



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:35 PM GMT

PDB ID : 3ZWV
Title : Crystal structure of ADP-ribosyl cyclase complexed with ara-2'F-ADP- ribose
at 2.3 angstrom
Authors : Kotaka, M.; Graeff, R.; Zhang, L.H.; Lee, H.C.; Hao, Q.
Deposited on : 2011-08-03
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

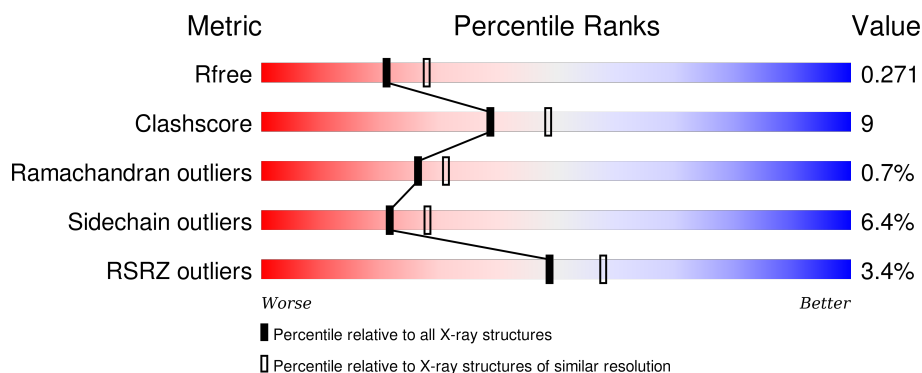
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
R_{free}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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.

Mol	Chain	Length	Quality of chain
1	A	260	<div> <div>5%</div> <div>75%</div> <div>20%</div> <div>• •</div> </div>
1	B	260	<div> <div>%</div> <div>80%</div> <div>15%</div> <div>• •</div> </div>
1	C	260	<div> <div>3%</div> <div>73%</div> <div>23%</div> <div>• •</div> </div>
1	D	260	<div> <div>2%</div> <div>78%</div> <div>17%</div> <div>• •</div> </div>
1	E	260	<div> <div>%</div> <div>81%</div> <div>15%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	260	<div><div>8%</div><div><div></div><div>66%</div><div>26%</div><div></div><div></div></div><div><div></div><div></div></div></div>
1	G	260	<div><div>%</div><div><div></div><div>77%</div><div>18%</div><div></div><div></div></div><div><div></div><div></div></div></div>
1	H	260	<div><div>6%</div><div><div></div><div>71%</div><div>23%</div><div></div><div></div></div><div><div></div><div></div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16687 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 ADP-RIBOSYL CYCLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	253	Total	C	N	O	S	0	0	1
			2018	1291	344	369	14			
1	B	252	Total	C	N	O	S	0	0	1
			2013	1288	343	368	14			
1	C	252	Total	C	N	O	S	0	0	1
			2013	1288	343	368	14			
1	D	252	Total	C	N	O	S	0	0	0
			2017	1291	343	369	14			
1	E	252	Total	C	N	O	S	0	0	1
			2013	1288	343	368	14			
1	F	252	Total	C	N	O	S	0	0	1
			2013	1288	343	368	14			
1	G	252	Total	C	N	O	S	0	0	1
			2013	1288	343	368	14			
1	H	252	Total	C	N	O	S	0	0	1
			2013	1288	343	368	14			

There are 16 discrepancies between the modelled and reference sequences:

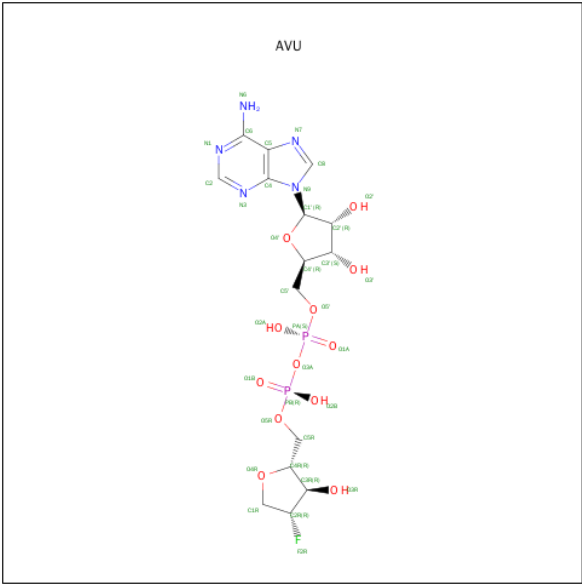
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ALA	-	EXPRESSION TAG	UNP P29241
A	0	ALA	-	EXPRESSION TAG	UNP P29241
B	-1	ALA	-	EXPRESSION TAG	UNP P29241
B	0	ALA	-	EXPRESSION TAG	UNP P29241
C	-1	ALA	-	EXPRESSION TAG	UNP P29241
C	0	ALA	-	EXPRESSION TAG	UNP P29241
D	-1	ALA	-	EXPRESSION TAG	UNP P29241
D	0	ALA	-	EXPRESSION TAG	UNP P29241
E	-1	ALA	-	EXPRESSION TAG	UNP P29241
E	0	ALA	-	EXPRESSION TAG	UNP P29241
F	-1	ALA	-	EXPRESSION TAG	UNP P29241
F	0	ALA	-	EXPRESSION TAG	UNP P29241
G	-1	ALA	-	EXPRESSION TAG	UNP P29241

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Chain	Residue	Modelled	Actual	Comment	Reference
G	0	ALA	-	EXPRESSION TAG	UNP P29241
H	-1	ALA	-	EXPRESSION TAG	UNP P29241
H	0	ALA	-	EXPRESSION TAG	UNP P29241

- Molecule 2 is [(2R,3S,4R,5R)-5-(6-AMINO-9H-PURIN-9-YL)-3,4-DIHYDROXYTETRAHYDROFURAN-2-YL]METHYL [(2R,3R,4R)-4-FLUORO-3-HYDROXYTETRAHYDROFURAN-2-YL]METHYL DIHYDROGEN DIPHOSPHATE (three-letter code: AVU) (formula: C₁₅H₂₂FN₅O₁₂P₂).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	F	N	O	P		
2	A	1	Total 35	15	1	5	12	2	0	0
2	B	1	Total 35	15	1	5	12	2	0	0
2	C	1	Total 35	15	1	5	12	2	0	0
2	D	1	Total 35	15	1	5	12	2	0	0
2	E	1	Total 35	15	1	5	12	2	0	0
2	F	1	Total 35	15	1	5	12	2	0	0
2	G	1	Total 35	15	1	5	12	2	0	0
2	H	1	Total 35	15	1	5	12	2	0	0

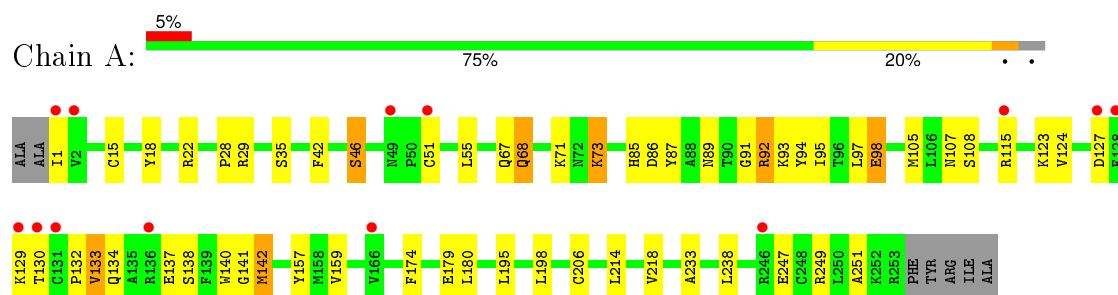
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	35	Total 35	O 35	0	0
3	B	47	Total 47	O 47	0	0
3	C	32	Total 32	O 32	0	0
3	D	49	Total 49	O 49	0	0
3	E	40	Total 40	O 40	0	0
3	F	19	Total 19	O 19	0	0
3	G	42	Total 42	O 42	0	0
3	H	30	Total 30	O 30	0	0

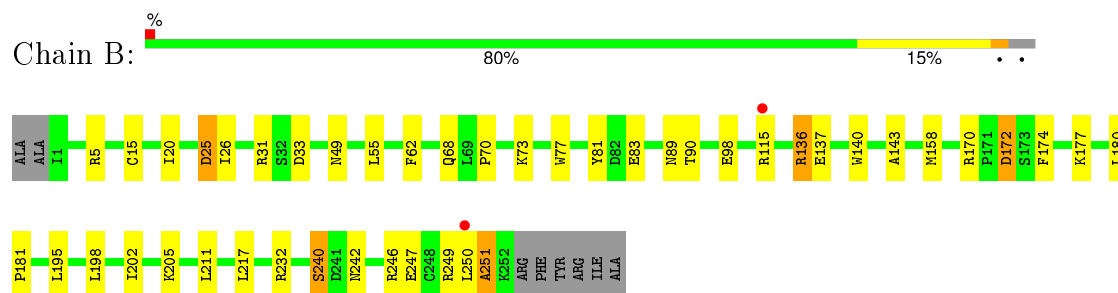
3 Residue-property plots [i](#)

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.

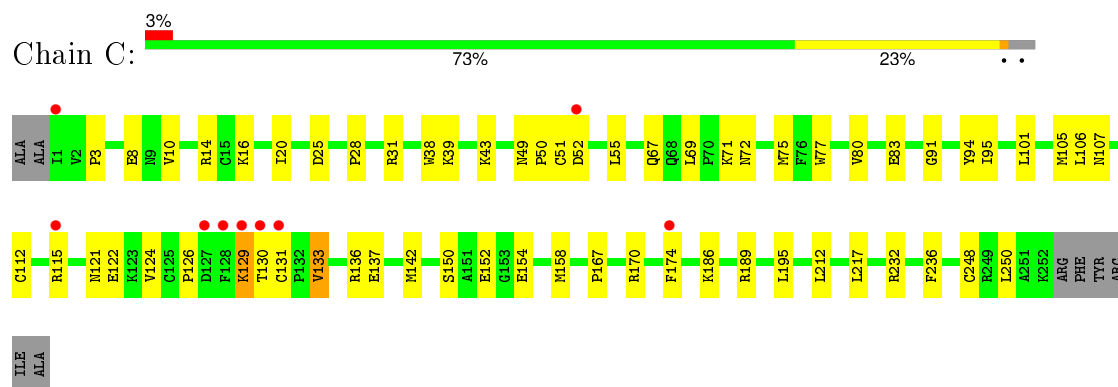
• Molecule 1: ADP-RIBOSYL CYCLASE



• Molecule 1: ADP-RIBOSYL CYCLASE

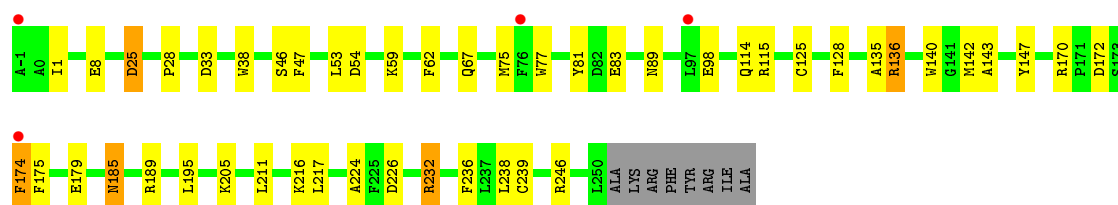


• Molecule 1: ADP-RIBOSYL CYCLASE

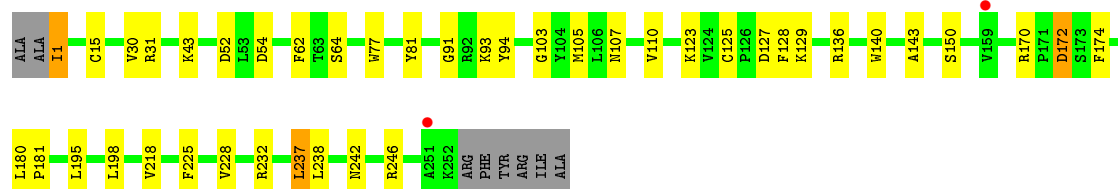
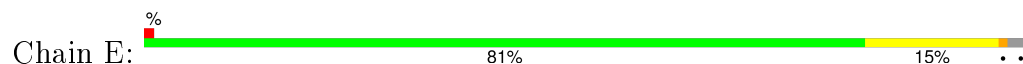


• Molecule 1: ADP-RIBOSYL CYCLASE

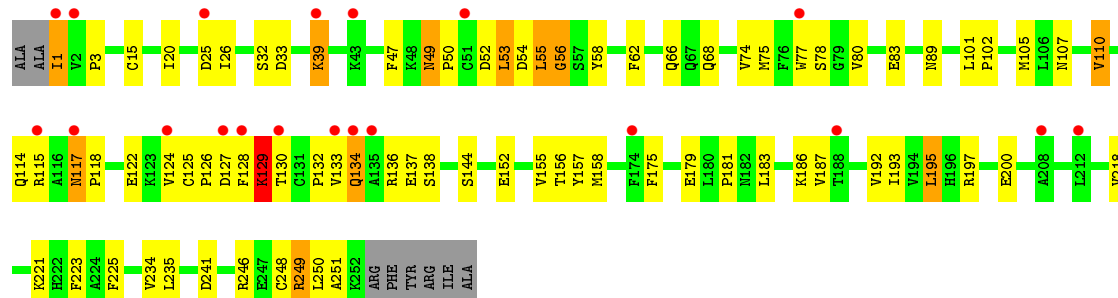




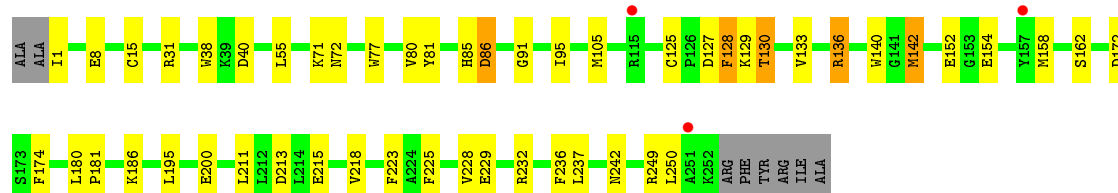
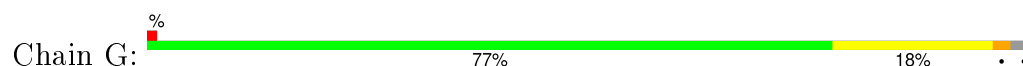
• Molecule 1: ADP-RIBOSYL CYCLASE



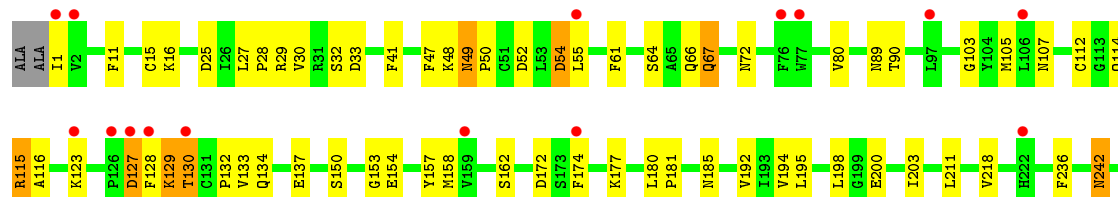
• Molecule 1: ADP-RIBOSYL CYCLASE

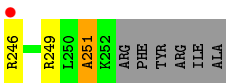


• Molecule 1: ADP-RIBOSYL CYCLASE



• Molecule 1: ADP-RIBOSYL CYCLASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	61.15Å 76.48Å 141.37Å 88.09° 90.88° 91.02°	Depositor
Resolution (Å)	30.00 – 2.30 29.78 – 2.30	Depositor EDS
% Data completeness (in resolution range)	87.5 (30.00-2.30) 68.1 (29.78-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.215 , 0.277 0.212 , 0.271	Depositor DCC
R_{free} test set	3933 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	44.1	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.1	EDS
Estimated twinning fraction	0.079 for h,-k,-l 0.022 for -h,k,-l 0.012 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 77905 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16687	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.92	2/2070 (0.1%)	0.85	0/2802
1	B	1.00	2/2065 (0.1%)	0.89	0/2795
1	C	0.83	0/2065	0.85	0/2795
1	D	1.01	2/2069 (0.1%)	0.90	0/2800
1	E	0.97	0/2065	0.90	0/2795
1	F	0.80	1/2065 (0.0%)	0.84	2/2795 (0.1%)
1	G	1.00	2/2065 (0.1%)	0.90	1/2795 (0.0%)
1	H	0.79	1/2065 (0.0%)	0.80	0/2795
All	All	0.92	10/16529 (0.1%)	0.87	3/22372 (0.0%)

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
1	E	0	1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	206	CYS	CB-SG	7.45	1.95	1.82
1	B	15	CYS	CB-SG	-5.66	1.72	1.81
1	D	174	PHE	CE2-CZ	5.62	1.48	1.37
1	G	172	ASP	CB-CG	5.54	1.63	1.51
1	B	140	TRP	CB-CG	5.51	1.60	1.50
1	G	86	ASP	CB-CG	-5.51	1.40	1.51
1	D	147	TYR	CD2-CE2	-5.33	1.31	1.39
1	H	251	ALA	C-N	-5.29	1.21	1.34
1	A	140	TRP	CB-CG	5.26	1.59	1.50
1	F	251	ALA	C-N	-5.01	1.22	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	20	ILE	CG1-CB-CG2	-6.29	97.56	111.40
1	F	110	VAL	CB-CA-C	-5.89	100.22	111.40
1	G	213	ASP	CB-CG-OD1	5.28	123.05	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	123	LYS	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	2018	0	1970	35	0
1	B	2013	0	1968	23	0
1	C	2013	0	1968	42	0
1	D	2017	0	1973	33	0
1	E	2013	0	1968	37	0
1	F	2013	0	1968	56	0
1	G	2013	0	1968	24	0
1	H	2013	0	1968	44	0
2	A	35	0	19	2	0
2	B	35	0	19	6	0
2	C	35	0	19	2	0
2	D	35	0	19	5	0
2	E	35	0	19	3	0
2	F	35	0	19	5	0
2	G	35	0	19	0	0
2	H	35	0	19	0	0
3	A	35	0	0	0	0
3	B	47	0	0	4	0
3	C	32	0	0	1	0
3	D	49	0	0	2	0
3	E	40	0	0	3	0
3	F	19	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	42	0	0	1	0
3	H	30	0	0	1	0
All	All	16687	0	15903	289	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 9.

All (289) 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)
1:A:28:PRO:HG3	1:A:68:GLN:NE2	1.56	1.19
1:D:128:PHE:O	1:D:136:ARG:HD2	1.43	1.17
1:H:115:ARG:HG3	1:H:115:ARG:HH11	1.11	1.10
2:B:301:AVU:C2	3:B:2035:HOH:O	1.98	1.09
1:A:28:PRO:CG	1:A:68:GLN:HE22	1.66	1.07
1:A:28:PRO:HG3	1:A:68:GLN:HE22	0.87	0.99
2:B:301:AVU:H2	3:B:2035:HOH:O	1.60	0.95
1:H:30:VAL:HG21	1:H:64:SER:O	1.66	0.95
1:A:89:ASN:O	1:A:92:ARG:HD3	1.68	0.94
1:C:137:GLU:HG2	1:C:174:PHE:CZ	2.03	0.94
1:E:1:ILE:HG23	1:E:110:VAL:HG11	1.47	0.93
1:H:115:ARG:HG3	1:H:115:ARG:NH1	1.86	0.89
1:A:68:GLN:H	1:A:68:GLN:CD	1.75	0.86
1:C:136:ARG:HG3	1:C:137:GLU:HG3	1.59	0.84
1:F:129:LYS:H	1:F:129:LYS:HD3	1.43	0.83
1:C:133:VAL:O	1:C:136:ARG:HG2	1.80	0.81
1:E:180:LEU:HB3	1:E:181:PRO:HD3	1.60	0.81
1:F:49:ASN:HB2	1:F:52:ASP:HB2	1.63	0.80
1:B:98:GLU:OE2	2:B:301:AVU:N6	2.16	0.79
1:C:137:GLU:HG2	1:C:174:PHE:HZ	1.45	0.78
1:F:127:ASP:HB3	1:F:130:THR:OG1	1.84	0.77
1:D:128:PHE:O	1:D:136:ARG:CD	2.30	0.77
2:B:301:AVU:N1	3:B:2035:HOH:O	2.12	0.75
1:H:246:ARG:HG3	1:H:249:ARG:HH21	1.50	0.75
1:F:74:VAL:HG22	3:F:2008:HOH:O	1.85	0.74
1:F:117:ASN:HD22	1:F:118:PRO:HA	1.50	0.74
1:H:137:GLU:HG2	1:H:174:PHE:CZ	2.22	0.74
1:F:15:CYS:HB2	1:F:105:MET:SD	2.29	0.73
1:E:1:ILE:HG22	1:E:125:CYS:HB2	1.69	0.73
1:C:49:ASN:HD22	1:C:115:ARG:HH21	1.34	0.73
1:C:174:PHE:CD2	2:C:301:AVU:O2'	2.43	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:54:ASP:HB2	3:D:2017:HOH:O	1.91	0.70
1:E:172:ASP:OD2	3:E:2034:HOH:O	2.11	0.69
1:D:195:LEU:HD11	1:D:238:LEU:HD11	1.76	0.68
1:F:74:VAL:HG23	1:F:155:VAL:HG12	1.76	0.68
1:H:30:VAL:HG22	1:H:32:SER:H	1.59	0.67
1:F:80:VAL:HG21	1:F:158:MET:HG2	1.75	0.67
1:F:133:VAL:O	1:F:137:GLU:HG3	1.95	0.66
1:G:127:ASP:HB3	1:G:130:THR:HG22	1.76	0.66
1:A:107:ASN:O	1:A:108:SER:HB2	1.96	0.66
1:G:128:PHE:N	1:G:128:PHE:CD1	2.63	0.66
1:C:250:LEU:HD11	1:D:236:PHE:CZ	2.30	0.65
1:C:25:ASP:OD2	3:C:2008:HOH:O	2.14	0.65
1:F:74:VAL:CG2	1:F:155:VAL:HG12	2.26	0.65
1:G:127:ASP:O	1:G:130:THR:HG23	1.96	0.65
1:E:140:TRP:CB	1:E:174:PHE:HE2	2.10	0.65
1:C:250:LEU:HD13	1:D:232:ARG:CZ	2.26	0.65
1:F:129:LYS:N	1:F:129:LYS:HD3	2.11	0.64
1:E:1:ILE:CG2	1:E:110:VAL:HG11	2.25	0.64
1:D:128:PHE:HB3	1:D:136:ARG:HG3	1.79	0.64
1:E:170:ARG:HD2	1:E:172:ASP:OD2	1.97	0.63
1:G:140:TRP:HB3	1:G:174:PHE:HE2	1.62	0.63
1:H:80:VAL:HG21	1:H:158:MET:HG2	1.81	0.63
1:C:49:ASN:ND2	1:C:115:ARG:HH21	1.97	0.63
1:B:205:LYS:O	1:B:211:LEU:HD12	1.99	0.62
1:F:152:GLU:OE2	1:F:186:LYS:HE3	2.00	0.62
1:F:246:ARG:O	1:F:249:ARG:HG3	1.99	0.62
1:E:128:PHE:O	1:E:136:ARG:HD2	1.99	0.62
1:F:3:PRO:HA	1:F:110:VAL:HB	1.81	0.62
1:H:137:GLU:HG2	1:H:174:PHE:HZ	1.63	0.61
1:E:91:GLY:HA2	1:E:94:TYR:O	2.01	0.61
1:C:232:ARG:HG3	1:C:232:ARG:HH11	1.65	0.61
1:A:15:CYS:HB2	1:A:105:MET:SD	2.41	0.60
1:E:140:TRP:HB2	1:E:174:PHE:HE2	1.65	0.60
1:E:127:ASP:OD2	1:E:129:LYS:HG2	2.01	0.60
1:G:77:TRP:CH2	1:G:81:TYR:HD2	2.20	0.60
1:H:128:PHE:H	1:H:130:THR:HG22	1.67	0.60
1:B:83:GLU:HG2	1:B:195:LEU:HD11	1.83	0.59
1:E:140:TRP:CH2	2:E:301:AVU:N6	2.70	0.59
1:B:49:ASN:HD21	1:B:115:ARG:HH21	1.49	0.59
1:A:133:VAL:HG13	1:A:134:GLN:H	1.67	0.58
1:F:49:ASN:HB2	1:F:52:ASP:CB	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:174:PHE:HB2	2:E:301:AVU:PB	2.45	0.57
1:F:80:VAL:HG12	1:F:197:ARG:NH2	2.19	0.57
1:F:181:PRO:HA	1:F:221:LYS:NZ	2.20	0.57
1:B:55:LEU:O	1:B:55:LEU:HG	2.06	0.56
1:H:198:LEU:HD12	1:H:242:ASN:HB3	1.87	0.56
1:H:174:PHE:N	1:H:174:PHE:HD1	2.03	0.56
1:A:87:TYR:CE1	1:A:93:LYS:HE2	2.41	0.56
1:B:136:ARG:HG2	1:B:137:GLU:HG3	1.88	0.56
1:G:15:CYS:HB2	1:G:105:MET:SD	2.46	0.56
1:G:55:LEU:HD12	1:G:142:MET:HG3	1.87	0.56
1:G:215:GLU:HG3	1:G:225:PHE:CD2	2.41	0.56
1:H:41:PHE:HB2	1:H:61:PHE:CD1	2.41	0.56
1:C:80:VAL:HB	1:C:83:GLU:HG2	1.85	0.56
1:H:72:ASN:OD1	1:H:153:GLY:HA3	2.05	0.55
1:A:85:HIS:NE2	1:A:98:GLU:HG2	2.21	0.55
1:E:198:LEU:HD12	1:E:242:ASN:HB3	1.88	0.55
1:E:15:CYS:HB2	1:E:105:MET:SD	2.47	0.55
1:F:54:ASP:O	1:F:56:GLY:N	2.37	0.55
1:H:174:PHE:CD1	1:H:174:PHE:N	2.73	0.55
1:D:170:ARG:HD2	1:D:172:ASP:OD1	2.07	0.54
1:E:30:VAL:HG21	1:E:64:SER:O	2.06	0.54
1:A:127:ASP:O	1:A:130:THR:HG22	2.07	0.54
1:A:28:PRO:HG2	1:A:67:GLN:HB3	1.89	0.54
1:C:167:PRO:HD2	1:C:170:ARG:HD3	1.91	0.54
1:H:47:PHE:HA	1:H:114:GLN:O	2.09	0.53
1:H:103:GLY:O	1:H:107:ASN:HB2	2.08	0.53
2:F:301:AVU:O1B	2:F:301:AVU:O3'	2.27	0.53
1:E:77:TRP:CH2	1:E:81:TYR:HD2	2.27	0.53
1:D:28:PRO:HG2	1:D:67:GLN:HB2	1.91	0.53
1:D:83:GLU:HG2	1:D:195:LEU:HD21	1.90	0.52
1:G:133:VAL:HA	1:G:136:ARG:HG3	1.90	0.52
1:A:68:GLN:N	1:A:68:GLN:CD	2.51	0.52
1:A:198:LEU:HD22	1:A:247:GLU:CG	2.40	0.52
1:H:246:ARG:HG3	1:H:249:ARG:NH2	2.21	0.52
1:F:144:SER:OG	2:F:301:AVU:H1RA	2.10	0.52
1:E:1:ILE:HG23	1:E:110:VAL:CG1	2.31	0.52
1:D:77:TRP:CH2	1:D:81:TYR:HD2	2.28	0.52
1:G:77:TRP:HA	1:G:158:MET:O	2.10	0.51
1:H:180:LEU:N	1:H:181:PRO:HD2	2.25	0.51
1:C:121:ASN:HD21	1:C:124:VAL:HG12	1.76	0.51
1:G:180:LEU:HB3	1:G:181:PRO:HD3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:ILE:HD11	3:B:2032:HOH:O	2.11	0.50
1:H:115:ARG:CG	1:H:115:ARG:HH11	1.99	0.50
1:B:77:TRP:CH2	1:B:81:TYR:HD2	2.29	0.50
1:H:66:GLN:HB3	3:H:2008:HOH:O	2.11	0.50
1:F:58:TYR:O	1:F:62:PHE:HD2	1.95	0.50
1:H:128:PHE:N	1:H:130:THR:HG22	2.25	0.50
1:C:16:LYS:HD3	1:C:20:ILE:HD12	1.94	0.50
1:C:152:GLU:HG2	1:C:186:LYS:HB3	1.92	0.50
1:C:137:GLU:OE2	1:C:174:PHE:CE1	2.64	0.49
1:F:33:ASP:OD1	1:F:33:ASP:C	2.51	0.49
1:C:3:PRO:HG2	1:C:122:GLU:O	2.12	0.49
1:B:90:THR:HG22	1:B:90:THR:O	2.12	0.49
1:G:152:GLU:HG2	1:G:186:LYS:HB3	1.95	0.49
1:H:128:PHE:C	1:H:130:THR:N	2.63	0.49
1:G:91:GLY:HA2	1:G:95:ILE:HD13	1.94	0.49
1:A:174:PHE:CD2	2:A:301:AVU:H2'	2.47	0.49
1:F:128:PHE:C	1:F:130:THR:H	2.15	0.48
1:A:42:PHE:O	1:A:46:SER:HB2	2.13	0.48
1:B:25:ASP:OD1	1:B:25:ASP:N	2.44	0.48
1:A:91:GLY:HA2	1:A:95:ILE:HD13	1.95	0.48
1:C:55:LEU:HG	1:C:142:MET:HE2	1.96	0.48
2:D:301:AVU:H5'	3:D:2048:HOH:O	2.13	0.48
1:E:1:ILE:CG2	1:E:110:VAL:CG1	2.92	0.48
1:C:154:GLU:HG2	1:C:189:ARG:HB3	1.96	0.48
1:E:103:GLY:O	1:E:107:ASN:HB2	2.14	0.48
1:C:137:GLU:CG	1:C:174:PHE:CZ	2.88	0.47
1:F:129:LYS:HG3	1:F:136:ARG:HD3	1.95	0.47
1:F:128:PHE:C	1:F:130:THR:N	2.67	0.47
1:E:140:TRP:HB2	1:E:174:PHE:CE2	2.48	0.47
1:B:70:PRO:HB2	1:B:73:LYS:HD3	1.95	0.47
1:F:181:PRO:HA	1:F:221:LYS:HZ1	1.79	0.47
1:H:49:ASN:OD1	1:H:50:PRO:HD2	2.15	0.47
1:A:137:GLU:O	1:A:141:GLY:HA3	2.14	0.47
1:F:1:ILE:HG22	1:F:125:CYS:O	2.15	0.47
1:C:49:ASN:HD22	1:C:115:ARG:NH2	2.09	0.47
1:F:39:LYS:NZ	1:F:39:LYS:HA	2.29	0.47
1:B:174:PHE:HB2	2:B:301:AVU:PB	2.54	0.47
1:C:236:PHE:CD2	1:D:239:CYS:HB3	2.49	0.47
1:F:101:LEU:N	1:F:102:PRO:HD2	2.30	0.47
1:H:127:ASP:HB2	1:H:130:THR:HG22	1.97	0.47
1:H:72:ASN:HA	1:H:154:GLU:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:218:VAL:HG11	1:E:225:PHE:HB2	1.95	0.47
1:G:229:GLU:OE2	3:G:2036:HOH:O	2.20	0.47
1:D:189:ARG:HB2	1:D:224:ALA:HB3	1.97	0.47
1:B:170:ARG:NH1	1:B:172:ASP:OD2	2.47	0.47
1:G:72:ASN:HA	1:G:154:GLU:O	2.15	0.47
1:H:129:LYS:H	1:H:129:LYS:HG2	1.52	0.47
1:A:133:VAL:HG13	1:A:134:GLN:N	2.30	0.47
1:A:233:ALA:O	1:B:240:SER:HB3	2.14	0.46
1:A:198:LEU:HD22	1:A:247:GLU:HG2	1.97	0.46
1:F:218:VAL:HG11	1:F:225:PHE:HB2	1.97	0.46
1:E:128:PHE:O	1:E:136:ARG:CD	2.62	0.46
1:A:174:PHE:HB2	2:A:301:AVU:PB	2.56	0.46
1:D:174:PHE:HB2	2:D:301:AVU:PB	2.55	0.46
1:F:39:LYS:HA	1:F:39:LYS:HZ3	1.80	0.46
3:E:2024:HOH:O	1:F:241:ASP:HA	2.16	0.46
1:H:133:VAL:HG13	1:H:134:GLN:N	2.30	0.46
1:C:71:LYS:O	1:C:72:ASN:HB2	2.16	0.46
1:A:214:LEU:O	1:A:218:VAL:HG23	2.15	0.46
1:G:127:ASP:O	1:G:130:THR:CG2	2.61	0.46
1:B:62:PHE:CZ	1:B:143:ALA:HB2	2.51	0.46
1:B:205:LYS:C	1:B:211:LEU:HD12	2.36	0.46
1:G:218:VAL:HG13	1:G:223:PHE:HB2	1.98	0.46
1:A:28:PRO:CB	1:A:68:GLN:HE22	2.26	0.46
1:C:248:CYS:O	1:C:250:LEU:HD12	2.16	0.46
1:F:83:GLU:HG2	1:F:195:LEU:HD21	1.97	0.46
1:E:93:LYS:HG2	1:E:94:TYR:CE1	2.51	0.45
1:D:62:PHE:CZ	1:D:143:ALA:HB2	2.52	0.45
1:D:8:GLU:HG3	1:D:38:TRP:CE2	2.52	0.45
1:F:175:PHE:HA	1:F:179:GLU:HB2	1.99	0.45
1:D:1:ILE:HB	1:D:125:CYS:O	2.16	0.45
1:F:249:ARG:NH2	3:F:2017:HOH:O	2.49	0.45
1:H:48:LYS:HE2	1:H:52:ASP:O	2.16	0.45
1:F:55:LEU:CD2	1:F:55:LEU:O	2.65	0.45
1:F:183:LEU:HD22	1:F:187:VAL:HG21	1.99	0.45
1:E:180:LEU:HB3	1:E:181:PRO:CD	2.39	0.45
1:A:93:LYS:HG2	1:A:94:TYR:CE1	2.51	0.45
1:A:51:CYS:HB3	1:A:132:PRO:HD2	1.99	0.45
1:D:216:LYS:HB2	1:D:216:LYS:NZ	2.31	0.45
1:H:128:PHE:H	1:H:130:THR:CG2	2.30	0.44
1:D:47:PHE:HA	1:D:114:GLN:O	2.17	0.44
1:B:77:TRP:HA	1:B:158:MET:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:140:TRP:CZ3	2:D:301:AVU:N6	2.85	0.44
1:F:75:MET:HE3	1:F:156:THR:CG2	2.48	0.44
1:B:198:LEU:HD22	1:B:247:GLU:HB2	1.99	0.44
1:E:140:TRP:HB3	1:E:174:PHE:HE2	1.80	0.44
1:H:15:CYS:HB2	1:H:105:MET:SD	2.58	0.44
1:D:98:GLU:OE2	2:D:301:AVU:N1	2.51	0.44
1:B:180:LEU:HB3	1:B:181:PRO:HD3	1.98	0.44
1:H:194:VAL:HG21	1:H:211:LEU:HD11	2.00	0.44
1:D:189:ARG:HD3	1:D:226:ASP:OD2	2.17	0.44
1:C:77:TRP:HA	1:C:158:MET:O	2.17	0.44
1:D:53:LEU:HB2	1:D:135:ALA:HA	1.99	0.44
1:D:170:ARG:NH1	1:D:172:ASP:OD1	2.50	0.43
1:G:250:LEU:O	1:H:251:ALA:HB2	2.18	0.43
1:G:80:VAL:O	1:G:81:TYR:C	2.56	0.43
1:F:78:SER:OG	2:F:301:AVU:O2B	2.21	0.43
1:D:174:PHE:CE2	2:D:301:AVU:O2'	2.71	0.43
1:A:127:ASP:O	1:A:130:THR:N	2.43	0.43
1:C:51:CYS:HB3	1:C:131:CYS:HB3	1.93	0.43
1:F:53:LEU:N	1:F:53:LEU:HD12	2.33	0.43
1:D:195:LEU:HA	1:D:195:LEU:HD12	1.64	0.43
1:D:217:LEU:HD11	1:E:129:LYS:HG3	1.99	0.43
1:H:114:GLN:HG3	1:H:116:ALA:O	2.18	0.43
1:C:39:LYS:O	1:C:43:LYS:HG2	2.19	0.43
1:F:75:MET:CE	1:F:156:THR:HG22	2.48	0.43
1:H:11:PHE:C	1:H:11:PHE:CD1	2.92	0.43
1:E:62:PHE:CZ	1:E:143:ALA:HB2	2.54	0.43
1:F:246:ARG:HA	1:F:246:ARG:NE	2.34	0.43
1:C:8:GLU:OE1	1:C:38:TRP:NE1	2.52	0.43
1:C:50:PRO:CB	1:C:126:PRO:HG2	2.49	0.42
1:C:236:PHE:O	1:D:236:PHE:HB3	2.20	0.42
1:F:144:SER:OG	2:F:301:AVU:C1R	2.67	0.42
1:H:50:PRO:HB3	1:H:112:CYS:HB2	2.01	0.42
1:F:47:PHE:O	1:F:115:ARG:HD3	2.19	0.42
1:A:71:LYS:HE3	1:A:71:LYS:HB3	1.88	0.42
1:B:217:LEU:HA	1:B:217:LEU:HD23	1.90	0.42
1:B:250:LEU:O	1:B:251:ALA:HB3	2.20	0.42
1:F:187:VAL:O	1:F:223:PHE:CD1	2.72	0.42
1:F:47:PHE:HA	1:F:114:GLN:O	2.19	0.42
1:C:10:VAL:O	1:C:14:ARG:HG3	2.19	0.42
1:C:28:PRO:HG2	1:C:67:GLN:HB3	2.02	0.42
1:B:172:ASP:O	1:B:177:LYS:HE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:CYS:SG	1:C:126:PRO:HD2	2.59	0.42
1:H:162:SER:HA	1:H:203:ILE:HG12	2.00	0.42
1:E:195:LEU:CD1	1:E:238:LEU:HD11	2.50	0.42
1:E:232:ARG:NH2	1:F:248:CYS:O	2.53	0.42
1:A:18:TYR:HA	1:A:22:ARG:CG	2.50	0.42
1:H:157:TYR:O	1:H:192:VAL:HA	2.20	0.42
1:G:8:GLU:HG3	1:G:38:TRP:CE2	2.55	0.42
1:C:174:PHE:CE2	2:C:301:AVU:O2'	2.72	0.42
1:E:180:LEU:CB	1:E:181:PRO:HD3	2.40	0.42
1:H:137:GLU:OE2	1:H:174:PHE:CE1	2.73	0.42
1:E:174:PHE:HB2	2:E:301:AVU:O1B	2.20	0.42
1:A:87:TYR:CZ	1:A:93:LYS:HE2	2.55	0.42
1:E:54:ASP:HB2	3:E:2015:HOH:O	2.20	0.42
1:F:187:VAL:O	1:F:223:PHE:HD1	2.03	0.42
1:C:106:LEU:O	1:C:107:ASN:C	2.58	0.42
1:G:140:TRP:CB	1:G:174:PHE:HE2	2.29	0.41
1:F:234:VAL:O	1:F:235:LEU:C	2.58	0.41
1:F:54:ASP:HB2	1:F:134:GLN:HG3	2.02	0.41
1:A:73:LYS:HD3	1:A:73:LYS:HA	1.85	0.41
1:E:237:LEU:HD12	1:E:237:LEU:HA	1.76	0.41
1:G:236:PHE:HB3	1:H:236:PHE:O	2.20	0.41
1:C:69:LEU:HG	1:C:150:SER:HB3	2.02	0.41
1:H:54:ASP:N	1:H:54:ASP:OD1	2.51	0.41
1:A:179:GLU:O	1:A:180:LEU:C	2.59	0.41
1:C:133:VAL:O	1:C:136:ARG:CG	2.60	0.41
1:E:128:PHE:HE1	1:E:129:LYS:HD3	1.85	0.41
1:E:30:VAL:CG2	1:E:64:SER:O	2.68	0.41
1:F:197:ARG:HD2	1:F:200:GLU:OE1	2.20	0.41
1:A:195:LEU:HD11	1:A:238:LEU:HD21	2.02	0.41
1:H:115:ARG:CG	1:H:115:ARG:NH1	2.66	0.41
1:B:174:PHE:HB2	2:B:301:AVU:O5R	2.21	0.41
1:C:91:GLY:HA2	1:C:94:TYR:O	2.21	0.41
1:F:157:TYR:O	1:F:192:VAL:HA	2.21	0.41
1:F:49:ASN:CB	1:F:52:ASP:HB2	2.43	0.41
1:H:177:LYS:O	1:H:181:PRO:HG3	2.22	0.41
1:D:46:SER:O	1:D:47:PHE:HB2	2.20	0.41
1:D:25:ASP:N	1:D:25:ASP:OD1	2.49	0.41
1:C:75:MET:HB3	1:C:95:ILE:O	2.21	0.41
1:F:77:TRP:O	2:F:301:AVU:H5RA	2.21	0.40
1:E:232:ARG:NH1	1:F:250:LEU:HD12	2.36	0.40
1:A:157:TYR:CE2	1:A:159:VAL:CG1	3.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:PRO:HB2	1:C:126:PRO:HG2	2.02	0.40
1:F:50:PRO:HB2	1:F:126:PRO:HG3	2.03	0.40
1:H:16:LYS:HB2	1:H:16:LYS:HE2	1.76	0.40
1:D:205:LYS:O	1:D:211:LEU:HD12	2.21	0.40
1:G:81:TYR:CE2	1:G:85:HIS:CE1	3.10	0.40
1:H:28:PRO:HG2	1:H:67:GLN:HB3	2.03	0.40
1:G:1:ILE:CG2	1:G:125:CYS:HB2	2.51	0.40
1:F:192:VAL:O	1:F:193:ILE:HD13	2.22	0.40
1:A:55:LEU:O	1:A:142:MET:HG2	2.21	0.40
1:D:175:PHE:HA	1:D:179:GLU:HB2	2.03	0.40
1:C:101:LEU:O	1:C:105:MET:HG3	2.21	0.40
1:D:59:LYS:HA	1:D:142:MET:HE2	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	251/260 (96%)	238 (95%)	11 (4%)	2 (1%)	24	27
1	B	250/260 (96%)	236 (94%)	13 (5%)	1 (0%)	39	48
1	C	250/260 (96%)	234 (94%)	15 (6%)	1 (0%)	39	48
1	D	250/260 (96%)	237 (95%)	12 (5%)	1 (0%)	39	48
1	E	250/260 (96%)	237 (95%)	13 (5%)	0	100	100
1	F	250/260 (96%)	230 (92%)	14 (6%)	6 (2%)	7	5
1	G	250/260 (96%)	245 (98%)	4 (2%)	1 (0%)	39	48
1	H	250/260 (96%)	233 (93%)	14 (6%)	3 (1%)	16	16
All	All	2001/2080 (96%)	1890 (94%)	96 (5%)	15 (1%)	26	31

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	251	ALA
1	F	55	LEU
1	A	251	ALA
1	C	129	LYS
1	F	129	LYS
1	G	129	LYS
1	H	185	ASN
1	A	133	VAL
1	D	185	ASN
1	F	56	GLY
1	F	107	ASN
1	F	132	PRO
1	H	132	PRO
1	F	26	ILE
1	H	218	VAL

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	220/226 (97%)	203 (92%)	17 (8%)	16	20
1	B	220/226 (97%)	205 (93%)	15 (7%)	20	25
1	C	220/226 (97%)	212 (96%)	8 (4%)	42	57
1	D	220/226 (97%)	211 (96%)	9 (4%)	37	50
1	E	220/226 (97%)	211 (96%)	9 (4%)	37	50
1	F	220/226 (97%)	203 (92%)	17 (8%)	16	20
1	G	220/226 (97%)	203 (92%)	17 (8%)	16	20
1	H	220/226 (97%)	199 (90%)	21 (10%)	11	12
All	All	1760/1808 (97%)	1647 (94%)	113 (6%)	22	28

All (113) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	ILE

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Mol	Chain	Res	Type
1	A	29	ARG
1	A	35	SER
1	A	46	SER
1	A	68	GLN
1	A	73	LYS
1	A	86	ASP
1	A	92	ARG
1	A	97	LEU
1	A	98	GLU
1	A	115	ARG
1	A	123	LYS
1	A	124	VAL
1	A	129	LYS
1	A	138	SER
1	A	142	MET
1	A	249	ARG
1	B	5	ARG
1	B	20	ILE
1	B	25	ASP
1	B	31	ARG
1	B	33	ASP
1	B	68	GLN
1	B	89	ASN
1	B	136	ARG
1	B	172	ASP
1	B	202	ILE
1	B	232	ARG
1	B	240	SER
1	B	242	ASN
1	B	246	ARG
1	B	249	ARG
1	C	31	ARG
1	C	52	ASP
1	C	129	LYS
1	C	130	THR
1	C	133	VAL
1	C	195	LEU
1	C	212	LEU
1	C	217	LEU
1	D	25	ASP
1	D	33	ASP
1	D	75	MET

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Mol	Chain	Res	Type
1	D	89	ASN
1	D	115	ARG
1	D	136	ARG
1	D	185	ASN
1	D	232	ARG
1	D	246	ARG
1	E	1	ILE
1	E	31	ARG
1	E	43	LYS
1	E	52	ASP
1	E	150	SER
1	E	172	ASP
1	E	228	VAL
1	E	237	LEU
1	E	246	ARG
1	F	1	ILE
1	F	25	ASP
1	F	32	SER
1	F	39	LYS
1	F	49	ASN
1	F	53	LEU
1	F	66	GLN
1	F	68	GLN
1	F	89	ASN
1	F	117	ASN
1	F	122	GLU
1	F	124	VAL
1	F	129	LYS
1	F	134	GLN
1	F	138	SER
1	F	195	LEU
1	F	249	ARG
1	G	31	ARG
1	G	40	ASP
1	G	71	LYS
1	G	86	ASP
1	G	128	PHE
1	G	130	THR
1	G	136	ARG
1	G	142	MET
1	G	162	SER
1	G	195	LEU

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Mol	Chain	Res	Type
1	G	200	GLU
1	G	211	LEU
1	G	228	VAL
1	G	232	ARG
1	G	237	LEU
1	G	242	ASN
1	G	249	ARG
1	H	1	ILE
1	H	25	ASP
1	H	27	LEU
1	H	29	ARG
1	H	33	ASP
1	H	49	ASN
1	H	54	ASP
1	H	55	LEU
1	H	67	GLN
1	H	89	ASN
1	H	90	THR
1	H	115	ARG
1	H	123	LYS
1	H	127	ASP
1	H	129	LYS
1	H	130	THR
1	H	150	SER
1	H	172	ASP
1	H	195	LEU
1	H	200	GLU
1	H	242	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	GLN
1	A	107	ASN
1	B	68	GLN
1	B	89	ASN
1	B	242	ASN
1	C	49	ASN
1	C	89	ASN
1	D	49	ASN
1	D	68	GLN
1	D	89	ASN

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Mol	Chain	Res	Type
1	D	185	ASN
1	D	242	ASN
1	E	89	ASN
1	F	89	ASN
1	F	114	GLN
1	F	117	ASN
1	F	134	GLN
1	G	89	ASN
1	G	242	ASN
1	H	68	GLN
1	H	89	ASN
1	H	242	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	AVU	A	301	1	31,38,38	0.85	0	38,58,58	2.10	6 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	AVU	B	301	1	31,38,38	1.03	2 (6%)	38,58,58	2.25	11 (28%)
2	AVU	C	301	1	31,38,38	0.85	1 (3%)	38,58,58	2.04	5 (13%)
2	AVU	D	301	1	31,38,38	0.88	1 (3%)	38,58,58	2.49	7 (18%)
2	AVU	E	301	1	31,38,38	1.00	1 (3%)	38,58,58	2.64	7 (18%)
2	AVU	F	301	1	31,38,38	0.96	1 (3%)	38,58,58	2.42	10 (26%)
2	AVU	G	301	1	31,38,38	0.94	2 (6%)	38,58,58	2.41	8 (21%)
2	AVU	H	301	1	31,38,38	0.94	1 (3%)	38,58,58	2.29	8 (21%)

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	AVU	A	301	1	-	0/18/51/51	0/4/4/4
2	AVU	B	301	1	-	0/18/51/51	0/4/4/4
2	AVU	C	301	1	-	0/18/51/51	0/4/4/4
2	AVU	D	301	1	-	0/18/51/51	0/4/4/4
2	AVU	E	301	1	-	0/18/51/51	0/4/4/4
2	AVU	F	301	1	-	0/18/51/51	0/4/4/4
2	AVU	G	301	1	-	0/18/51/51	0/4/4/4
2	AVU	H	301	1	-	0/18/51/51	0/4/4/4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	301	AVU	O4R-C4R	-2.08	1.40	1.44
2	B	301	AVU	PB-O5R	2.42	1.70	1.59
2	G	301	AVU	O4'-C1'	2.46	1.44	1.41
2	F	301	AVU	O4'-C1'	2.66	1.44	1.41
2	H	301	AVU	O4'-C1'	2.67	1.44	1.41
2	C	301	AVU	O4'-C1'	2.93	1.44	1.41
2	D	301	AVU	O4'-C1'	3.05	1.45	1.41
2	B	301	AVU	O4'-C1'	3.40	1.45	1.41
2	E	301	AVU	O4'-C1'	3.75	1.45	1.41

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	301	AVU	N3-C2-N1	-11.60	120.01	128.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	AVU	N3-C2-N1	-11.37	120.19	128.89
2	F	301	AVU	N3-C2-N1	-10.44	120.90	128.89
2	G	301	AVU	N3-C2-N1	-10.40	120.93	128.89
2	C	301	AVU	N3-C2-N1	-10.38	120.95	128.89
2	B	301	AVU	N3-C2-N1	-9.98	121.25	128.89
2	H	301	AVU	N3-C2-N1	-9.20	121.85	128.89
2	A	301	AVU	N3-C2-N1	-8.94	122.05	128.89
2	E	301	AVU	F2R-C2R-C1R	-7.33	97.39	108.89
2	F	301	AVU	F2R-C2R-C3R	-6.37	104.62	108.81
2	E	301	AVU	F2R-C2R-C3R	-5.79	105.01	108.81
2	G	301	AVU	C2'-C1'-N9	-4.84	106.89	114.29
2	D	301	AVU	F2R-C2R-C3R	-4.69	105.72	108.81
2	A	301	AVU	C1'-N9-C4	-4.61	119.99	126.94
2	D	301	AVU	F2R-C2R-C1R	-4.01	102.60	108.89
2	G	301	AVU	O3A-PB-O5R	-3.99	92.35	102.94
2	A	301	AVU	C4-C5-N7	-3.82	105.97	109.48
2	F	301	AVU	C1'-N9-C4	-3.78	121.23	126.94
2	H	301	AVU	PB-O3A-PA	-3.72	122.28	132.73
2	H	301	AVU	O3R-C3R-C2R	-3.65	99.25	111.71
2	G	301	AVU	PB-O3A-PA	-3.56	122.74	132.73
2	H	301	AVU	C1'-N9-C4	-3.34	121.89	126.94
2	D	301	AVU	PB-O3A-PA	-3.19	123.76	132.73
2	B	301	AVU	C1'-N9-C4	-3.19	122.13	126.94
2	F	301	AVU	F2R-C2R-C1R	-3.17	103.91	108.89
2	A	301	AVU	PB-O3A-PA	-3.17	123.83	132.73
2	B	301	AVU	C4-C5-N7	-3.05	106.68	109.48
2	D	301	AVU	O3A-PB-O5R	-3.04	94.88	102.94
2	H	301	AVU	C4-C5-N7	-2.83	106.88	109.48
2	C	301	AVU	O3R-C3R-C2R	-2.73	102.39	111.71
2	A	301	AVU	O3R-C3R-C2R	-2.71	102.44	111.71
2	B	301	AVU	O3A-PB-O5R	-2.55	96.16	102.94
2	B	301	AVU	F2R-C2R-C1R	-2.55	104.89	108.89
2	G	301	AVU	F2R-C2R-C1R	-2.51	104.95	108.89
2	D	301	AVU	C2'-C1'-N9	-2.42	110.60	114.29
2	F	301	AVU	O5'-C5'-C4'	-2.40	100.28	109.12
2	E	301	AVU	C4-C5-N7	-2.30	107.36	109.48
2	C	301	AVU	C2'-C1'-N9	-2.24	110.86	114.29
2	F	301	AVU	C4-C5-N7	-2.20	107.46	109.48
2	B	301	AVU	C2'-C1'-N9	-2.18	110.96	114.29
2	B	301	AVU	PB-O3A-PA	-2.18	126.61	132.73
2	F	301	AVU	C1R-C2R-C3R	-2.14	101.70	104.12
2	F	301	AVU	O2'-C2'-C3'	-2.04	105.20	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	301	AVU	C2'-C1'-N9	-2.01	111.22	114.29
2	F	301	AVU	O4R-C1R-C2R	-2.00	101.82	104.80
2	E	301	AVU	C4'-O4'-C1'	2.01	111.93	109.72
2	B	301	AVU	C1R-O4R-C4R	2.05	113.22	108.08
2	E	301	AVU	C1R-O4R-C4R	2.06	113.26	108.08
2	G	301	AVU	O2A-PA-O1A	2.12	124.02	112.53
2	C	301	AVU	O3A-PA-O5'	2.13	108.60	102.94
2	D	301	AVU	O2A-PA-O1A	2.16	124.25	112.53
2	E	301	AVU	C2'-C3'-C4'	2.16	107.06	102.61
2	G	301	AVU	O5R-C5R-C4R	2.17	117.12	109.12
2	F	301	AVU	O4'-C4'-C5'	2.20	117.19	109.32
2	B	301	AVU	O4R-C1R-C2R	2.29	108.21	104.80
2	G	301	AVU	C1R-O4R-C4R	2.41	114.14	108.08
2	C	301	AVU	O4'-C1'-N9	2.51	113.34	108.10
2	H	301	AVU	F2R-C2R-C3R	3.00	110.77	108.81
2	B	301	AVU	O4'-C1'-N9	3.06	114.50	108.10
2	B	301	AVU	O5R-C5R-C4R	3.07	120.43	109.12
2	A	301	AVU	F2R-C2R-C3R	3.10	110.84	108.81
2	H	301	AVU	O4'-C1'-N9	5.19	118.96	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	AVU	2	0
2	B	301	AVU	6	0
2	C	301	AVU	2	0
2	D	301	AVU	5	0
2	E	301	AVU	3	0
2	F	301	AVU	5	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	253/260 (97%)	0.18	13 (5%) 32 41	29, 47, 80, 95	0
1	B	252/260 (96%)	-0.10	2 (0%) 87 90	27, 41, 61, 76	0
1	C	252/260 (96%)	0.10	9 (3%) 46 55	33, 51, 80, 96	0
1	D	252/260 (96%)	-0.10	4 (1%) 74 80	27, 41, 59, 67	0
1	E	252/260 (96%)	-0.07	2 (0%) 87 90	28, 42, 64, 73	0
1	F	252/260 (96%)	0.39	20 (7%) 15 22	42, 67, 98, 107	0
1	G	252/260 (96%)	-0.06	3 (1%) 81 85	28, 44, 65, 75	0
1	H	252/260 (96%)	0.31	16 (6%) 23 31	36, 57, 82, 93	0
All	All	2017/2080 (96%)	0.08	69 (3%) 49 58	27, 48, 81, 107	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	128	PHE	8.1
1	C	128	PHE	7.0
1	H	128	PHE	6.7
1	F	133	VAL	5.1
1	A	1	ILE	4.3
1	F	174	PHE	4.3
1	A	129	LYS	4.2
1	C	130	THR	4.0
1	C	127	ASP	3.9
1	F	51	CYS	3.9
1	F	134	GLN	3.8
1	G	251	ALA	3.6
1	F	124	VAL	3.5
1	A	130	THR	3.4
1	C	1	ILE	3.4
1	A	115	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	115	ARG	3.3
1	H	127	ASP	3.3
1	F	188	THR	3.2
1	H	77	TRP	3.2
1	A	2	VAL	3.2
1	F	43	LYS	3.2
1	C	52	ASP	3.2
1	H	130	THR	3.2
1	A	131	CYS	3.1
1	B	250	LEU	3.1
1	F	39	LYS	3.0
1	H	1	ILE	2.9
1	E	251	ALA	2.8
1	F	128	PHE	2.8
1	D	-1	ALA	2.8
1	A	49	ASN	2.7
1	A	51	CYS	2.7
1	H	126	PRO	2.7
1	C	115	ARG	2.6
1	F	25	ASP	2.6
1	C	174	PHE	2.6
1	H	246	ARG	2.6
1	A	127	ASP	2.5
1	F	115	ARG	2.5
1	A	246	ARG	2.5
1	F	130	THR	2.4
1	F	208	ALA	2.4
1	H	97	LEU	2.4
1	G	157	TYR	2.4
1	F	1	ILE	2.4
1	H	55	LEU	2.4
1	F	212	LEU	2.3
1	H	2	VAL	2.3
1	H	106	LEU	2.3
1	H	222	HIS	2.3
1	F	127	ASP	2.3
1	F	117	ASN	2.3
1	E	159	VAL	2.2
1	C	131	CYS	2.2
1	H	123	LYS	2.2
1	C	129	LYS	2.1
1	A	166	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	174	PHE	2.1
1	A	136	ARG	2.1
1	F	2	VAL	2.1
1	G	115	ARG	2.1
1	H	174	PHE	2.1
1	F	77	TRP	2.1
1	F	135	ALA	2.1
1	H	76	PHE	2.1
1	D	97	LEU	2.0
1	H	159	VAL	2.0
1	D	76	PHE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	AVU	E	301	35/35	0.94	0.17	0.68	36,66,109,110	0
2	AVU	B	301	35/35	0.91	0.16	0.59	34,63,102,103	0
2	AVU	G	301	35/35	0.94	0.16	0.33	39,64,102,102	0
2	AVU	H	301	35/35	0.95	0.20	0.30	55,64,77,79	0
2	AVU	A	301	35/35	0.91	0.15	0.27	50,74,83,85	0
2	AVU	C	301	35/35	0.92	0.14	-0.12	54,77,95,96	0
2	AVU	F	301	35/35	0.96	0.17	-0.25	58,70,85,87	0
2	AVU	D	301	35/35	0.94	0.14	-0.45	38,66,101,101	0

6.5 Other polymers

There are no such residues in this entry.