2.45	0.46
1:A:19:ALA:CB	1:A:58:MET:HE2	2.45	0.46
1:A:79:VAL:HG22	1:A:242:ARG:HG2	1.96	0.46
1:A:215:ILE:HG23	2:B:131:GLN:HE22	1.81	0.46
2:B:20:ASN:H	2:B:20:ASN:ND2	2.14	0.46
2:B:255:ARG:CG	2:B:256:SER:H	2.28	0.46
3:C:161:ILE:HG23	3:C:161:ILE:O	2.16	0.46
1:A:87:ALA:O	1:A:90:ILE:HG12	2.16	0.46
2:B:30:ASN:HD22	2:B:31:ALA:H	1.63	0.46
1:A:85:LYS:HB3	1:A:236:LYS:HG3	1.98	0.46
1:A:282:ARG:HD2	1:A:285:ASP:O	2.15	0.45
3:C:47:ILE:HG22	3:C:100:ILE:HD12	1.98	0.45
2:B:20:ASN:HD22	2:B:20:ASN:N	2.14	0.45
2:B:202:PRO:O	2:B:204:ILE:HG13	2.16	0.45
3:C:214:CYS:C	3:C:216:ASP:N	2.75	0.45
3:C:106:HIS:HB2	3:C:220:ARG:HG3	1.99	0.45
2:B:172:ASN:OD1	2:B:178:LEU:HA	2.16	0.45
2:B:172:ASN:HB2	2:B:174:ASN:HD21	1.81	0.45
1:A:248:LYS:O	1:A:249:HIS:HB2	2.16	0.44
1:A:152:TYR:O	1:A:154:PRO:HD3	2.16	0.44
3:C:104:TYR:O	3:C:176:THR:HG21	2.17	0.44
2:B:20:ASN:C	2:B:20:ASN:ND2	2.75	0.44
1:A:130:ILE:HB	1:A:184:SER:HB3	1.99	0.44
1:A:151:MET:CA	1:A:175:SER:HB3	2.45	0.44
2:B:9:TYR:N	2:B:9:TYR:CD1	2.84	0.44
2:B:30:ASN:ND2	2:B:31:ALA:H	2.16	0.44
1:A:30:LYS:O	1:A:32:PRO:HD3	2.17	0.44
1:A:200:PHE:CE1	1:A:256:ARG:HD2	2.52	0.44
3:C:109:GLY:O	3:C:167:TRP:HB2	2.17	0.44
6:A:7011:HOH:O	2:B:168:ASP:HB2	2.18	0.44
1:A:261:LEU:HD11	2:B:171:TYR:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:163:MET:HG2	3:C:164:THR:N	2.32	0.43
3:C:179:ASP:OD1	3:C:182:THR:HB	2.18	0.43
3:C:190:TRP:CD1	3:C:190:TRP:N	2.85	0.43
2:B:174:ASN:ND2	2:B:174:ASN:C	2.75	0.43
3:C:175:TYR:H	3:C:182:THR:HG21	1.84	0.43
1:A:165:ASP:HB3	1:A:167:THR:H	1.84	0.43
1:A:204:TYR:CE2	1:A:213:TYR:HB2	2.53	0.43
3:C:50:ASP:CB	3:C:212:SER:HB3	2.45	0.43
1:A:102:TRP:HZ3	1:A:224:MET:HE3	1.83	0.43
1:A:214:GLY:C	1:A:216:THR:N	2.76	0.43
1:A:82:ILE:CD1	1:A:228:ILE:HG13	2.48	0.43
1:A:148:VAL:O	1:A:177:PHE:HA	2.19	0.43
2:B:174:ASN:HD22	2:B:174:ASN:H	1.66	0.43
2:B:76:LYS:HB3	2:B:169:VAL:CG2	2.49	0.42
2:B:18:LEU:HD12	2:B:18:LEU:HA	1.89	0.42
3:C:192:GLN:HA	3:C:192:GLN:NE2	2.34	0.42
2:B:91:VAL:HG12	2:B:95:ASN:ND2	2.35	0.42
2:B:137:GLY:O	2:B:138:GLY:C	2.63	0.42
1:A:258:PRO:HG2	3:C:99:GLU:HG2	2.01	0.42
2:B:93:GLY:O	2:B:96:MET:HB2	2.20	0.42
3:C:21:SER:O	4:D:37:SER:CB	2.68	0.42
3:C:18:ASP:C	3:C:19:ARG:HD2	2.44	0.42
3:C:105:THR:HB	3:C:221:LEU:HB3	2.02	0.42
3:C:172:GLN:HE21	3:C:172:GLN:HB3	1.71	0.42
1:A:38:GLU:OE1	3:C:159:SER:OG	2.31	0.42
2:B:228:LEU:HD12	2:B:229:THR:N	2.32	0.41
1:A:38:GLU:HA	2:B:188:PHE:HB2	2.02	0.41
1:A:92:ASN:HD21	1:A:94:ARG:HB3	1.84	0.41
1:A:99:PHE:CD2	1:A:99:PHE:C	2.97	0.41
3:C:113:PHE:CE1	3:C:115:LEU:HD13	2.54	0.41
3:C:219:LEU:HD12	3:C:219:LEU:HA	1.87	0.41
1:A:106:LEU:HD22	5:A:7001:W11:C3	2.51	0.41
2:B:125:VAL:HG22	2:B:126:VAL:N	2.35	0.41
1:A:179:LYS:O	1:A:182:ASP:HB2	2.19	0.41
2:B:81:LYS:CG	2:B:219:SER:HB3	2.48	0.41
2:B:174:ASN:HD22	2:B:174:ASN:C	2.29	0.41
2:B:125:VAL:CG2	2:B:126:VAL:N	2.84	0.41
3:C:55:MET:HE1	3:C:83:THR:HG21	2.02	0.41
1:A:282:ARG:HG3	3:C:57:ASN:ND2	2.36	0.41
1:A:37:ASN:CB	3:C:116:MET:HE3	2.49	0.41
1:A:38:GLU:CD	3:C:116:MET:HE2	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:167:LYS:HE3	6:B:307:HOH:O	2.21	0.41
3:C:198:PRO:HG2	3:C:201:THR:OG1	2.21	0.41
3:C:210:PHE:N	3:C:210:PHE:CD1	2.88	0.41
3:C:195:LEU:HD12	3:C:196:ILE:N	2.37	0.41
1:A:118:LEU:HA	1:A:118:LEU:HD12	1.83	0.40
3:C:109:GLY:C	3:C:167:TRP:HB2	2.46	0.40
1:A:123:ARG:HD2	1:A:195:SER:O	2.20	0.40
2:B:172:ASN:CG	2:B:181:LEU:HD13	2.47	0.40
1:A:43:MET:HA	1:A:43:MET:HE3	2.03	0.40
1:A:58:MET:O	1:A:59:HIS:HB2	2.21	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	271/289 (94%)	252 (93%)	16 (6%)	3 (1%)	11	39
2	B	253/262 (97%)	225 (89%)	23 (9%)	5 (2%)	6	27
3	C	234/236 (99%)	219 (94%)	11 (5%)	4 (2%)	7	30
4	D	38/68 (56%)	33 (87%)	5 (13%)	0	100	100
All	All	796/855 (93%)	729 (92%)	55 (7%)	12 (2%)	8	32

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	139	SER
3	C	58	THR
2	B	257	LYS
3	C	78	GLU
2	B	191	LEU

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Mol	Chain	Res	Type
2	B	195	ASN
1	A	249	HIS
2	B	90	GLY
3	C	194	SER
3	C	167	TRP
2	B	114	ALA
1	A	25	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	239/253 (94%)	224 (94%)	15 (6%)	16	44
2	B	223/229 (97%)	202 (91%)	21 (9%)	8	30
3	C	209/209 (100%)	194 (93%)	15 (7%)	13	40
4	D	33/57 (58%)	31 (94%)	2 (6%)	17	45
All	All	704/748 (94%)	651 (92%)	53 (8%)	12	38

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	SER
1	A	29	GLN
1	A	43	MET
1	A	63	SER
1	A	106	LEU
1	A	122	VAL
1	A	136	GLN
1	A	138	ASP
1	A	144	SER
1	A	145	ASN
1	A	176	VAL
1	A	221	MET
1	A	268	ARG
1	A	278	VAL

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Mol	Chain	Res	Type
1	A	287	LYS
2	B	12	ARG
2	B	18	LEU
2	B	20	ASN
2	B	30	ASN
2	B	69	LYS
2	B	87	LYS
2	B	96	MET
2	B	106	TYR
2	B	111	GLN
2	B	119	SER
2	B	143	LYS
2	B	145	THR
2	B	169	VAL
2	B	170	LEU
2	B	174	ASN
2	B	181	LEU
2	B	192	ARG
2	B	195	ASN
2	B	238	LEU
2	B	255	ARG
2	B	262	GLN
3	C	16	THR
3	C	21	SER
3	C	56	ASN
3	C	66	SER
3	C	115	LEU
3	C	157	LEU
3	C	158	GLN
3	C	164	THR
3	C	172	GLN
3	C	182	THR
3	C	204	GLN
3	C	216	ASP
3	C	217	PHE
3	C	219	LEU
3	C	228	ILE
4	D	47	MET
4	D	61	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	HIS
1	A	61	ASN
1	A	84	ASN
1	A	92	ASN
1	A	100	ASN
1	A	136	GLN
1	A	198	ASN
1	A	212	GLN
1	A	219	ASN
1	A	235	HIS
2	B	15	GLN
2	B	20	ASN
2	B	30	ASN
2	B	94	GLN
2	B	111	GLN
2	B	131	GLN
2	B	135	HIS
2	B	174	ASN
2	B	187	GLN
2	B	190	ASN
2	B	194	ASN
2	B	195	ASN
2	B	205	ASN
2	B	216	ASN
2	B	217	ASN
3	C	20	GLN
3	C	42	ASN
3	C	56	ASN
3	C	57	ASN
3	C	72	ASN
3	C	74	ASN
3	C	77	ASN
3	C	84	ASN
3	C	102	GLN
3	C	140	GLN
3	C	172	GLN
3	C	192	GLN
3	C	204	GLN
3	C	226	GLN
4	D	30	ASN
4	D	43	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	W11	A	7001	-	29,29,29	4.14	16 (55%)	41,42,42	3.72	26 (63%)

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
5	W11	A	7001	-	-	2/17/17/17	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	7001	W11	C3A-N3A	12.10	1.41	1.28
5	A	7001	W11	O1A-N1A	8.45	1.57	1.42
5	A	7001	W11	C2A-N1A	6.78	1.41	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	7001	W11	C5B-C6B	6.15	1.48	1.39
5	A	7001	W11	C1B-C6B	5.23	1.49	1.40
5	A	7001	W11	C5B-C4B	5.10	1.47	1.39
5	A	7001	W11	C1B-C2B	4.89	1.48	1.40
5	A	7001	W11	O1A-C3A	4.58	1.39	1.33
5	A	7001	W11	O1-N2	4.34	1.50	1.42
5	A	7001	W11	O1-C5	4.09	1.41	1.35
5	A	7001	W11	C3B-C4B	3.48	1.44	1.39
5	A	7001	W11	C2A-N3A	2.98	1.43	1.37
5	A	7001	W11	C4-C5	2.75	1.38	1.34
5	A	7001	W11	CM2-C2B	-2.16	1.47	1.51
5	A	7001	W11	CM4-C3A	-2.11	1.47	1.51
5	A	7001	W11	C4B-C2A	-2.07	1.44	1.47

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	7001	W11	C3A-O1A-N1A	-12.29	100.23	105.34
5	A	7001	W11	F1-CM4-C3A	-6.67	98.64	111.67
5	A	7001	W11	O1A-C3A-N3A	6.00	119.09	115.25
5	A	7001	W11	CM4-C3A-N3A	-5.80	121.63	127.22
5	A	7001	W11	C4B-C2A-N1A	5.57	126.98	122.22
5	A	7001	W11	CM3-C3-N2	-5.11	115.78	119.97
5	A	7001	W11	C3B-C2B-C1B	4.94	125.09	117.94
5	A	7001	W11	C3B-C4B-C2A	-4.91	113.68	120.27
5	A	7001	W11	C5B-C4B-C3B	4.18	124.54	119.65
5	A	7001	W11	C4B-C3B-C2B	-4.11	116.68	121.74
5	A	7001	W11	C4B-C5B-C6B	-4.03	116.78	121.74
5	A	7001	W11	O1A-N1A-C2A	3.96	105.60	103.28
5	A	7001	W11	O1B-C1B-C6B	3.60	125.56	118.90
5	A	7001	W11	CM6-C6B-C5B	-3.42	113.55	119.57
5	A	7001	W11	F2-CM4-C3A	3.37	118.27	111.67
5	A	7001	W11	CM6-C6B-C1B	3.31	126.11	120.80
5	A	7001	W11	F3-CM4-C3A	-3.22	105.38	111.67
5	A	7001	W11	CM3-C3-C4	2.88	132.93	128.62
5	A	7001	W11	C5-O1-N2	-2.63	106.32	108.30
5	A	7001	W11	C6B-C1B-C2B	-2.62	116.57	122.07
5	A	7001	W11	C2C-C1C-C5	-2.60	108.17	113.49
5	A	7001	W11	O1A-C3A-CM4	2.48	119.27	116.75
5	A	7001	W11	O1B-C3C-C2C	-2.27	99.95	108.30
5	A	7001	W11	CM2-C2B-C3B	-2.23	115.65	119.57
5	A	7001	W11	F3-CM4-F1	2.19	115.66	106.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	7001	W11	C1C-C2C-C3C	-2.08	108.15	112.33

There are no chirality outliers.

All (2) torsion outliers are listed below:

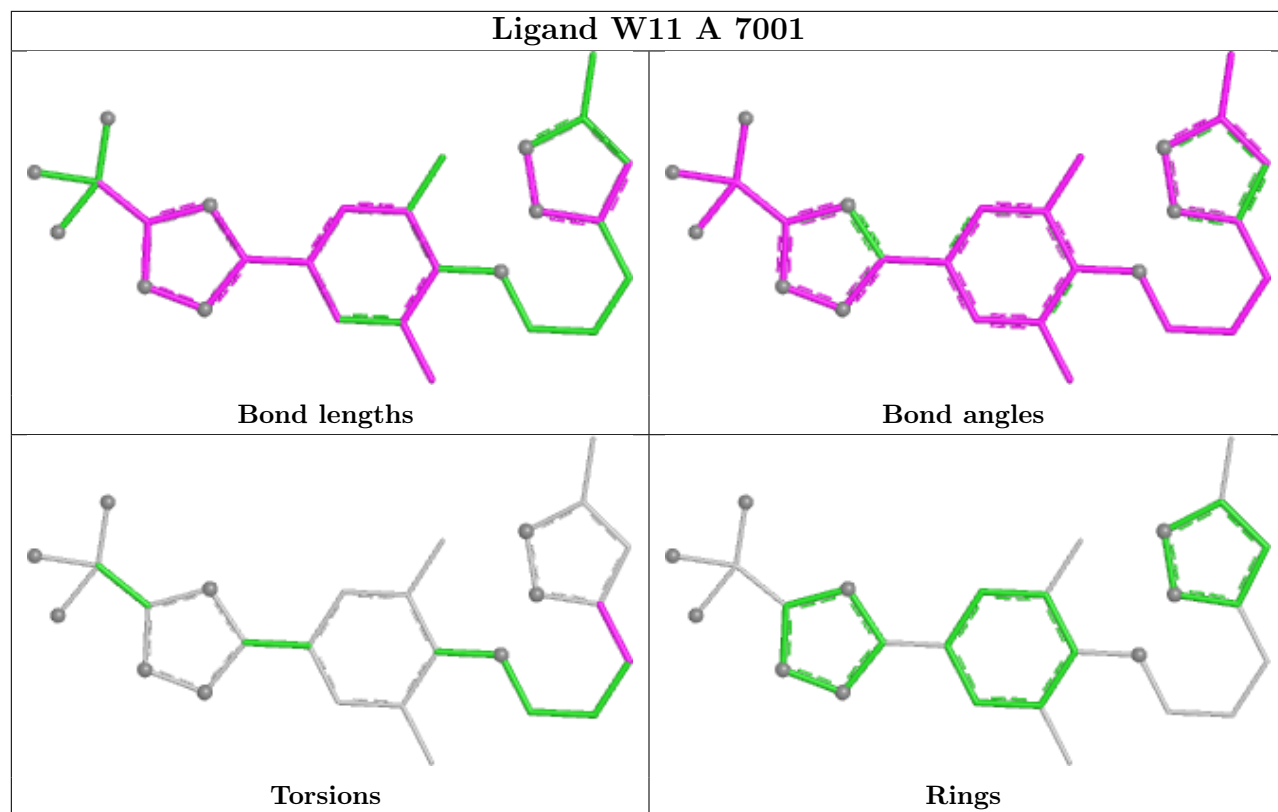
Mol	Chain	Res	Type	Atoms
5	A	7001	W11	C2C-C1C-C5-O1
5	A	7001	W11	C2C-C1C-C5-C4

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	7001	W11	4	0

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