



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 05:41 PM UTC

PDB ID : 5HVP / pdb_00005hvp
Title : CRYSTALLOGRAPHIC ANALYSIS OF A COMPLEX BETWEEN HUMAN IMMUNODEFICIENCY VIRUS TYPE 1 PROTEASE AND ACETYL-PEPSTATIN AT 2.0-ANGSTROMS RESOLUTION
Authors : Fitzgerald, P.M.D.; McKeever, B.M.; Vanmiddlesworth, J.F.; Springer, J.P.
Deposited on : 1990-04-30
Resolution : 2.00 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
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

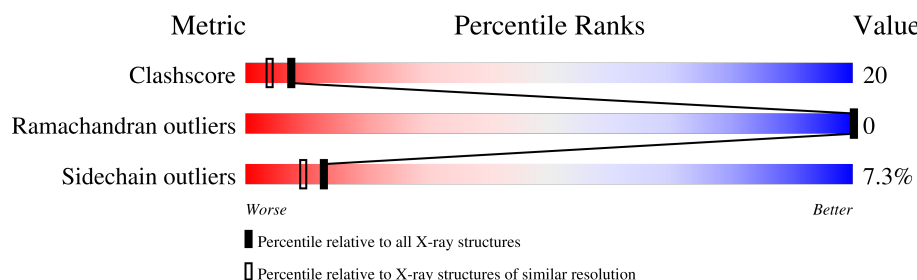
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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	99	49% 37% 13%
1	B	99	55% 36% 9%
2	C	6	33% 33% 17% 17%

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
2	STA	C	4[A]	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 1802 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 HIV-1 PROTEASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	99	Total	C	N	O	S	0	8	0
			759	485	129	140	5			
1	B	99	Total	C	N	O	S	0	9	0
			781	502	134	140	5			

- Molecule 2 is a protein called ACETYL-*PEPSTATIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	6	0
			90	62	10	18			

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	67	Total	O	0	7
			67	67		
4	B	99	Total	O	0	7
			99	99		
4	C	4	Total	O	0	2
			4	4		

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.

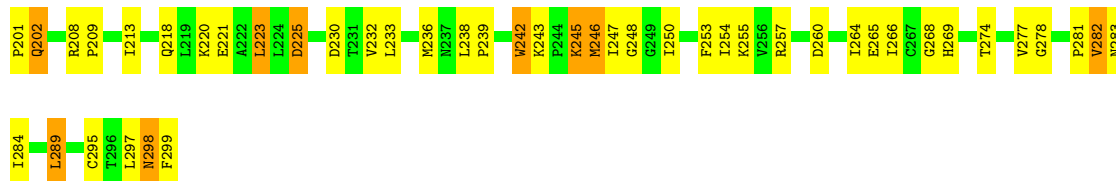
• Molecule 1: HIV-1 PROTEASE

Chain A: 

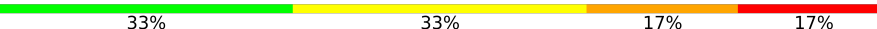


• Molecule 1: HIV-1 PROTEASE

Chain B: 



• Molecule 2: ACETYL-*PEPSTATIN

Chain C: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	58.39Å 86.70Å 46.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.176 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1802	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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: CL, STA, ACE

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	1.36	2/822 (0.2%)	2.27	37/1114 (3.3%)
1	B	1.50	3/844 (0.4%)	2.12	26/1143 (2.3%)
2	C	0.62	0/36	1.08	0/46
All	All	1.42	5/1702 (0.3%)	2.18	63/2303 (2.7%)

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
2	C	0	3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	40	GLY	N-CA	6.56	1.51	1.45
1	B	248	GLY	N-CA	5.46	1.51	1.45
1	A	54	ILE	CA-C	5.04	1.58	1.52
1	B	243[A]	LYS	N-CA	5.03	1.50	1.45
1	B	243[B]	LYS	N-CA	5.03	1.50	1.45

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	58	GLN	OE1-CD-NE2	9.74	132.34	122.60
1	A	8	ARG	CD-NE-CZ	-9.57	111.00	124.40
1	B	269	HIS	CA-CB-CG	-8.68	105.12	113.80
1	B	260	ASP	CA-CB-CG	-8.24	104.36	112.60
1	B	243[A]	LYS	CB-CA-C	8.12	121.56	109.22

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	243[B]	LYS	CB-CA-C	8.12	121.56	109.22
1	A	37	ASN	CA-CB-CG	-7.62	104.98	112.60
1	B	274	THR	CA-CB-OG1	-7.31	98.63	109.60
1	A	40	GLY	N-CA-C	-7.26	103.54	112.33
1	A	85	ILE	N-CA-CB	7.20	119.20	111.00
1	A	85	ILE	O-C-N	7.18	131.70	122.94
1	A	49	GLY	O-C-N	7.06	129.66	123.80
1	A	85	ILE	CA-C-O	-6.88	112.60	120.65
1	A	22	ALA	CA-C-N	6.61	130.95	121.50
1	A	22	ALA	C-N-CA	6.61	130.95	121.50
1	B	278	GLY	O-C-N	6.59	128.36	121.77
1	A	75	VAL	N-CA-CB	-6.55	102.29	111.99
1	A	6	TRP	N-CA-C	-6.54	104.24	111.82
1	B	208	ARG	O-C-N	6.50	126.71	121.38
1	A	55[A]	LYS	CA-C-O	-6.37	113.34	120.54
1	A	55[B]	LYS	CA-C-O	-6.37	113.34	120.54
1	A	37	ASN	OD1-CG-ND2	6.35	128.95	122.60
1	A	75	VAL	CB-CA-C	6.35	118.76	110.12
1	A	7	GLN	OE1-CD-NE2	-6.28	116.32	122.60
1	B	298[A]	ASN	CA-CB-CG	-6.01	106.59	112.60
1	B	298[B]	ASN	CA-CB-CG	-6.01	106.59	112.60
1	A	7	GLN	CB-CG-CD	5.91	122.65	112.60
1	B	230[A]	ASP	N-CA-CB	-5.89	101.93	110.70
1	B	230[B]	ASP	N-CA-CB	-5.89	101.93	110.70
1	B	225	ASP	N-CA-CB	-5.81	101.40	110.87
1	B	245	LYS	CA-CB-CG	-5.75	102.60	114.10
1	A	60[A]	ASP	CB-CA-C	5.72	119.16	109.50
1	A	60[B]	ASP	CB-CA-C	5.72	119.16	109.50
1	A	32	VAL	O-C-N	5.63	129.04	123.18
1	A	25	ASP	N-CA-CB	-5.60	101.17	110.47
1	A	11	VAL	CA-C-O	-5.52	116.23	121.64
1	B	243[A]	LYS	CA-C-N	5.52	125.28	119.76
1	B	243[A]	LYS	C-N-CA	5.52	125.28	119.76
1	B	243[B]	LYS	CA-C-N	5.52	125.28	119.76
1	B	243[B]	LYS	C-N-CA	5.52	125.28	119.76
1	A	83	ASN	CA-C-N	-5.50	115.34	122.99
1	A	83	ASN	C-N-CA	-5.50	115.34	122.99
1	B	238	LEU	CB-CA-C	5.46	117.43	109.08
1	A	38	LEU	O-C-N	5.45	126.78	121.66
1	B	223	LEU	N-CA-CB	-5.41	101.23	110.16
1	B	282	VAL	O-C-N	5.39	128.28	122.67
1	A	75	VAL	CA-C-O	-5.38	114.93	120.25

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	ASN	CA-CB-CG	5.37	117.97	112.60
1	A	88	ASN	OD1-CG-ND2	-5.34	117.26	122.60
1	A	21	GLU	O-C-N	5.34	129.59	123.29
1	B	277	VAL	N-CA-CB	-5.33	103.73	111.52
1	A	38	LEU	N-CA-C	-5.33	102.09	110.14
1	A	3	ILE	N-CA-CB	-5.29	104.64	110.99
1	B	242	TRP	O-C-N	5.27	130.17	123.11
1	B	281	PRO	CA-C-N	5.25	128.00	122.11
1	B	281	PRO	C-N-CA	5.25	128.00	122.11
1	A	64	ILE	O-C-N	5.20	128.66	123.10
1	B	221	GLU	O-C-N	5.15	129.22	123.19
1	B	289	LEU	CB-CA-C	5.13	118.22	109.24
1	A	87	ARG	CA-C-N	5.09	127.36	120.38
1	A	87	ARG	C-N-CA	5.09	127.36	120.38
1	A	82	VAL	CA-C-O	-5.08	116.37	121.45
1	A	39	PRO	N-CD-CG	-5.02	95.67	103.20

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	4[A]	STA	Mainchain
2	C	4[B]	STA	Mainchain,Peptide

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	759	0	767	38	0
1	B	781	0	825	38	0
2	C	90	0	108	14	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	67	0	0	2	0
4	B	99	0	0	2	0
4	C	4	0	0	0	0
All	All	1802	0	1700	69	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 20.

All (69) 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:C:3[A]:VAL:O	2:C:4[A]:STA:HD22	1.77	0.85
1:A:25:ASP:OD1	2:C:4[A]:STA:OH	1.99	0.81
1:A:84:ILE:HD11	2:C:4[B]:STA:HD12	1.63	0.81
1:A:2:GLN:HE21	1:B:298[B]:ASN:HD21	1.29	0.78
1:A:50:ILE:HG23	1:B:254:ILE:HD13	1.67	0.75
1:A:98[A]:ASN:HD21	1:B:202:GLN:NE2	1.87	0.73
1:B:250:ILE:HD13	2:C:3[A]:VAL:HG11	1.72	0.72
1:A:98[A]:ASN:ND2	1:B:202:GLN:HE21	1.90	0.69
1:B:246:MET:SD	1:B:255:LYS:HD2	2.33	0.68
1:B:246:MET:HG2	1:B:255:LYS:HD2	1.76	0.68
1:A:26:THR:HG21	1:A:97[B]:LEU:HD21	1.76	0.67
1:A:65[B]:GLU:OE1	4:A:305[B]:HOH:O	2.12	0.66
1:A:98[B]:ASN:OD1	1:B:202:GLN:HB2	1.95	0.66
1:B:245:LYS:NZ	4:B:319[B]:HOH:O	2.23	0.64
1:A:98[A]:ASN:ND2	1:B:202:GLN:HB2	2.12	0.64
1:A:98[A]:ASN:HD21	1:B:202:GLN:HE21	1.44	0.63
1:A:12[B]:THR:HG21	1:A:19:LEU:HD22	1.80	0.62
1:B:246:MET:CG	1:B:255:LYS:HD2	2.29	0.62
1:A:30:ASP:OD2	2:C:6[B]:STA:OXT	2.19	0.61
1:B:282:VAL:O	1:B:284[A]:ILE:HG23	2.01	0.60
1:B:264[A]:ILE:HG21	1:B:289:LEU:HD13	1.84	0.59
1:A:2:GLN:HE21	1:B:298[B]:ASN:ND2	2.00	0.58
1:A:12[A]:THR:HG23	1:A:21:GLU:OE2	2.05	0.57
1:B:213:ILE:HD12	1:B:264[A]:ILE:HD11	1.87	0.57
1:A:12[B]:THR:CG2	1:A:19:LEU:HD22	2.34	0.57
1:B:265:GLU:OE1	1:B:268:GLY:HA2	2.07	0.55
1:B:257:ARG:HD3	4:B:414:HOH:O	2.07	0.54
2:C:3[A]:VAL:C	2:C:4[A]:STA:HG	2.33	0.54
1:A:2:GLN:HB2	1:B:298[B]:ASN:ND2	2.23	0.54
1:B:247:ILE:O	1:B:253:PHE:HA	2.08	0.54
1:A:25:ASP:OD1	2:C:4[B]:STA:HC	2.07	0.53
1:A:58:GLN:HG3	1:A:74:THR:HG23	1.91	0.52
1:A:69:HIS:N	1:A:69:HIS:ND1	2.58	0.52
1:B:242:TRP:CE2	1:B:257:ARG:HD2	2.45	0.50
2:C:2[A]:VAL:CG1	2:C:4[A]:STA:HD21	2.42	0.50
1:B:233:LEU:HD23	1:B:283:ASN:HB3	1.94	0.50
1:A:15:ILE:HG23	1:A:15:ILE:O	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:ARG:CZ	2:C:6[A]:STA:HD22	2.43	0.49
1:A:99:PHE:OXT	1:B:201:PRO:N	2.46	0.49
1:A:65[B]:GLU:OE1	1:A:70:LYS:HE3	2.13	0.49
1:B:265:GLU:HG2	1:B:268:GLY:CA	2.44	0.48
1:A:1:PRO:N	1:B:299:PHE:OXT	2.46	0.48
1:B:265:GLU:HG2	1:B:268:GLY:HA2	1.95	0.48
1:B:232:VAL:HB	1:B:284[A]:ILE:HG12	1.97	0.47
2:C:2[A]:VAL:CG1	2:C:3[A]:VAL:N	2.78	0.46
1:B:233:LEU:CD2	1:B:283:ASN:HB3	2.46	0.46
1:B:264[A]:ILE:HG21	1:B:289:LEU:CD1	2.46	0.46
1:A:13:ILE:HG21	1:A:85:ILE:HD11	1.97	0.46
1:A:34:GLU:HG2	1:A:80:THR:O	2.17	0.45
1:A:66:ILE:O	1:A:67:CYS:HB2	2.17	0.45
1:B:236:MET:HE3	1:B:236:MET:HB2	1.64	0.44
1:B:246:MET:HE1	1:B:253:PHE:HB3	1.99	0.44
1:A:97[B]:LEU:HD23	1:B:297[B]:LEU:HD13	2.00	0.44
2:C:2[A]:VAL:HG13	2:C:4[A]:STA:HD21	2.01	0.43
1:A:38:LEU:HA	1:A:39:PRO:HD3	1.80	0.42
1:A:87:ARG:HD3	1:B:209:PRO:HD3	2.00	0.42
1:B:264[B]:ILE:HG22	1:B:266:ILE:HG13	2.01	0.42
1:A:8:ARG:HH21	1:A:8:ARG:HD2	1.40	0.41
1:B:223:LEU:CD2	2:C:4[A]:STA:HD13	2.50	0.41
1:A:25:ASP:CG	2:C:4[A]:STA:OH	2.62	0.41
1:A:98[B]:ASN:ND2	1:B:298[B]:ASN:OD1	2.43	0.41
1:B:225:ASP:HB3	1:B:284[B]:ILE:HG23	2.01	0.41
2:C:6[A]:STA:HD13	2:C:6[A]:STA:N	2.36	0.41
1:A:15:ILE:HD11	1:A:38:LEU:HD21	2.02	0.40
1:A:56:VAL:HG12	1:A:78:GLY:HA3	2.04	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	105/99 (106%)	102 (97%)	3 (3%)	0	100	100
1	B	106/99 (107%)	104 (98%)	2 (2%)	0	100	100
2	C	6/6 (100%)	6 (100%)	0	0	100	100
All	All	217/204 (106%)	212 (98%)	5 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	87/83 (105%)	80 (92%)	7 (8%)	11	8
1	B	92/83 (111%)	85 (92%)	7 (8%)	12	9
2	C	4/2 (200%)	4 (100%)	0	100	100
All	All	183/168 (109%)	169 (92%)	14 (8%)	13	8

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12[A]	THR
1	A	12[B]	THR
1	A	15	ILE
1	A	33	LEU
1	A	69	HIS
1	A	75	VAL
1	A	84	ILE
1	B	202	GLN
1	B	218	GLN
1	B	220	LYS
1	B	239	PRO
1	B	246	MET
1	B	295[A]	CYS
1	B	295[B]	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	37	ASN
1	A	92	GLN
1	B	202	GLN
1	B	237	ASN
1	B	288	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	STA	C	4[B]	2	10,10,11	1.38	1 (10%)	9,12,14	1.35	2 (22%)
2	STA	C	6[B]	2	11,11,11	1.39	1 (9%)	11,14,14	0.85	0
2	STA	C	4[A]	2	10,10,11	1.46	1 (10%)	9,12,14	1.16	1 (11%)
2	STA	C	6[A]	2	11,11,11	1.54	3 (27%)	11,14,14	0.69	0

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	STA	C	4[B]	2	-	5/11/11/12	-
2	STA	C	6[B]	2	-	2/12/12/12	-
2	STA	C	4[A]	2	-	7/11/11/12	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	STA	C	6[A]	2	-	2/12/12/12	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	4[A]	STA	CM-CH	3.03	1.58	1.53
2	C	4[B]	STA	CH-CA	2.68	1.56	1.53
2	C	6[B]	STA	CM-CH	2.64	1.56	1.53
2	C	6[A]	STA	CB-CA	2.62	1.57	1.53
2	C	6[A]	STA	CM-CH	2.41	1.56	1.53
2	C	6[A]	STA	CH-CA	2.38	1.55	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4[B]	STA	CM-CH-CA	2.58	116.92	112.93
2	C	4[A]	STA	CM-CH-CA	2.16	116.26	112.93
2	C	4[B]	STA	CH-CM-C	2.15	116.69	113.02

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	4[A]	STA	N-CA-CB-CG
2	C	4[A]	STA	N-CA-CH-OH
2	C	4[A]	STA	N-CA-CH-CM
2	C	4[A]	STA	CB-CA-CH-OH
2	C	4[A]	STA	CB-CA-CH-CM
2	C	4[B]	STA	N-CA-CH-OH
2	C	4[B]	STA	N-CA-CH-CM
2	C	4[B]	STA	CB-CA-CH-OH
2	C	4[B]	STA	CB-CA-CH-CM
2	C	6[A]	STA	CA-CH-CM-C
2	C	6[A]	STA	OH-CH-CM-C
2	C	6[B]	STA	CA-CH-CM-C
2	C	4[B]	STA	O-C-CM-CH
2	C	4[A]	STA	O-C-CM-CH
2	C	4[A]	STA	CA-CB-CG-CD1
2	C	6[B]	STA	OH-CH-CM-C

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	4[B]	STA	2	0
2	C	6[B]	STA	1	0
2	C	4[A]	STA	7	0
2	C	6[A]	STA	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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.