



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:05 PM GMT

PDB ID : 1NCQ  
Title : The structure of HRV14 when complexed with pleconaril, an antiviral compound  
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 : 2002-12-05  
Resolution : 2.50 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

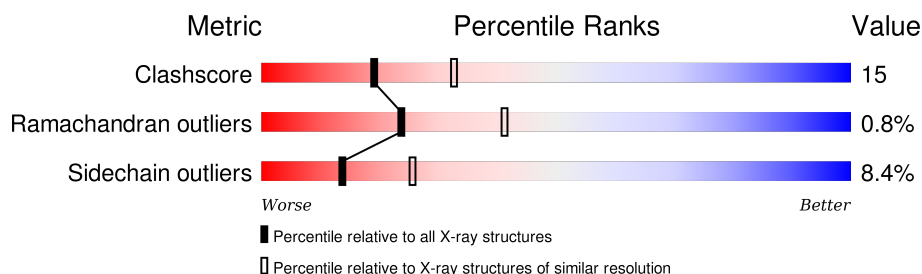
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	289	
2	B	262	
3	C	236	
4	D	68	

## 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

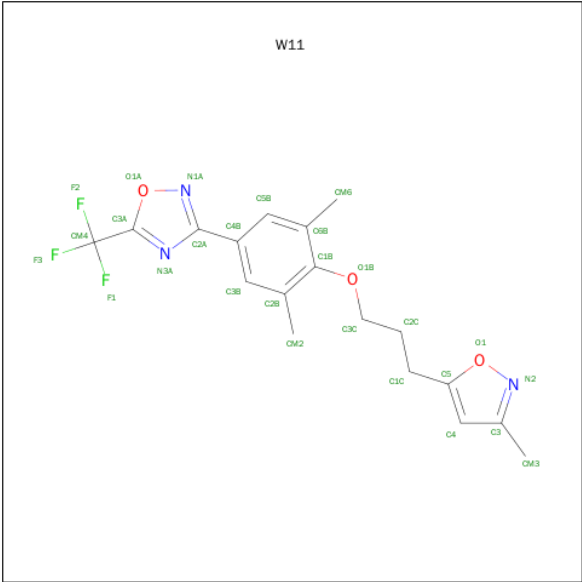
- Molecule 3 is a protein called COAT PROTEIN VP3.

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]-PHE NYL}-5-TRIFLUOROMETHYL-[1,2,4]OXADIAZOLE (three-letter code: W11) (formula: C<sub>18</sub>H<sub>18</sub>F<sub>3</sub>N<sub>3</sub>O<sub>3</sub>).



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.

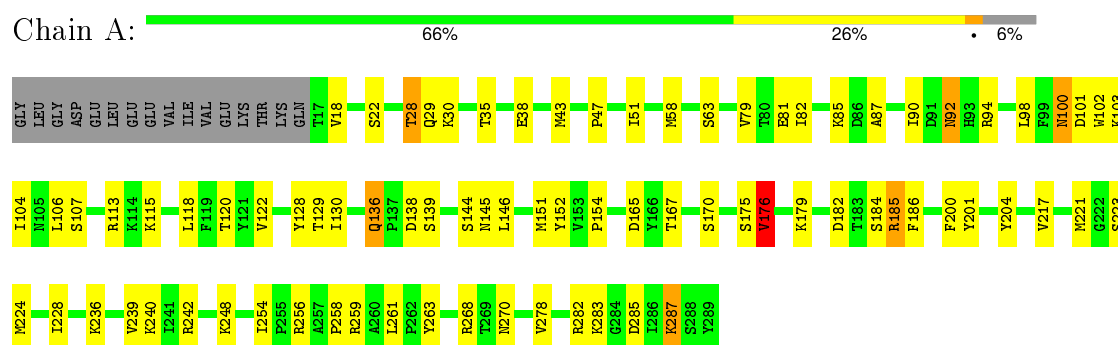
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	83	Total	O	0	0
			83	83		
6	B	69	Total	O	0	0
			69	69		
6	C	83	Total	O	0	0
			83	83		
6	D	7	Total	O	0	0
			7	7		

### 3 Residue-property plots [i](#)

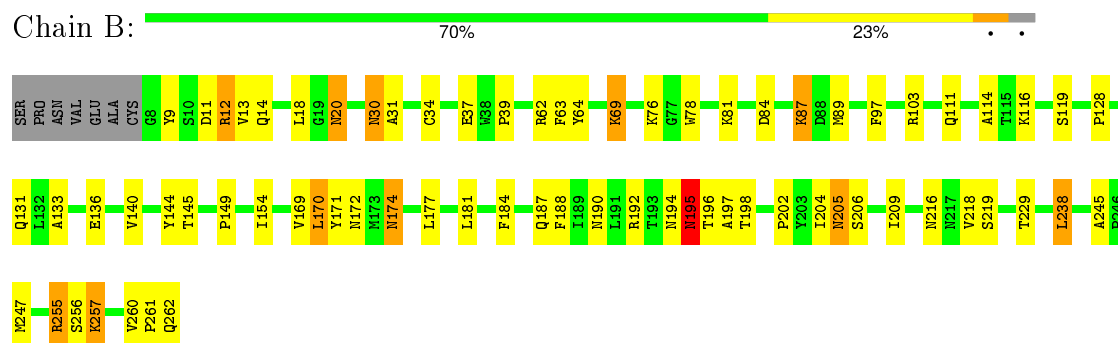
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

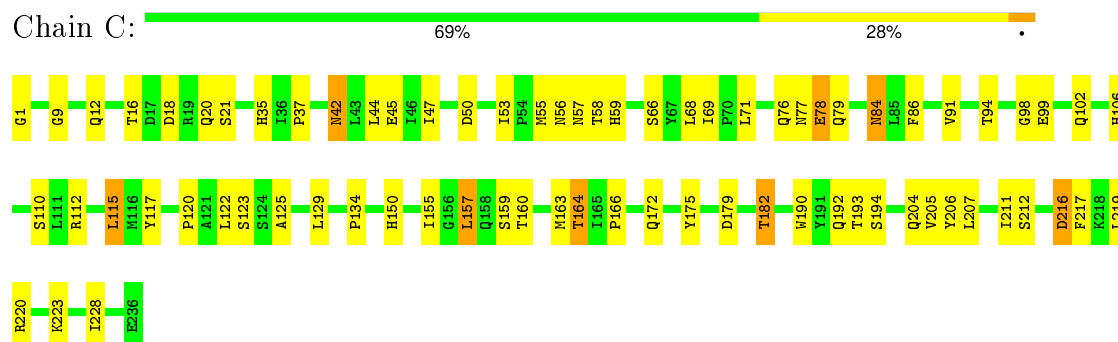
#### • Molecule 1: COAT PROTEIN VP1



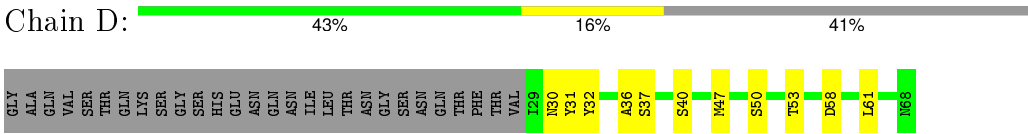
#### • Molecule 2: COAT PROTEIN VP2



#### • Molecule 3: COAT PROTEIN VP3



● Molecule 4: COAT PROTEIN VP4



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	438.72Å 438.72Å 438.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.50)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	unknown	Depositor
R, $R_{free}$	0.216 , 0.217	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6537	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	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	0.41	0/2228	0.70	2/3031 (0.1%)
2	B	0.40	0/2001	0.70	0/2735
3	C	0.39	0/1898	0.72	0/2597
4	D	0.52	0/302	0.71	0/406
All	All	0.41	0/6429	0.71	2/8769 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	176	VAL	CB-CA-C	-6.02	99.97	111.40
1	A	176	VAL	CG1-CB-CG2	5.63	119.91	110.90

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	2170	0	2107	74	0
2	B	1952	0	1927	70	0
3	C	1849	0	1833	78	0
4	D	297	0	294	12	0
5	A	27	0	18	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	83	0	0	1	0
6	B	69	0	0	1	0
6	C	83	0	0	3	0
6	D	7	0	0	0	0
All	All	6537	0	6179	190	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 15.

All (190) 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:N1A	5:A:7001:W11:O1A	1.59	1.18
1:A:47:PRO:HA	3:C:164:THR:HG21	1.45	0.97
3:C:42:ASN:HD22	3:C:44:LEU:H	1.13	0.94
1:A:282:ARG:HG3	3:C:57:ASN:HB2	1.50	0.92
2:B:255:ARG:HG2	2:B:256:SER:N	1.88	0.88
2:B:12:ARG:HH11	2:B:12:ARG:HG3	1.40	0.87
1:A:282:ARG:HE	3:C:57:ASN:HD22	1.26	0.83
2:B:256:SER:O	2:B:257:LYS:HB3	1.76	0.82
1:A:285:ASP:HB3	1:A:287:LYS:H	1.45	0.80
2:B:12:ARG:NH1	2:B:12:ARG:HG3	1.95	0.78
2:B:187:GLN:NE2	2:B:198:THR:H	1.82	0.77
2:B:205:ASN:HD22	2:B:206:SER:H	1.34	0.74
3:C:21:SER:O	4:D:37:SER:HB2	1.88	0.74
2:B:195:ASN:ND2	2:B:196:THR:HG23	2.03	0.74
1:A:176:VAL:HG22	5:A:7001:W11:F2	1.79	0.73
3:C:57:ASN:HD21	3:C:91:VAL:HG12	1.54	0.73
1:A:28:THR:HB	1:A:30:LYS:H	1.55	0.72
2:B:195:ASN:HD22	2:B:195:ASN:C	1.93	0.72
1:A:100:ASN:HD22	1:A:101:ASP:H	1.36	0.71
1:A:152:TYR:CE1	1:A:154:PRO:HG3	2.26	0.69
1:A:201:TYR:H	2:B:131:GLN:HE21	1.38	0.69
3:C:84:ASN:HD22	3:C:86:PHE:H	1.41	0.69
3:C:79:GLN:HB2	3:C:190:TRP:CZ3	2.27	0.69
2:B:255:ARG:CG	2:B:256:SER:N	2.57	0.68
1:A:258:PRO:CB	3:C:99:GLU:HG2	2.24	0.68
3:C:55:MET:HE3	3:C:91:VAL:HG11	1.76	0.67
2:B:30:ASN:HD21	4:D:58:ASP:H	1.43	0.67
3:C:50:ASP:HB3	3:C:212:SER:HB3	1.77	0.67
3:C:192:GLN:HA	3:C:192:GLN:HE21	1.58	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:170:LEU:HD13	2:B:170:LEU:O	1.95	0.67
3:C:57:ASN:HD21	3:C:91:VAL:CG1	2.07	0.66
3:C:20:GLN:HE22	4:D:31:TYR:H	1.42	0.66
2:B:255:ARG:CG	2:B:256:SER:H	2.10	0.65
1:A:100:ASN:ND2	1:A:101:ASP:H	1.94	0.65
2:B:12:ARG:NH2	3:C:157:LEU:CD2	2.60	0.65
1:A:38:GLU:O	2:B:187:GLN:HG3	1.97	0.65
1:A:175:SER:HA	5:A:7001:W11:F3	1.88	0.64
3:C:125:ALA:HB3	3:C:155:ILE:HD12	1.80	0.64
3:C:42:ASN:ND2	3:C:44:LEU:H	1.91	0.64
1:A:47:PRO:CA	3:C:164:THR:HG21	2.24	0.63
3:C:9:GLY:O	3:C:12:GLN:HG2	1.98	0.63
2:B:205:ASN:ND2	2:B:206:SER:H	1.96	0.63
2:B:12:ARG:NH2	3:C:157:LEU:HD21	2.14	0.63
1:A:151:MET:HA	1:A:175:SER:HB3	1.81	0.62
2:B:30:ASN:HD22	2:B:31:ALA:H	1.45	0.62
1:A:228:ILE:HD11	1:A:239:VAL:HG21	1.82	0.62
2:B:195:ASN:ND2	2:B:195:ASN:C	2.53	0.61
3:C:53:ILE:HD11	3:C:211:ILE:HB	1.82	0.61
2:B:187:GLN:HE21	2:B:197:ALA:HA	1.65	0.61
1:A:201:TYR:H	2:B:131:GLN:NE2	1.98	0.61
1:A:107:SER:HB2	1:A:113:ARG:HD2	1.82	0.61
2:B:195:ASN:HD22	2:B:196:THR:N	1.98	0.60
2:B:30:ASN:ND2	2:B:31:ALA:H	2.00	0.60
1:A:217:VAL:HG11	6:A:7003:HOH:O	2.01	0.60
2:B:30:ASN:HD22	2:B:31:ALA:N	1.99	0.60
3:C:164:THR:HG22	3:C:166:PRO:HD3	1.84	0.59
3:C:204:GLN:N	6:C:240:HOH:O	2.35	0.59
1:A:282:ARG:HD2	1:A:285:ASP:O	2.02	0.59
1:A:79:VAL:HG22	1:A:242:ARG:HG2	1.85	0.59
2:B:204:ILE:HG12	3:C:37:PRO:HG2	1.85	0.58
2:B:114:ALA:HB2	2:B:238:LEU:HD21	1.84	0.58
1:A:58:MET:HE1	3:C:216:ASP:O	2.02	0.58
2:B:128:PRO:HA	2:B:218:VAL:HG12	1.86	0.57
3:C:150:HIS:HA	3:C:163:MET:CE	2.34	0.57
1:A:270:ASN:HA	2:B:133:ALA:HB1	1.86	0.57
1:A:85:LYS:HB3	1:A:236:LYS:HG3	1.87	0.56
2:B:256:SER:O	2:B:257:LYS:CB	2.47	0.56
1:A:92:ASN:HD21	1:A:94:ARG:HB3	1.69	0.56
1:A:94:ARG:HH11	1:A:94:ARG:HG2	1.70	0.56
3:C:192:GLN:HA	3:C:192:GLN:NE2	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:103:ARG:HB3	2:B:209:ILE:HG13	1.87	0.56
3:C:42:ASN:HD22	3:C:44:LEU:N	1.95	0.56
2:B:20:ASN:HD22	2:B:20:ASN:C	2.09	0.56
1:A:100:ASN:HD22	1:A:101:ASP:N	2.04	0.55
3:C:115:LEU:HD22	3:C:129:LEU:HD21	1.88	0.55
1:A:129:THR:OG1	1:A:185:ARG:NH1	2.39	0.55
2:B:136:GLU:HB3	2:B:140:VAL:HG21	1.89	0.55
2:B:205:ASN:HD22	2:B:206:SER:N	2.03	0.55
1:A:102:TRP:O	1:A:223:SER:HB2	2.06	0.55
3:C:20:GLN:HE22	4:D:30:ASN:HA	1.70	0.55
3:C:55:MET:HE1	3:C:91:VAL:HG21	1.89	0.55
2:B:84:ASP:OD1	2:B:87:LYS:HE3	2.07	0.55
1:A:259:ARG:HD2	1:A:263:TYR:CE2	2.42	0.54
2:B:255:ARG:HG2	2:B:256:SER:H	1.64	0.54
2:B:63:PHE:CD2	2:B:245:ALA:HB2	2.43	0.54
3:C:110:SER:H	3:C:216:ASP:HB3	1.73	0.54
3:C:42:ASN:O	3:C:45:GLU:HB2	2.08	0.53
3:C:55:MET:CE	3:C:91:VAL:HG11	2.37	0.53
1:A:38:GLU:HA	2:B:188:PHE:HB2	1.91	0.53
3:C:106:HIS:HB2	3:C:220:ARG:HG3	1.90	0.53
2:B:34:CYS:HB2	2:B:202:PRO:HD2	1.90	0.53
1:A:94:ARG:NH1	1:A:94:ARG:HG2	2.24	0.53
2:B:20:ASN:HD21	2:B:62:ARG:HH21	1.56	0.52
3:C:57:ASN:O	3:C:59:HIS:N	2.43	0.52
3:C:68:LEU:HB3	3:C:206:TYR:CD2	2.45	0.52
1:A:261:LEU:HD11	2:B:171:TYR:CD1	2.44	0.52
3:C:179:ASP:OD1	3:C:182:THR:HB	2.10	0.52
3:C:182:THR:HG23	3:C:182:THR:O	2.09	0.52
3:C:18:ASP:OD1	4:D:40:SER:HB2	2.10	0.52
2:B:12:ARG:NH2	3:C:157:LEU:HD22	2.25	0.51
1:A:248:LYS:HE3	4:D:36:ALA:O	2.11	0.51
2:B:97:PHE:O	2:B:257:LYS:HA	2.11	0.51
3:C:1:GLY:N	6:C:250:HOH:O	2.34	0.51
2:B:119:SER:OG	2:B:229:THR:HB	2.10	0.50
3:C:76:GLN:O	3:C:78:GLU:N	2.35	0.50
1:A:228:ILE:HD11	1:A:239:VAL:CG2	2.41	0.50
2:B:37:GLU:CD	3:C:35:HIS:HE2	2.15	0.50
2:B:12:ARG:HH11	2:B:12:ARG:CG	2.14	0.50
3:C:98:GLY:O	3:C:102:GLN:HG3	2.11	0.50
1:A:120:THR:HB	1:A:254:ILE:HB	1.94	0.50
1:A:256:ARG:HB3	2:B:184:PHE:CZ	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:76:LYS:HB3	2:B:169:VAL:CG2	2.42	0.49
1:A:224:MET:HE1	5:A:7001:W11:H3B	1.94	0.49
1:A:104:ILE:HG13	1:A:223:SER:HA	1.93	0.49
1:A:51:ILE:HD13	3:C:166:PRO:HG3	1.94	0.49
2:B:81:LYS:HG3	2:B:219:SER:HB3	1.96	0.48
1:A:87:ALA:O	1:A:90:ILE:HG12	2.13	0.48
1:A:165:ASP:HB3	1:A:167:THR:H	1.77	0.48
1:A:92:ASN:C	1:A:92:ASN:HD22	2.16	0.48
1:A:200:PHE:CE1	1:A:256:ARG:HD2	2.49	0.48
1:A:259:ARG:HD2	1:A:263:TYR:CZ	2.49	0.48
3:C:50:ASP:HB3	3:C:212:SER:CB	2.42	0.48
3:C:57:ASN:ND2	3:C:91:VAL:HG12	2.25	0.47
2:B:12:ARG:HH21	3:C:157:LEU:CD2	2.27	0.47
3:C:117:TYR:CD2	3:C:155:ILE:HD13	2.49	0.47
1:A:104:ILE:HD13	5:A:7001:W11:HM22	1.96	0.47
1:A:47:PRO:HB3	3:C:166:PRO:HB3	1.97	0.46
1:A:282:ARG:HE	3:C:57:ASN:ND2	2.05	0.46
3:C:21:SER:O	4:D:37:SER:CB	2.61	0.46
3:C:44:LEU:HD23	3:C:47:ILE:HD11	1.96	0.46
1:A:258:PRO:HB2	3:C:99:GLU:HG2	1.97	0.46
1:A:82:ILE:HD12	1:A:228:ILE:HG13	1.98	0.46
2:B:34:CYS:HB2	2:B:202:PRO:CD	2.45	0.46
2:B:116:LYS:HB3	3:C:122:LEU:HG	1.97	0.46
2:B:204:ILE:HG12	3:C:37:PRO:CG	2.44	0.46
2:B:205:ASN:ND2	2:B:206:SER:N	2.63	0.46
3:C:193:THR:O	3:C:194:SER:HB3	2.16	0.46
1:A:259:ARG:CD	1:A:263:TYR:CZ	2.98	0.45
1:A:146:LEU:HD13	1:A:228:ILE:HD13	1.99	0.45
1:A:130:ILE:HB	1:A:184:SER:HB3	1.99	0.45
2:B:172:ASN:HB2	2:B:174:ASN:ND2	2.31	0.45
1:A:128:TYR:CZ	5:A:7001:W11:HM23	2.52	0.45
1:A:152:TYR:O	1:A:154:PRO:HD3	2.16	0.45
1:A:282:ARG:HG3	3:C:57:ASN:CB	2.36	0.44
1:A:38:GLU:HB2	2:B:194:ASN:HD21	1.82	0.44
3:C:50:ASP:CB	3:C:212:SER:HB3	2.45	0.44
3:C:175:TYR:H	3:C:182:THR:HG21	1.83	0.44
3:C:84:ASN:ND2	3:C:86:PHE:H	2.11	0.44
1:A:115:LYS:O	1:A:118:LEU:HB2	2.18	0.44
2:B:84:ASP:HB2	2:B:216:ASN:HD21	1.83	0.44
1:A:38:GLU:OE1	3:C:159:SER:OG	2.36	0.44
2:B:69:LYS:HG2	2:B:78:TRP:CD1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:GLN:HB2	1:A:136:GLN:HE21	1.66	0.44
3:C:20:GLN:HE22	4:D:31:TYR:N	2.12	0.44
2:B:177:LEU:CD1	3:C:94:THR:HB	2.48	0.44
2:B:9:TYR:CD1	2:B:9:TYR:N	2.84	0.44
1:A:256:ARG:NH2	2:B:128:PRO:O	2.51	0.43
1:A:152:TYR:CD1	1:A:154:PRO:HG3	2.53	0.43
1:A:258:PRO:CG	3:C:99:GLU:HG2	2.47	0.43
2:B:170:LEU:HD22	2:B:170:LEU:HA	1.75	0.43
1:A:282:ARG:NE	3:C:57:ASN:HD22	2.05	0.43
1:A:283:LYS:HG2	1:A:283:LYS:O	2.19	0.43
2:B:64:TYR:CD2	2:B:89:MET:HB3	2.54	0.43
1:A:35:THR:HG23	3:C:160:THR:HB	2.00	0.43
3:C:129:LEU:N	3:C:129:LEU:HD12	2.34	0.42
2:B:39:PRO:HG3	2:B:247:MET:HG3	2.00	0.42
2:B:20:ASN:ND2	2:B:20:ASN:C	2.70	0.42
2:B:169:VAL:HG23	6:B:297:HOH:O	2.20	0.42
1:A:258:PRO:HG2	3:C:99:GLU:HG2	2.00	0.42
3:C:50:ASP:CA	3:C:212:SER:HB3	2.50	0.42
3:C:69:ILE:HB	3:C:207:LEU:HB2	2.00	0.42
2:B:13:VAL:O	2:B:14:GLN:HG2	2.20	0.42
3:C:150:HIS:HA	3:C:163:MET:HE1	2.02	0.42
3:C:120:PRO:HG2	3:C:123:SER:HB3	2.01	0.42
2:B:174:ASN:HD22	2:B:174:ASN:C	2.22	0.41
2:B:149:PRO:HG3	2:B:154:ILE:HG13	2.01	0.41
1:A:186:PHE:N	1:A:186:PHE:CD1	2.88	0.41
4:D:50:SER:HA	4:D:53:THR:OG1	2.21	0.41
1:A:98:LEU:HD23	1:A:98:LEU:HA	1.90	0.41
1:A:103:LYS:HA	1:A:223:SER:HB3	2.03	0.41
4:D:32:TYR:HB2	4:D:37:SER:OG	2.20	0.41
3:C:20:GLN:NE2	4:D:30:ASN:HA	2.36	0.41
3:C:71:LEU:HB3	3:C:205:VAL:HG22	2.02	0.41
3:C:134:PRO:HG2	6:C:245:HOH:O	2.20	0.41
1:A:204:TYR:CZ	2:B:144:TYR:HA	2.56	0.41
1:A:100:ASN:ND2	1:A:101:ASP:N	2.64	0.41
3:C:20:GLN:NE2	4:D:31:TYR:H	2.14	0.41
1:A:81:GLU:HG2	1:A:82:ILE:N	2.36	0.41
1:A:179:LYS:O	1:A:182:ASP:HB2	2.20	0.40
2:B:260:VAL:HB	2:B:261:PRO:HD2	2.03	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%)	258 (95%)	12 (4%)	1 (0%)	39	61
2	B	253/262 (97%)	230 (91%)	21 (8%)	2 (1%)	24	41
3	C	234/236 (99%)	222 (95%)	9 (4%)	3 (1%)	15	26
4	D	38/68 (56%)	35 (92%)	3 (8%)	0	100	100
All	All	796/855 (93%)	745 (94%)	45 (6%)	6 (1%)	24	41

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	58	THR
1	A	139	SER
2	B	257	LYS
2	B	195	ASN
3	C	77	ASN
3	C	78	GLU

### 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%)	217 (91%)	22 (9%)	11	21
2	B	223/229 (97%)	204 (92%)	19 (8%)	13	25
3	C	209/209 (100%)	193 (92%)	16 (8%)	16	30
4	D	33/57 (58%)	31 (94%)	2 (6%)	23	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	704/748 (94%)	645 (92%)	59 (8%)	14	25

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	VAL
1	A	22	SER
1	A	28	THR
1	A	29	GLN
1	A	43	MET
1	A	63	SER
1	A	92	ASN
1	A	100	ASN
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	170	SER
1	A	176	VAL
1	A	185	ARG
1	A	221	MET
1	A	240	LYS
1	A	268	ARG
1	A	278	VAL
1	A	287	LYS
2	B	11	ASP
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	111	GLN
2	B	145	THR
2	B	170	LEU
2	B	174	ASN
2	B	181	LEU
2	B	190	ASN
2	B	192	ARG
2	B	195	ASN

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Mol	Chain	Res	Type
2	B	205	ASN
2	B	238	LEU
2	B	255	ARG
2	B	262	GLN
3	C	16	THR
3	C	42	ASN
3	C	56	ASN
3	C	66	SER
3	C	84	ASN
3	C	112	ARG
3	C	115	LEU
3	C	157	LEU
3	C	164	THR
3	C	172	GLN
3	C	182	THR
3	C	216	ASP
3	C	217	PHE
3	C	219	LEU
3	C	223	LYS
3	C	228	ILE
4	D	47	MET
4	D	61	LEU

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

Mol	Chain	Res	Type
1	A	61	ASN
1	A	92	ASN
1	A	100	ASN
1	A	136	GLN
1	A	198	ASN
2	B	15	GLN
2	B	20	ASN
2	B	30	ASN
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

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Mol	Chain	Res	Type
2	B	216	ASN
3	C	20	GLN
3	C	41	HIS
3	C	42	ASN
3	C	56	ASN
3	C	57	ASN
3	C	84	ASN
3	C	140	GLN
3	C	172	GLN
3	C	192	GLN
3	C	226	GLN
4	D	30	ASN
4	D	43	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	-	22,29,29	2.75	8 (36%)	29,42,42	2.65	12 (41%)

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	-	-	0/13/17/17	0/1/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	7001	W11	CM4-C3A	-4.29	1.44	1.50
5	A	7001	W11	CM6-C6B	2.15	1.55	1.51
5	A	7001	W11	C2A-N3A	2.75	1.39	1.35
5	A	7001	W11	C3B-C4B	2.79	1.44	1.39
5	A	7001	W11	C5B-C4B	4.00	1.47	1.39
5	A	7001	W11	C5B-C6B	5.18	1.47	1.39
5	A	7001	W11	C1B-C2B	5.68	1.50	1.40
5	A	7001	W11	C1B-C6B	6.13	1.51	1.40

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	7001	W11	C4B-C2A-N3A	-4.80	118.12	123.83
5	A	7001	W11	C4B-C3B-C2B	-4.05	116.74	122.52
5	A	7001	W11	C4B-C5B-C6B	-3.25	117.88	122.52
5	A	7001	W11	C5B-C4B-C2A	-2.88	116.32	120.17
5	A	7001	W11	CM6-C6B-C5B	-2.85	114.18	119.49
5	A	7001	W11	C1C-C2C-C3C	-2.24	105.37	113.29
5	A	7001	W11	C6B-C1B-C2B	-2.16	117.49	122.12
5	A	7001	W11	CM6-C6B-C1B	3.51	126.36	120.89
5	A	7001	W11	O1B-C1B-C6B	3.59	124.50	118.97
5	A	7001	W11	C3B-C2B-C1B	3.79	123.42	117.78
5	A	7001	W11	C5B-C4B-C3B	4.37	125.01	118.29
5	A	7001	W11	C4B-C2A-N1A	6.74	127.61	119.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	7001	W11	6	0

## 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 will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.