



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 12:56 AM GMT

PDB ID : 1FPN
Title : HUMAN RHINOVIRUS SEROTYPE 2 (HRV2)
Authors : Verdaguer, N.; Blaas, D.; Fita, I.
Deposited on : 2000-08-31
Resolution : 2.60 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

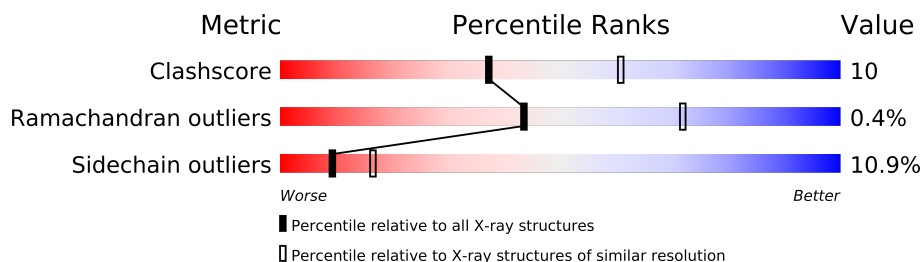
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.60 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	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	1	289	
2	2	261	
3	3	237	
4	4	68	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6316 atoms, of which 0 are hydrogen and 0 are deuterium.

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	1	269	Total	C	N	O	S	0	0	0
			2154	1356	378	409	11			

- Molecule 2 is a protein called COAT PROTEIN VP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	2	250	Total	C	N	O	S	8	0	0
			1961	1245	339	369	8			

- Molecule 3 is a protein called COAT PROTEIN VP3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	3	237	Total	C	N	O	S	0	0	0
			1834	1172	304	346	12			

- Molecule 4 is a protein called COAT PROTEIN VP4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	4	25	Total	C	N	O	0	0	0
			194	122	35	37			

- Molecule 5 is LAURIC ACID (three-letter code: DAO) (formula: C₁₂H₂₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	1	1	Total	C	O	0	0
			14	12	2		

- Molecule 6 is water.

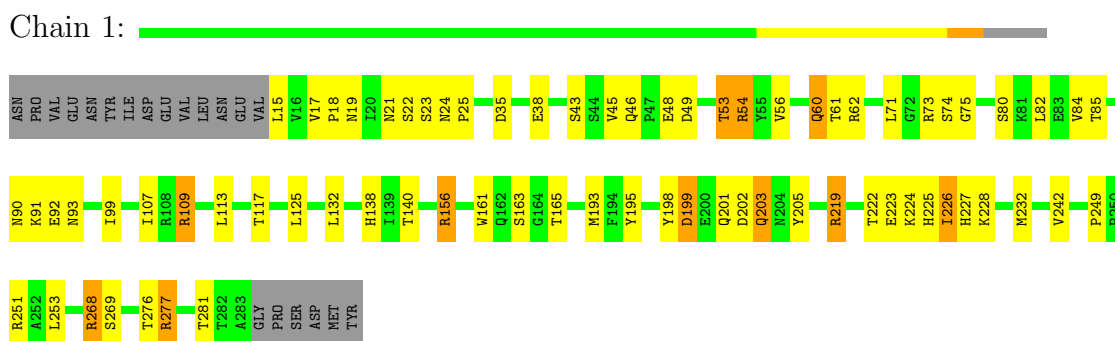
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	1	51	Total	O	0	0
			51	51		
6	2	49	Total	O	0	0
			49	49		
6	3	56	Total	O	0	0
			56	56		
6	4	3	Total	O	0	0
			3	3		

3 Residue-property plots

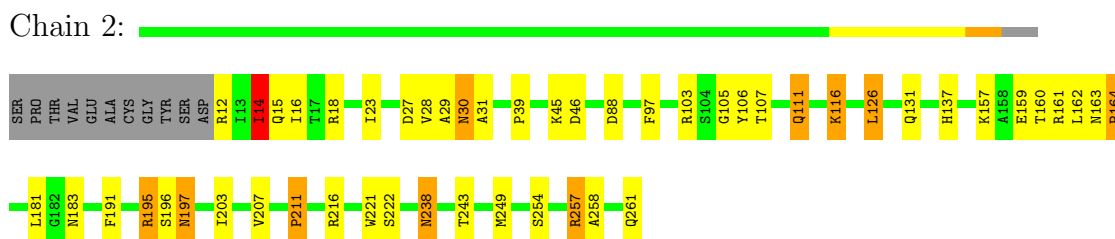
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

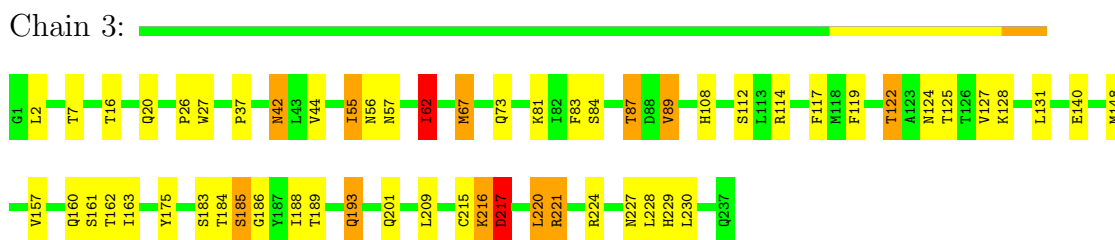
• Molecule 1: COAT PROTEIN VP1



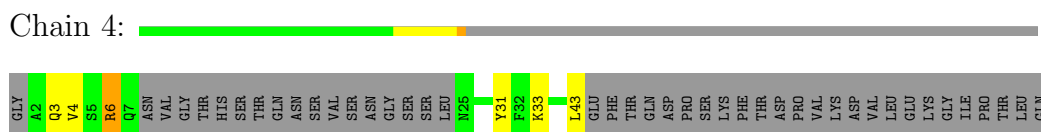
• Molecule 2: COAT PROTEIN VP2



• 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	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	308.68Å 352.98Å 380.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.60)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.180 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6316	wwPDB-VP
Average B, all atoms (Å ²)	28.0	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: DAO

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	1	0.69	1/2210 (0.0%)	0.80	1/3010 (0.0%)
2	2	0.63	2/2016 (0.1%)	0.94	4/2752 (0.1%)
3	3	0.67	1/1884 (0.1%)	0.85	3/2579 (0.1%)
4	4	0.80	0/196	0.89	0/261
All	All	0.67	4/6306 (0.1%)	0.87	8/8602 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	2	261	GLN	CG-CD	5.87	1.64	1.51
1	1	199	ASP	CB-CG	-5.52	1.40	1.51
2	2	14	ILE	CA-CB	5.51	1.67	1.54
3	3	217	ASP	CB-CG	-5.39	1.40	1.51

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	2	261	GLN	N-CA-CB	-15.18	83.27	110.60
2	2	29	ALA	N-CA-C	-9.48	85.40	111.00
3	3	220	LEU	CA-CB-CG	6.51	130.28	115.30
2	2	181	LEU	C-N-CA	-6.24	109.20	122.30
1	1	199	ASP	CB-CA-C	-6.00	98.39	110.40
3	3	89	VAL	CB-CA-C	-5.40	101.14	111.40
2	2	105	GLY	N-CA-C	-5.02	100.56	113.10
3	3	62	ILE	N-CA-C	5.01	124.53	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1	2154	0	2077	63	0
2	2	1961	0	1902	35	0
3	3	1834	0	1817	47	0
4	4	194	0	180	2	0
5	1	14	0	23	1	0
6	1	51	0	0	0	0
6	2	49	0	0	0	0
6	3	56	0	0	1	0
6	4	3	0	0	0	0
All	All	6316	0	5999	122	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 10.

All (122) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:1:268:ARG:HB3	1:1:268:ARG:HH11	1.10	1.09
1:1:268:ARG:CB	1:1:268:ARG:HH11	1.69	1.04
3:3:87:THR:HG22	3:3:185:SER:HB2	1.47	0.97
1:1:199:ASP:HB3	1:1:201:GLN:H	1.32	0.94
1:1:268:ARG:HB3	1:1:268:ARG:NH1	1.87	0.90
1:1:48:GLU:HG3	1:1:53:THR:HG21	1.51	0.89
3:3:224:ARG:HH12	3:3:227:ASN:HD22	1.18	0.87
1:1:43:SER:HB3	3:3:114:ARG:HD2	1.56	0.87
1:1:85:THR:H	1:1:93:ASN:HD21	1.23	0.87
3:3:42:ASN:HD22	3:3:44:VAL:H	1.19	0.86
2:2:207:VAL:HG22	3:3:37:PRO:HG2	1.62	0.81
1:1:199:ASP:HB2	1:1:203:GLN:H	1.46	0.80
3:3:20:GLN:HE22	4:4:31:TYR:H	1.30	0.79
3:3:122:THR:CG2	3:3:124:ASN:H	1.98	0.77
1:1:195:TYR:H	2:2:131:GLN:HE21	1.33	0.74
3:3:224:ARG:HH12	3:3:227:ASN:ND2	1.86	0.72
3:3:193:GLN:HE21	3:3:193:GLN:HA	1.53	0.72
3:3:87:THR:CG2	3:3:185:SER:HB2	2.22	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:3:184:THR:HG22	3:3:186:GLY:H	1.58	0.68
2:2:14:ILE:HB	2:2:28:VAL:HG11	1.74	0.68
1:1:23:SER:HB3	1:1:53:THR:H	1.58	0.68
2:2:197:ASN:HD22	2:2:197:ASN:N	1.91	0.68
3:3:42:ASN:ND2	3:3:44:VAL:H	1.90	0.67
1:1:21:ASN:O	1:1:53:THR:CG2	2.42	0.67
1:1:38:GLU:HA	2:2:191:PHE:HB2	1.76	0.66
2:2:203:ILE:HG23	2:2:249:MET:HE1	1.79	0.65
3:3:122:THR:HG22	3:3:124:ASN:H	1.60	0.65
1:1:46:GLN:HG3	3:3:216:LYS:HE3	1.79	0.63
1:1:276:THR:CG2	3:3:62:ILE:HG13	2.29	0.62
2:2:12:ARG:O	2:2:28:VAL:HG22	1.99	0.61
1:1:21:ASN:O	1:1:53:THR:HG23	2.01	0.61
1:1:18:PRO:HG2	3:3:216:LYS:HB2	1.83	0.60
2:2:238:ASN:H	2:2:238:ASN:HD22	1.49	0.60
3:3:62:ILE:HA	3:3:67:MET:HE3	1.83	0.60
3:3:122:THR:HG23	3:3:124:ASN:H	1.67	0.59
1:1:161:TRP:CE2	1:1:219:ARG:HG2	2.38	0.59
2:2:39:PRO:HG2	2:2:249:MET:HE3	1.84	0.59
3:3:55:ILE:HD11	3:3:83:PHE:CE2	2.38	0.59
2:2:106:TYR:C	2:2:249:MET:HE2	2.23	0.59
1:1:109:ARG:NH1	1:1:251:ARG:O	2.35	0.59
1:1:225:HIS:HD2	1:1:227:HIS:H	1.52	0.57
1:1:281:THR:O	3:3:81:LYS:HE2	2.05	0.57
2:2:16:ILE:HD12	2:2:23:ILE:HB	1.87	0.56
3:3:122:THR:HG22	3:3:124:ASN:N	2.20	0.56
2:2:162:LEU:C	2:2:164:PRO:CD	2.74	0.56
1:1:202:ASP:OD2	2:2:216:ARG:NH2	2.38	0.56
2:2:162:LEU:C	2:2:164:PRO:HD3	2.25	0.56
1:1:276:THR:HG22	3:3:62:ILE:HG13	1.86	0.56
2:2:97:PHE:HE2	2:2:216:ARG:HG3	1.71	0.56
1:1:24:ASN:HB3	1:1:25:PRO:HD2	1.88	0.56
1:1:38:GLU:HG2	2:2:197:ASN:HD21	1.71	0.55
2:2:197:ASN:ND2	2:2:197:ASN:N	2.54	0.55
2:2:18:ARG:NH1	2:2:107:THR:OG1	2.40	0.55
1:1:249:PRO:HB3	2:2:183:ASN:HB3	1.89	0.54
3:3:119:PHE:CD1	3:3:157:VAL:HG11	2.42	0.54
1:1:268:ARG:CB	1:1:268:ARG:NH1	2.54	0.54
1:1:80:SER:O	1:1:232:MET:HE3	2.08	0.53
3:3:227:ASN:H	3:3:227:ASN:ND2	2.07	0.52
3:3:62:ILE:HD13	3:3:62:ILE:O	2.09	0.52
2:2:162:LEU:O	2:2:164:PRO:N	2.41	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:1:23:SER:CB	1:1:53:THR:HG22	2.40	0.52
1:1:46:GLN:HA	1:1:46:GLN:OE1	2.09	0.51
1:1:277:ARG:HD3	3:3:57:ASN:OD1	2.10	0.51
1:1:71:LEU:HD12	1:1:242:VAL:HG11	1.91	0.51
1:1:268:ARG:HH11	1:1:268:ARG:CA	2.23	0.51
1:1:80:SER:C	1:1:232:MET:CE	2.80	0.50
2:2:238:ASN:HD22	2:2:238:ASN:N	2.07	0.50
1:1:85:THR:N	1:1:93:ASN:HD21	2.03	0.50
1:1:138:HIS:O	1:1:222:THR:HG21	2.12	0.49
3:3:20:GLN:NE2	6:3:5133:HOH:O	2.45	0.49
1:1:17:VAL:HG21	1:1:60:GLN:NE2	2.27	0.49
1:1:15:LEU:O	1:1:61:THR:HA	2.14	0.48
3:3:193:GLN:NE2	3:3:193:GLN:HA	2.27	0.48
1:1:35:ASP:O	3:3:162:THR:HB	2.13	0.48
2:2:111:GLN:HG2	2:2:243:THR:O	2.13	0.48
1:1:18:PRO:CG	3:3:216:LYS:HB2	2.43	0.48
4:4:6:ARG:O	4:4:6:ARG:HG2	2.14	0.48
1:1:90:ASN:OD1	1:1:156:ARG:NH1	2.47	0.48
1:1:225:HIS:CD2	1:1:227:HIS:H	2.31	0.47
3:3:26:PRO:O	3:3:27:TRP:HB2	2.13	0.47
1:1:80:SER:C	1:1:232:MET:HE3	2.35	0.47
1:1:18:PRO:HG2	3:3:216:LYS:HD3	1.96	0.47
1:1:84:VAL:HA	1:1:93:ASN:ND2	2.29	0.47
1:1:195:TYR:H	2:2:131:GLN:NE2	2.06	0.46
1:1:161:TRP:CD2	1:1:219:ARG:HG2	2.50	0.46
3:3:112:SER:H	3:3:217:ASP:HB3	1.80	0.46
1:1:99:ILE:CG2	5:1:6001:DAO:H81	2.45	0.46
2:2:126:LEU:HD22	2:2:221:TRP:HB3	1.96	0.46
1:1:73:ARG:HG2	3:3:16:THR:HG22	1.97	0.46
1:1:199:ASP:HB3	1:1:201:GLN:N	2.13	0.45
1:1:226:ILE:O	1:1:226:ILE:HG13	2.16	0.45
1:1:117:THR:OG1	1:1:193:MET:CE	2.65	0.45
2:2:162:LEU:C	2:2:164:PRO:N	2.68	0.45
1:1:46:GLN:OE1	3:3:215:CYS:HB3	2.16	0.45
2:2:14:ILE:HD12	2:2:30:ASN:ND2	2.31	0.44
2:2:257:ARG:HG2	2:2:258:ALA:N	2.33	0.44
1:1:45:VAL:HG13	1:1:49:ASP:HB2	2.00	0.43
1:1:253:LEU:O	3:3:229:HIS:HE1	2.01	0.43
2:2:45:LYS:HE3	2:2:46:ASP:OD2	2.19	0.43
2:2:116:LYS:O	3:3:122:THR:HG23	2.19	0.43
1:1:75:GLY:HA3	1:1:107:ILE:HD11	2.01	0.43
3:3:108:HIS:HB2	3:3:221:ARG:HG3	2.00	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:3:84:SER:HB2	3:3:189:THR:HG22	2.01	0.42
1:1:277:ARG:NH2	3:3:84:SER:O	2.49	0.42
1:1:17:VAL:HG12	1:1:62:ARG:NE	2.35	0.42
2:2:15:GLN:HA	2:2:23:ILE:O	2.20	0.42
2:2:103:ARG:HD2	2:2:211:PRO:O	2.20	0.42
1:1:17:VAL:HG12	1:1:62:ARG:CZ	2.50	0.42
3:3:117:PHE:O	3:3:162:THR:HG23	2.20	0.42
3:3:122:THR:HG22	3:3:125:THR:N	2.34	0.41
3:3:162:THR:HG22	3:3:163:ILE:N	2.34	0.41
1:1:54:ARG:HD3	1:1:56:VAL:HG22	2.02	0.41
2:2:195:ARG:HD3	3:3:161:SER:OG	2.21	0.41
3:3:127:VAL:HG22	3:3:128:LYS:N	2.36	0.41
1:1:140:THR:H	1:1:222:THR:HG23	1.85	0.41
1:1:198:TYR:CE1	1:1:205:TYR:HB2	2.56	0.41
1:1:92:GLU:HB3	1:1:93:ASN:H	1.61	0.41
3:3:62:ILE:O	3:3:67:MET:HE3	2.21	0.41
2:2:238:ASN:H	2:2:238:ASN:ND2	2.18	0.41
1:1:60:GLN:HG2	3:3:175:TYR:OH	2.21	0.41
2:2:137:HIS:CD2	2:2:137:HIS:H	2.39	0.41
2:2:196:SER:HB2	2:2:197:ASN:HD22	1.86	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	1	267/289 (92%)	256 (96%)	11 (4%)	0	100	100
2	2	248/261 (95%)	228 (92%)	17 (7%)	3 (1%)	19	39
3	3	235/237 (99%)	222 (94%)	13 (6%)	0	100	100
4	4	21/68 (31%)	21 (100%)	0	0	100	100
All	All	771/855 (90%)	727 (94%)	41 (5%)	3 (0%)	43	72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	2	160	THR
2	2	31	ALA
2	2	164	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	1	237/257 (92%)	213 (90%)	24 (10%)	11	20
2	2	216/226 (96%)	198 (92%)	18 (8%)	16	30
3	3	210/210 (100%)	183 (87%)	27 (13%)	6	11
4	4	19/59 (32%)	14 (74%)	5 (26%)	1	1
All	All	682/752 (91%)	608 (89%)	74 (11%)	9	16

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	19	ASN
1	1	22	SER
1	1	53	THR
1	1	54	ARG
1	1	60	GLN
1	1	74	SER
1	1	82	LEU
1	1	91	LYS
1	1	109	ARG
1	1	113	LEU
1	1	125	LEU
1	1	132	LEU
1	1	156	ARG
1	1	163	SER
1	1	165	THR
1	1	203	GLN
1	1	219	ARG
1	1	223	GLU
1	1	224	LYS
1	1	226	ILE

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Mol	Chain	Res	Type
1	1	228	LYS
1	1	268	ARG
1	1	269	SER
1	1	277	ARG
2	2	14	ILE
2	2	27	ASP
2	2	30	ASN
2	2	88	ASP
2	2	111	GLN
2	2	116	LYS
2	2	126	LEU
2	2	157	LYS
2	2	159	GLU
2	2	161	ARG
2	2	163	ASN
2	2	195	ARG
2	2	197	ASN
2	2	211	PRO
2	2	222	SER
2	2	238	ASN
2	2	254	SER
2	2	257	ARG
3	3	2	LEU
3	3	7	THR
3	3	42	ASN
3	3	55	ILE
3	3	56	ASN
3	3	62	ILE
3	3	67	MET
3	3	73	GLN
3	3	87	THR
3	3	89	VAL
3	3	122	THR
3	3	131	LEU
3	3	140	GLU
3	3	148	MET
3	3	160	GLN
3	3	183	SER
3	3	185	SER
3	3	188	ILE
3	3	193	GLN
3	3	201	GLN

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Mol	Chain	Res	Type
3	3	209	LEU
3	3	216	LYS
3	3	217	ASP
3	3	220	LEU
3	3	221	ARG
3	3	228	LEU
3	3	230	LEU
4	4	3	GLN
4	4	4	VAL
4	4	6	ARG
4	4	33	LYS
4	4	43	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	19	ASN
1	1	93	ASN
1	1	162	GLN
1	1	225	HIS
2	2	15	GLN
2	2	30	ASN
2	2	40	HIS
2	2	131	GLN
2	2	193	ASN
2	2	197	ASN
2	2	219	ASN
2	2	220	ASN
2	2	238	ASN
3	3	20	GLN
3	3	42	ASN
3	3	48	GLN
3	3	56	ASN
3	3	193	GLN
3	3	227	ASN
3	3	229	HIS
4	4	30	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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	DAO	1	6001	-	13,13,13	1.93	3 (23%)	13,13,13	0.88	1 (7%)

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	DAO	1	6001	-	-	0/11/11/11	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	1	6001	DAO	C2-C1	4.82	1.62	1.50
5	1	6001	DAO	O1-C1	2.20	1.30	1.22
5	1	6001	DAO	C3-C2	2.08	1.60	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	1	6001	DAO	O2-C1-C2	2.09	121.61	114.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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 will therefore be empty.

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

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

6.3 Carbohydrates ⓘ

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

6.4 Ligands ⓘ

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

6.5 Other polymers ⓘ

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