



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 09:42 PM GMT

PDB ID : 1R15
Title : Aplysia ADP ribosyl cyclase with bound nicotinamide and R5P
Authors : Love, M.L.; Szebenyi, D.M.E.; Kriksunov, I.A.; Thiel, D.J.; Munshi, C.; Graeff, R.; Lee, H.C.; Hao, Q.
Deposited on : 2003-09-23
Resolution : 2.40 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>

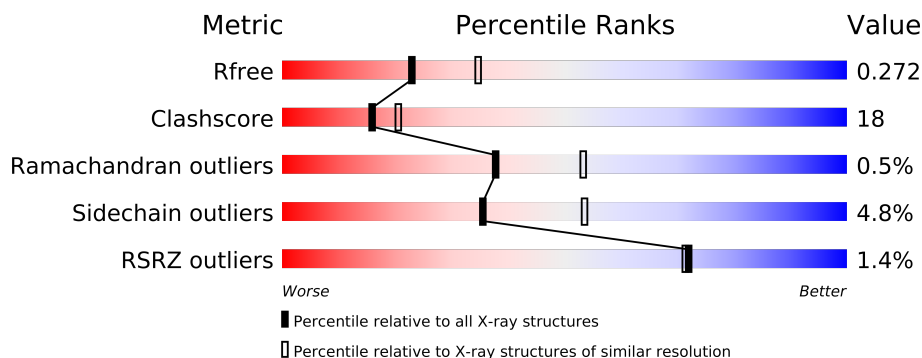
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
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 2.40 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}	66092	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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.

Mol	Chain	Length	Quality of chain
1	A	258	
1	B	258	
1	C	258	
1	D	258	
1	E	258	
1	F	258	
1	G	258	
1	H	258	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	N	A	1179	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
2	N	B	1179	-	X
2	N	C	1179	-	X
2	N	E	1179	-	X
2	N	F	1179	-	X
2	N	G	1179	-	X
2	N	H	1179	-	X
3	NCA	A	319	-	X
3	NCA	A	419	-	X
3	NCA	B	329	-	X
3	NCA	B	429	-	X
3	NCA	C	339	-	X
3	NCA	C	439	-	X
3	NCA	D	349	-	X
3	NCA	D	449	-	X
3	NCA	E	359	-	X
3	NCA	E	459	-	X
3	NCA	F	369	-	X
3	NCA	F	469	X	X
3	NCA	G	379	-	X
3	NCA	G	479	X	X
3	NCA	H	389	-	X
3	NCA	H	489	-	X

2 Entry composition

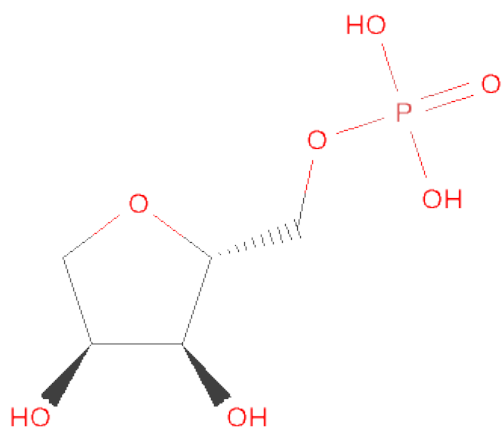
There are 3 unique types of molecules in this entry. The entry contains 16344 atoms, of which 0 are hydrogen and 0 are deuterium.

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 ADP-ribosyl cyclase.

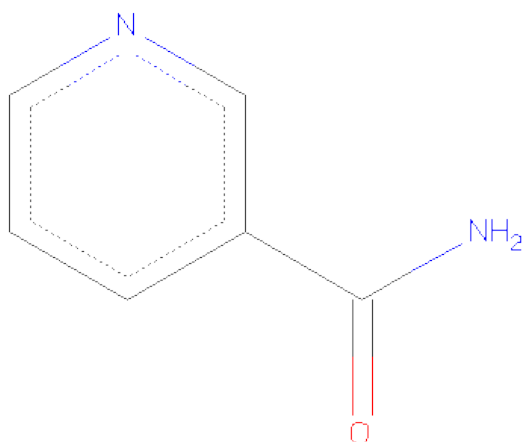
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total 2012	C 1288	N 342	O 368	S 14	0	0	0
1	B	251	Total 2012	C 1288	N 342	O 368	S 14	0	0	0
1	C	251	Total 2012	C 1288	N 342	O 368	S 14	0	0	0
1	D	251	Total 2012	C 1288	N 342	O 368	S 14	0	0	0
1	E	251	Total 2012	C 1288	N 342	O 368	S 14	0	0	0
1	F	251	Total 2012	C 1288	N 342	O 368	S 14	0	0	0
1	G	251	Total 2012	C 1288	N 342	O 368	S 14	0	0	0
1	H	251	Total 2012	C 1288	N 342	O 368	S 14	0	0	0

- Molecule 2 is ANY 5'-MONOPHOSPHATE NUCLEOTIDE (three-letter code: N) (formula: C₅H₁₁O₇P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			13	5	7	1		
2	B	1	Total	C	O	P	0	0
			13	5	7	1		
2	C	1	Total	C	O	P	0	0
			13	5	7	1		
2	D	1	Total	C	O	P	0	0
			13	5	7	1		
2	E	1	Total	C	O	P	0	0
			13	5	7	1		
2	F	1	Total	C	O	P	0	0
			13	5	7	1		
2	G	1	Total	C	O	P	0	0
			13	5	7	1		
2	H	1	Total	C	O	P	0	0
			13	5	7	1		

- Molecule 3 is NICOTINAMIDE (three-letter code: NCA) (formula: C₆H₆N₂O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	6	2	1		
3	B	1	Total	C	N	O	0	0
			9	6	2	1		
3	C	1	Total	C	N	O	0	0
			9	6	2	1		
3	D	1	Total	C	N	O	0	0
			9	6	2	1		
3	E	1	Total	C	N	O	0	0
			9	6	2	1		
3	F	1	Total	C	N	O	0	0
			9	6	2	1		
3	G	1	Total	C	N	O	0	0
			9	6	2	1		
3	H	1	Total	C	N	O	0	0
			9	6	2	1		
3	A	1	Total	C	N	O	0	0
			9	6	2	1		
3	B	1	Total	C	N	O	0	0
			9	6	2	1		
3	C	1	Total	C	N	O	0	0
			9	6	2	1		
3	D	1	Total	C	N	O	0	0
			9	6	2	1		
3	E	1	Total	C	N	O	0	0
			9	6	2	1		
3	F	1	Total	C	N	O	0	0
			9	6	2	1		

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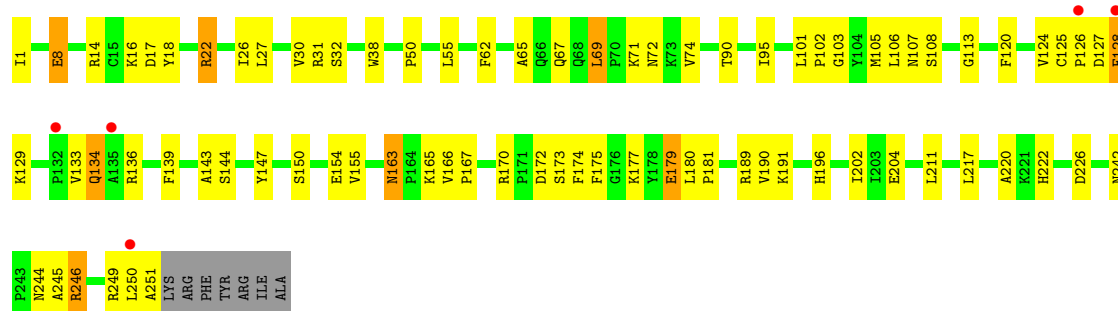
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	G	1	Total	C	N	O	0	0
			9	6	2	1		
3	H	1	Total	C	N	O	0	0
			9	6	2	1		

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.

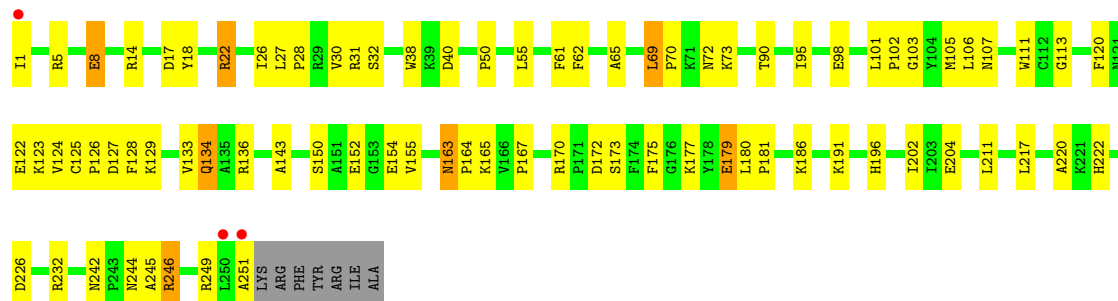
• Molecule 1: ADP-ribosyl cyclase

Chain A: 



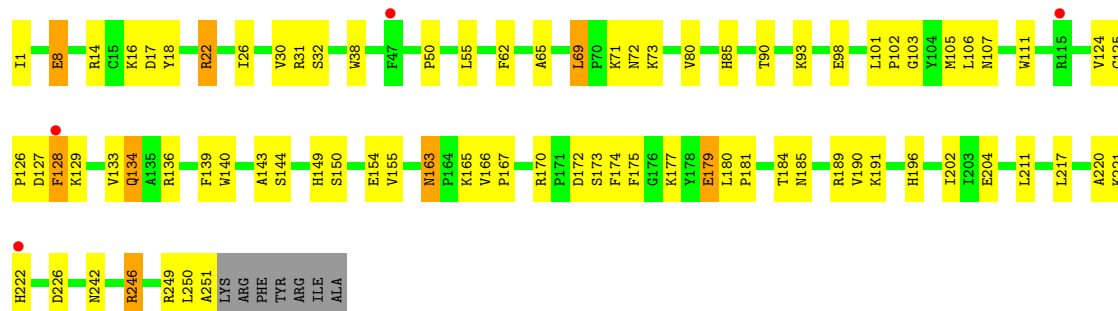
• Molecule 1: ADP-ribosyl cyclase

Chain B: 

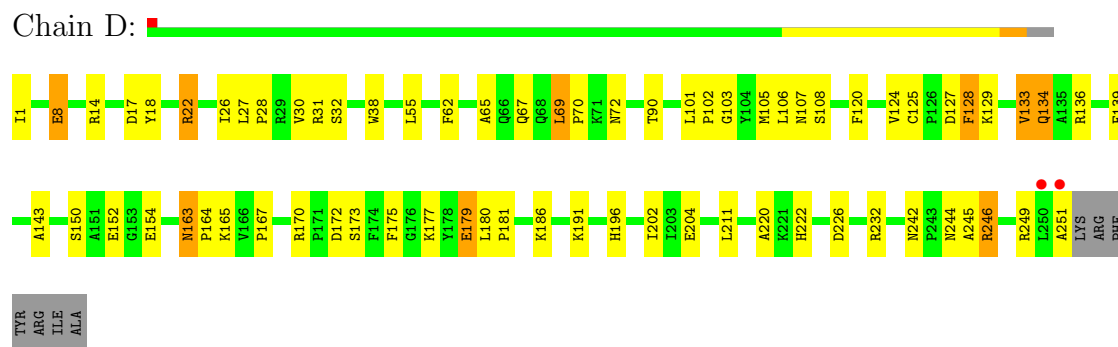


• Molecule 1: ADP-ribosyl cyclase

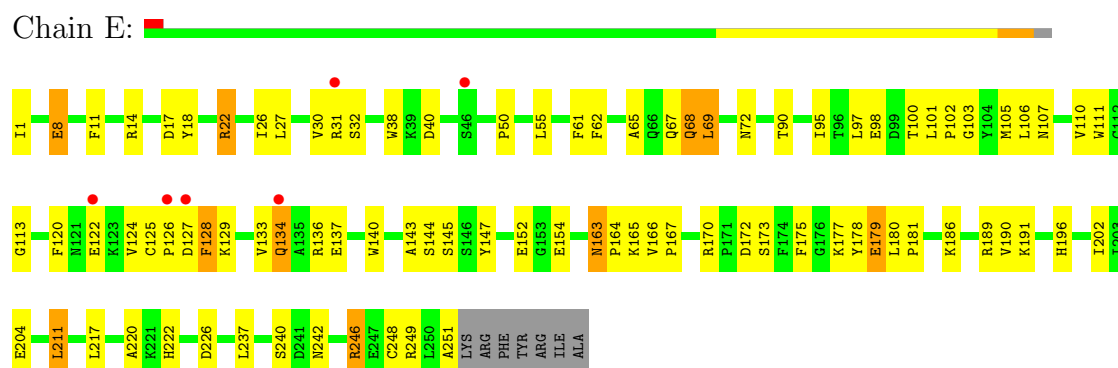
Chain C: 



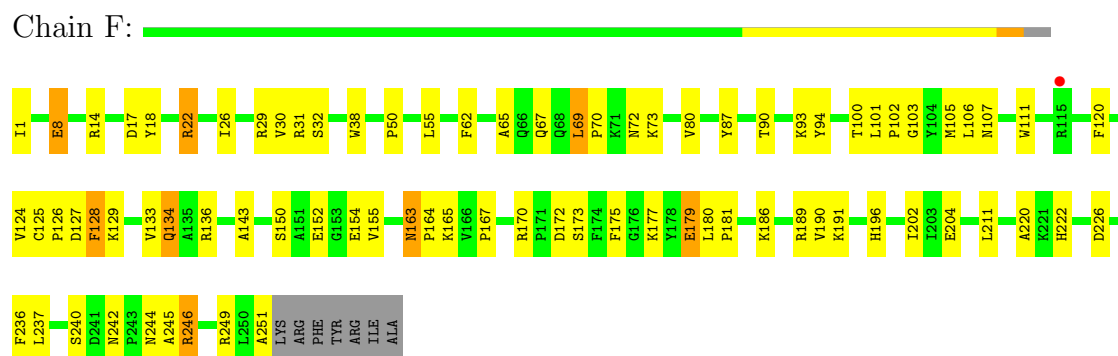
- Molecule 1: ADP-ribosyl cyclase



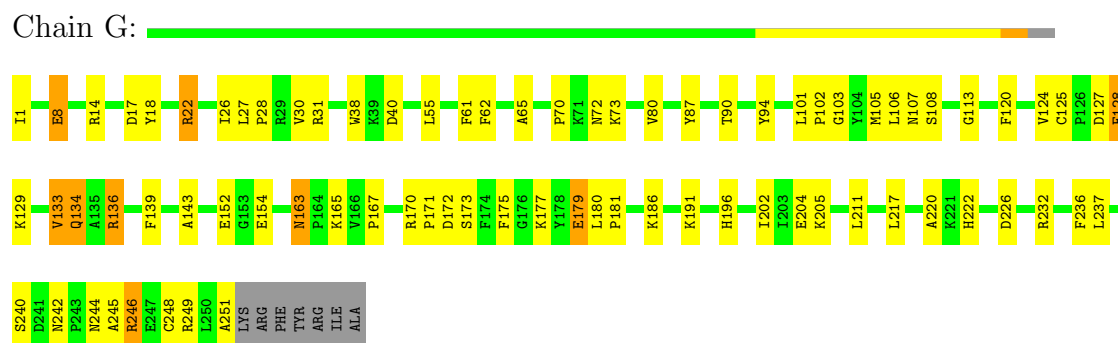
- Molecule 1: ADP-ribosyl cyclase



- Molecule 1: ADP-ribosyl cyclase

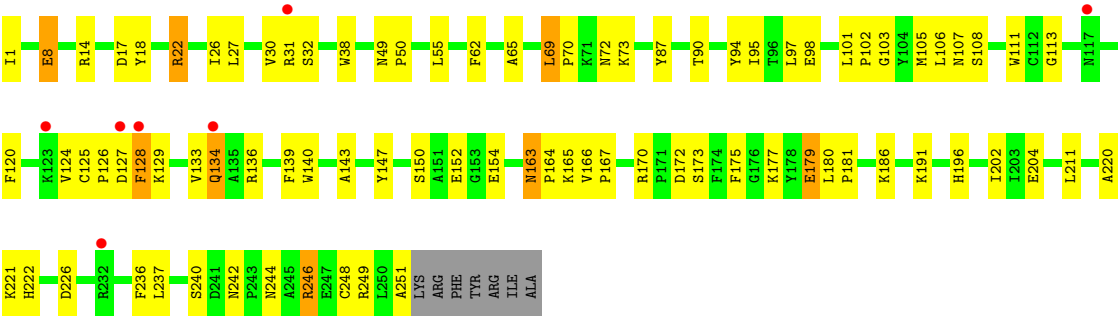


- Molecule 1: ADP-ribosyl cyclase



- Molecule 1: ADP-ribosyl cyclase

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.40Å 75.32Å 138.13Å 88.16° 89.22° 89.09°	Depositor
Resolution (Å)	33.00 – 2.40 32.95 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (33.00-2.40) 60.1 (32.95-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.244 , 0.269 0.246 , 0.272	Depositor DCC
R_{free} test set	2870 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 12.8	EDS
Estimated twinning fraction	0.085 for h,-k,-l 0.022 for -h,k,-l 0.008 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 57205 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	16344	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.84	1/2064 (0.0%)	0.77	2/2793 (0.1%)
1	B	0.85	2/2064 (0.1%)	0.79	2/2793 (0.1%)
1	C	0.84	1/2064 (0.0%)	0.77	2/2793 (0.1%)
1	D	0.84	1/2064 (0.0%)	0.80	2/2793 (0.1%)
1	E	0.82	1/2064 (0.0%)	0.78	2/2793 (0.1%)
1	F	0.86	2/2064 (0.1%)	0.79	2/2793 (0.1%)
1	G	0.87	2/2064 (0.1%)	1.15	6/2793 (0.2%)
1	H	0.80	1/2064 (0.0%)	0.77	2/2793 (0.1%)
All	All	0.84	11/16512 (0.1%)	0.84	20/22344 (0.1%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	179	GLU	CD-OE2	23.53	1.51	1.25
1	E	179	GLU	CD-OE2	23.07	1.51	1.25
1	C	179	GLU	CD-OE2	23.00	1.50	1.25
1	H	179	GLU	CD-OE2	22.84	1.50	1.25
1	G	179	GLU	CD-OE2	22.59	1.50	1.25
1	F	179	GLU	CD-OE2	22.37	1.50	1.25
1	D	179	GLU	CD-OE2	21.57	1.49	1.25
1	B	179	GLU	CD-OE2	21.44	1.49	1.25
1	G	136	ARG	CD-NE	-5.83	1.36	1.46
1	F	179	GLU	CB-CG	5.71	1.63	1.52
1	B	179	GLU	CB-CG	5.47	1.62	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	136	ARG	NE-CZ-NH1	-29.87	105.36	120.30
1	G	136	ARG	NE-CZ-NH2	29.33	134.96	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	136	ARG	CD-NE-CZ	13.46	142.44	123.60
1	G	136	ARG	CG-CD-NE	8.30	129.23	111.80
1	B	22	ARG	N-CA-C	6.93	129.72	111.00
1	D	22	ARG	N-CA-C	6.84	129.47	111.00
1	E	22	ARG	N-CA-C	6.68	129.03	111.00
1	H	22	ARG	N-CA-C	6.66	128.99	111.00
1	A	22	ARG	N-CA-C	6.57	128.75	111.00
1	G	22	ARG	N-CA-C	6.54	128.67	111.00
1	C	22	ARG	N-CA-C	6.52	128.59	111.00
1	F	22	ARG	N-CA-C	6.41	128.31	111.00
1	B	173	SER	N-CA-C	-5.65	95.75	111.00
1	E	173	SER	N-CA-C	-5.50	96.16	111.00
1	C	173	SER	N-CA-C	-5.36	96.53	111.00
1	A	173	SER	N-CA-C	-5.33	96.60	111.00
1	F	173	SER	N-CA-C	-5.29	96.73	111.00
1	D	173	SER	N-CA-C	-5.28	96.75	111.00
1	H	173	SER	N-CA-C	-5.19	96.98	111.00
1	G	173	SER	N-CA-C	-5.01	97.48	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	A	2012	0	1968	80	0
1	B	2012	0	1968	82	0
1	C	2012	0	1968	86	0
1	D	2012	0	1968	79	0
1	E	2012	0	1968	82	1
1	F	2012	0	1968	68	0
1	G	2012	0	1968	83	0
1	H	2012	0	1968	84	1
2	A	13	0	8	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	13	0	8	0	0
2	C	13	0	8	2	0
2	D	13	0	8	0	0
2	E	13	0	8	1	0
2	F	13	0	8	1	0
2	G	13	0	8	3	0
2	H	13	0	8	0	0
3	A	18	0	12	3	0
3	B	18	0	12	3	0
3	C	18	0	12	4	0
3	D	18	0	12	2	0
3	E	18	0	12	7	0
3	F	18	0	12	3	0
3	G	18	0	12	5	0
3	H	18	0	12	7	0
All	All	16344	0	15904	592	1

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 18.

All (592) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:123:LYS:HD2	1:C:149:HIS:CE1	1.86	1.10
1:B:123:LYS:HD2	1:C:149:HIS:ND1	1.76	0.99
1:D:152:GLU:HG2	1:D:186:LYS:HB3	1.40	0.97
1:A:139:PHE:HB3	3:A:419:NCA:H6	1.42	0.97
1:G:152:GLU:HG2	1:G:186:LYS:HB3	1.46	0.96
1:B:111:TRP:H	3:B:429:NCA:HN72	1.11	0.96
1:B:123:LYS:HG2	1:C:184:THR:HG22	1.50	0.93
1:A:17:ASP:OD1	1:B:14:ARG:HD2	1.70	0.91
1:B:123:LYS:CD	1:C:149:HIS:CE1	2.54	0.90
1:H:140:TRP:CD1	3:H:489:NCA:H5	2.06	0.90
1:G:17:ASP:OD1	1:H:14:ARG:HD2	1.73	0.88
1:D:133:VAL:HG22	1:G:177:LYS:HB3	1.56	0.87
2:F:1179:N:H1'	3:F:369:NCA:H6	1.54	0.87
1:C:69:LEU:CD2	1:C:150:SER:HB2	2.06	0.86
1:D:69:LEU:HD22	1:D:150:SER:HB2	1.57	0.86
1:H:108:SER:H	3:H:389:NCA:HN72	1.23	0.85
1:F:69:LEU:HD22	1:F:150:SER:HB2	1.58	0.85
1:H:152:GLU:HG2	1:H:186:LYS:HB3	1.57	0.85
1:E:128:PHE:HB3	1:E:129:LYS:HD2	1.58	0.85
1:E:17:ASP:OD1	1:F:14:ARG:HD2	1.77	0.84

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:128:PHE:HB3	1:A:129:LYS:HD2	1.60	0.84
1:H:128:PHE:HB3	1:H:129:LYS:HD2	1.59	0.83
1:G:14:ARG:HD2	1:H:17:ASP:OD1	1.79	0.83
1:E:129:LYS:CD	1:E:129:LYS:H	1.91	0.83
1:A:128:PHE:CE2	1:A:136:ARG:HD3	2.12	0.83
1:G:128:PHE:HB3	1:G:129:LYS:HD2	1.62	0.82
1:C:17:ASP:OD1	1:D:14:ARG:HD2	1.80	0.82
1:E:128:PHE:CE2	1:E:136:ARG:HD3	2.14	0.82
1:E:129:LYS:H	1:E:129:LYS:HD2	1.45	0.81
1:C:128:PHE:HB3	1:C:129:LYS:HD2	1.62	0.81
1:A:129:LYS:H	1:A:129:LYS:CD	1.94	0.81
1:F:128:PHE:HB3	1:F:129:LYS:HD2	1.63	0.81
1:H:129:LYS:H	1:H:129:LYS:HD2	1.46	0.80
1:H:129:LYS:H	1:H:129:LYS:CD	1.94	0.80
1:C:128:PHE:CE2	1:C:136:ARG:HD3	2.15	0.79
1:C:129:LYS:CD	1:C:129:LYS:H	1.96	0.79
1:F:128:PHE:CE2	1:F:136:ARG:HD3	2.18	0.79
1:B:128:PHE:HB3	1:B:129:LYS:HD2	1.65	0.79
1:B:123:LYS:CG	1:C:184:THR:HG22	2.13	0.78
1:H:128:PHE:CE2	1:H:136:ARG:HD3	2.17	0.78
1:G:129:LYS:CD	1:G:129:LYS:H	1.97	0.78
1:B:128:PHE:CE2	1:B:136:ARG:HD3	2.19	0.78
1:A:129:LYS:HD2	1:A:129:LYS:H	1.49	0.77
1:C:129:LYS:HD2	1:C:129:LYS:H	1.49	0.77
1:F:129:LYS:H	1:F:129:LYS:CD	1.97	0.77
1:E:14:ARG:HD2	1:F:17:ASP:OD1	1.85	0.77
1:C:71:LYS:HG2	1:C:72:ASN:ND2	2.00	0.77
1:B:5:ARG:HD2	1:C:185:ASN:O	1.85	0.77
1:D:133:VAL:HG13	1:G:177:LYS:HG2	1.67	0.77
1:D:133:VAL:CG2	1:G:177:LYS:HB3	2.14	0.76
1:B:129:LYS:H	1:B:129:LYS:CD	1.99	0.76
1:D:128:PHE:HB3	1:D:129:LYS:HD2	1.67	0.76
1:F:111:TRP:O	3:F:469:NCA:N7	2.19	0.76
1:A:167:PRO:HG2	1:A:170:ARG:HG2	1.68	0.76
1:B:123:LYS:CD	1:C:149:HIS:HE1	1.97	0.75
1:H:8:GLU:HG2	1:H:38:TRP:CZ2	2.22	0.75
1:B:111:TRP:N	3:B:429:NCA:HN72	1.84	0.75
1:H:69:LEU:H	1:H:69:LEU:HD13	1.52	0.74
1:E:167:PRO:HG2	1:E:170:ARG:HG2	1.70	0.74
1:F:129:LYS:HD2	1:F:129:LYS:H	1.53	0.74
1:D:128:PHE:CE2	1:D:136:ARG:HD3	2.23	0.73
1:G:129:LYS:HD2	1:G:129:LYS:H	1.52	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:129:LYS:HD2	1:B:129:LYS:H	1.53	0.73
1:C:167:PRO:HG2	1:C:170:ARG:HG2	1.70	0.72
1:H:167:PRO:HG2	1:H:170:ARG:HG2	1.71	0.72
1:E:163:ASN:HD22	1:E:163:ASN:C	1.91	0.72
1:A:69:LEU:CD2	1:A:150:SER:HB2	2.19	0.72
1:E:240:SER:OG	1:F:237:LEU:HD13	1.90	0.72
1:B:101:LEU:HB3	1:B:102:PRO:HD3	1.72	0.71
1:C:8:GLU:HG2	1:C:38:TRP:CZ2	2.25	0.71
1:A:8:GLU:HG2	1:A:38:TRP:CZ2	2.26	0.71
1:G:246:ARG:O	1:G:249:ARG:HG3	1.91	0.71
1:D:129:LYS:H	1:D:129:LYS:CD	2.04	0.71
1:C:101:LEU:HB3	1:C:102:PRO:HD3	1.72	0.70
1:E:8:GLU:HG2	1:E:38:TRP:CZ2	2.26	0.70
1:F:246:ARG:O	1:F:249:ARG:HG3	1.91	0.70
1:D:129:LYS:H	1:D:129:LYS:HD2	1.57	0.70
1:G:237:LEU:HD13	1:H:240:SER:OG	1.92	0.70
1:C:14:ARG:HD2	1:D:17:ASP:OD1	1.92	0.70
1:F:72:ASN:HA	1:F:154:GLU:O	1.92	0.70
1:E:69:LEU:HG	1:E:95:ILE:HD11	1.73	0.69
1:C:31:ARG:HG3	1:C:31:ARG:HH11	1.57	0.69
1:D:129:LYS:HG3	1:G:171:PRO:HG3	1.74	0.69
1:E:101:LEU:HB3	1:E:102:PRO:HD3	1.74	0.69
1:G:240:SER:OG	1:H:237:LEU:HD13	1.93	0.68
1:B:8:GLU:HG2	1:B:38:TRP:CZ2	2.28	0.68
1:D:129:LYS:HB3	1:G:217:LEU:HD21	1.75	0.68
1:B:167:PRO:HG2	1:B:170:ARG:HG2	1.74	0.68
1:F:8:GLU:HG2	1:F:38:TRP:CZ2	2.29	0.68
1:E:134:GLN:H	1:E:134:GLN:HE21	1.42	0.68
1:D:101:LEU:HB3	1:D:102:PRO:HD3	1.76	0.68
1:H:134:GLN:HE21	1:H:134:GLN:H	1.42	0.67
1:D:167:PRO:HG2	1:D:170:ARG:HG2	1.75	0.67
1:A:69:LEU:HD22	1:A:150:SER:HB2	1.76	0.67
1:A:31:ARG:HH11	1:A:31:ARG:HG3	1.59	0.67
1:H:70:PRO:HG2	1:H:73:LYS:HB2	1.77	0.67
1:E:237:LEU:HD13	1:F:240:SER:OG	1.95	0.67
1:A:31:ARG:NH2	1:C:191:LYS:NZ	2.42	0.67
1:F:101:LEU:HB3	1:F:102:PRO:HD3	1.76	0.67
1:E:111:TRP:O	3:E:459:NCA:N7	2.27	0.67
1:D:246:ARG:O	1:D:249:ARG:HG3	1.95	0.67
1:B:163:ASN:HD22	1:B:163:ASN:C	1.98	0.67
1:A:251:ALA:HA	1:B:251:ALA:HA	1.75	0.66
1:G:8:GLU:HG2	1:G:38:TRP:CZ2	2.30	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:8:GLU:HG2	1:D:38:TRP:CZ2	2.30	0.66
1:A:14:ARG:HD2	1:B:17:ASP:OD1	1.95	0.66
1:H:107:ASN:OD1	3:H:389:NCA:N7	2.27	0.66
1:G:134:GLN:H	1:G:134:GLN:HE21	1.43	0.66
1:H:101:LEU:HB3	1:H:102:PRO:HD3	1.77	0.66
1:H:31:ARG:HH11	1:H:31:ARG:HG3	1.60	0.66
1:G:191:LYS:HE3	1:G:226:ASP:OD2	1.96	0.65
1:A:65:ALA:HB1	1:A:102:PRO:HG3	1.79	0.65
1:B:127:ASP:OD1	1:B:129:LYS:HD3	1.97	0.65
1:C:111:TRP:H	3:C:439:NCA:HN72	1.44	0.65
1:B:191:LYS:HE3	1:B:226:ASP:OD2	1.96	0.65
1:A:101:LEU:HB3	1:A:102:PRO:HD3	1.78	0.65
1:F:191:LYS:HE3	1:F:226:ASP:OD2	1.96	0.65
1:H:163:ASN:HD22	1:H:163:ASN:C	1.99	0.65
1:B:65:ALA:HB1	1:B:102:PRO:HG3	1.79	0.65
1:A:163:ASN:HD22	1:A:163:ASN:C	1.98	0.65
1:E:31:ARG:HH11	1:E:31:ARG:HG3	1.61	0.65
1:C:163:ASN:C	1:C:163:ASN:HD22	1.99	0.65
1:A:69:LEU:HD22	1:A:150:SER:CB	2.27	0.65
1:F:167:PRO:HG2	1:F:170:ARG:HG2	1.79	0.65
1:D:163:ASN:C	1:D:163:ASN:HD22	2.00	0.65
1:B:246:ARG:O	1:B:249:ARG:HG3	1.97	0.64
1:C:31:ARG:HG3	1:C:31:ARG:NH1	2.12	0.64
1:F:134:GLN:H	1:F:134:GLN:HE21	1.43	0.64
1:F:134:GLN:H	1:F:134:GLN:NE2	1.96	0.64
1:G:167:PRO:HG2	1:G:170:ARG:HG2	1.79	0.64
1:B:69:LEU:HG	1:B:95:ILE:HD11	1.79	0.64
1:E:110:VAL:HA	3:E:459:NCA:O7	1.98	0.64
1:B:134:GLN:H	1:B:134:GLN:HE21	1.45	0.64
1:C:134:GLN:HE21	1:C:134:GLN:H	1.45	0.64
1:A:31:ARG:NH1	1:A:31:ARG:HG3	2.12	0.64
1:H:31:ARG:HG3	1:H:31:ARG:NH1	2.13	0.64
1:G:70:PRO:HG2	1:G:73:LYS:HB2	1.80	0.63
1:A:134:GLN:HE21	1:A:134:GLN:H	1.44	0.63
1:H:108:SER:N	3:H:389:NCA:HN72	1.95	0.63
1:G:101:LEU:HB3	1:G:102:PRO:HD3	1.79	0.63
1:E:140:TRP:CD1	3:E:459:NCA:H5	2.34	0.63
1:H:127:ASP:OD1	1:H:129:LYS:HD3	1.98	0.63
1:E:251:ALA:HA	1:F:251:ALA:HA	1.79	0.63
1:D:191:LYS:HE3	1:D:226:ASP:OD2	1.99	0.62
1:D:133:VAL:HG11	1:G:177:LYS:CD	2.30	0.62
1:B:26:ILE:N	1:B:26:ILE:HD12	2.13	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:152:GLU:HG2	1:G:186:LYS:CB	2.26	0.62
1:H:69:LEU:N	1:H:69:LEU:HD13	2.14	0.62
1:D:133:VAL:CG1	1:G:177:LYS:HG2	2.28	0.62
1:H:65:ALA:HB1	1:H:102:PRO:HG3	1.82	0.62
1:H:191:LYS:HE3	1:H:226:ASP:OD2	2.00	0.62
1:C:65:ALA:HB1	1:C:102:PRO:HG3	1.82	0.62
1:D:26:ILE:HD12	1:D:26:ILE:N	2.15	0.62
1:F:163:ASN:C	1:F:163:ASN:HD22	2.03	0.61
1:D:65:ALA:HB1	1:D:102:PRO:HG3	1.83	0.61
1:E:65:ALA:HB1	1:E:102:PRO:HG3	1.82	0.61
1:H:134:GLN:NE2	1:H:134:GLN:H	1.98	0.61
1:E:31:ARG:NH1	1:E:31:ARG:HG3	2.14	0.61
1:H:246:ARG:O	1:H:249:ARG:HG3	2.01	0.61
1:C:72:ASN:HA	1:C:154:GLU:O	2.01	0.60
1:C:127:ASP:OD1	1:C:129:LYS:HD3	2.01	0.60
1:B:152:GLU:HG2	1:B:186:LYS:HB3	1.83	0.60
1:C:172:ASP:HA	1:C:177:LYS:HG3	1.84	0.60
1:D:31:ARG:HH11	1:D:31:ARG:HG3	1.64	0.60
1:E:127:ASP:OD1	1:E:129:LYS:HD3	2.00	0.60
1:F:111:TRP:H	3:F:469:NCA:HN72	1.49	0.60
1:A:108:SER:H	3:A:319:NCA:HN72	1.48	0.60
1:F:31:ARG:HG3	1:F:31:ARG:HH11	1.66	0.60
1:D:134:GLN:HE21	1:D:134:GLN:H	1.49	0.60
1:A:246:ARG:O	1:A:249:ARG:HG3	2.01	0.60
1:E:134:GLN:H	1:E:134:GLN:NE2	1.99	0.60
1:G:134:GLN:NE2	1:G:134:GLN:H	2.00	0.60
1:A:31:ARG:NH2	1:C:191:LYS:HZ1	1.99	0.60
1:G:163:ASN:C	1:G:163:ASN:HD22	2.05	0.60
1:E:191:LYS:HE3	1:E:226:ASP:OD2	2.02	0.60
1:D:127:ASP:OD1	1:D:129:LYS:HD3	2.02	0.60
1:C:69:LEU:CD2	1:C:150:SER:CB	2.79	0.60
1:G:127:ASP:OD1	1:G:129:LYS:HD3	2.02	0.60
1:C:251:ALA:HA	1:D:251:ALA:HA	1.84	0.59
1:H:26:ILE:N	1:H:26:ILE:HD12	2.17	0.59
1:H:140:TRP:HD1	3:H:489:NCA:H5	1.66	0.59
1:G:65:ALA:HB1	1:G:102:PRO:HG3	1.83	0.59
1:B:55:LEU:HD22	1:B:55:LEU:H	1.68	0.59
1:A:191:LYS:HE3	1:A:226:ASP:OD2	2.02	0.59
1:F:26:ILE:HD12	1:F:26:ILE:N	2.17	0.59
1:B:134:GLN:H	1:B:134:GLN:NE2	1.99	0.59
1:E:1:ILE:HG22	1:E:125:CYS:O	2.02	0.59
1:E:129:LYS:N	1:E:129:LYS:HD2	2.17	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:172:ASP:HA	1:E:177:LYS:HG3	1.85	0.59
1:F:127:ASP:OD1	1:F:129:LYS:HD3	2.03	0.59
1:H:172:ASP:HA	1:H:177:LYS:HG3	1.85	0.59
1:G:31:ARG:HG3	1:G:31:ARG:HH11	1.68	0.59
1:A:71:LYS:O	1:A:72:ASN:HB2	2.03	0.59
1:B:122:GLU:O	1:C:185:ASN:OD1	2.20	0.59
1:B:31:ARG:HH11	1:B:31:ARG:HG3	1.66	0.58
1:A:127:ASP:OD1	1:A:129:LYS:HD3	2.03	0.58
1:H:111:TRP:O	3:H:489:NCA:N7	2.36	0.58
1:H:129:LYS:HD2	1:H:129:LYS:N	2.18	0.58
1:C:246:ARG:O	1:C:249:ARG:HG3	2.02	0.58
1:A:172:ASP:HA	1:A:177:LYS:HG3	1.86	0.58
1:C:128:PHE:CD2	1:C:136:ARG:HD3	2.38	0.58
1:B:55:LEU:HD22	1:B:55:LEU:N	2.19	0.58
2:C:1179:N:HI'	3:C:339:NCA:H6	1.85	0.58
1:D:177:LYS:HB3	1:G:133:VAL:HG22	1.86	0.58
1:B:72:ASN:HA	1:B:154:GLU:O	2.04	0.58
1:D:152:GLU:HG2	1:D:186:LYS:CB	2.26	0.57
1:A:134:GLN:NE2	1:A:134:GLN:H	2.02	0.57
1:D:196:HIS:CE1	1:D:202:ILE:HG23	2.39	0.57
1:D:134:GLN:NE2	1:D:134:GLN:H	2.02	0.57
1:B:31:ARG:NH1	1:B:31:ARG:HG3	2.19	0.57
1:E:26:ILE:N	1:E:26:ILE:HD12	2.19	0.57
1:D:129:LYS:CB	1:G:217:LEU:HD21	2.35	0.57
1:H:69:LEU:N	1:H:69:LEU:CD1	2.67	0.57
1:C:69:LEU:HD22	1:C:150:SER:CB	2.35	0.57
1:D:31:ARG:NH1	1:D:31:ARG:HG3	2.18	0.57
1:A:128:PHE:CD2	1:A:136:ARG:HD3	2.39	0.57
1:C:134:GLN:NE2	1:C:134:GLN:H	2.01	0.56
1:F:31:ARG:NH1	1:F:31:ARG:HG3	2.19	0.56
1:D:129:LYS:HG3	1:G:171:PRO:CG	2.35	0.56
1:G:108:SER:H	3:G:379:NCA:HN72	1.52	0.55
1:G:31:ARG:NH1	1:G:31:ARG:HG3	2.20	0.55
1:C:191:LYS:HE3	1:C:226:ASP:OD2	2.06	0.55
1:D:177:LYS:HB3	1:G:133:VAL:CG2	2.36	0.55
1:E:144:SER:HB3	1:E:179:GLU:HG3	1.88	0.55
1:B:101:LEU:HG	1:B:105:MET:HE1	1.88	0.55
1:H:175:PHE:HA	1:H:179:GLU:HB2	1.89	0.55
1:B:196:HIS:CE1	1:B:202:ILE:HG23	2.42	0.55
1:C:163:ASN:ND2	1:C:165:LYS:H	2.05	0.55
1:H:167:PRO:HG2	1:H:170:ARG:CG	2.37	0.54
1:E:27:LEU:HD21	1:E:95:ILE:HD13	1.89	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:196:HIS:CE1	1:E:202:ILE:HG23	2.41	0.54
1:E:175:PHE:HA	1:E:179:GLU:HB2	1.90	0.54
1:F:175:PHE:HA	1:F:179:GLU:HB2	1.88	0.54
1:E:69:LEU:N	1:E:69:LEU:CD1	2.70	0.54
1:E:152:GLU:HG2	1:E:186:LYS:HB3	1.90	0.54
1:D:55:LEU:HD22	1:D:55:LEU:N	2.22	0.54
1:G:26:ILE:HD12	1:G:26:ILE:N	2.22	0.54
1:A:26:ILE:HD12	1:A:26:ILE:N	2.22	0.54
1:E:67:GLN:OE1	1:E:100:THR:HB	2.08	0.54
1:C:26:ILE:HD12	1:C:26:ILE:N	2.22	0.54
1:H:55:LEU:N	1:H:55:LEU:HD22	2.23	0.54
1:E:246:ARG:O	1:E:249:ARG:HG3	2.08	0.54
1:E:55:LEU:N	1:E:55:LEU:HD22	2.23	0.54
1:B:62:PHE:CZ	1:B:143:ALA:HB2	2.43	0.53
1:C:129:LYS:HD2	1:C:129:LYS:N	2.20	0.53
1:F:65:ALA:HB1	1:F:102:PRO:HG3	1.89	0.53
1:D:55:LEU:H	1:D:55:LEU:HD22	1.74	0.53
1:D:72:ASN:HA	1:D:154:GLU:O	2.08	0.53
1:E:128:PHE:CD2	1:E:136:ARG:HD3	2.42	0.53
1:H:128:PHE:CD2	1:H:136:ARG:HD3	2.42	0.53
1:H:69:LEU:H	1:H:69:LEU:CD1	2.21	0.53
1:D:101:LEU:HG	1:D:105:MET:CE	2.38	0.53
1:G:251:ALA:HA	1:H:251:ALA:HA	1.88	0.53
1:C:18:TYR:HA	1:C:22:ARG:CG	2.39	0.53
1:B:172:ASP:HA	1:B:177:LYS:HG3	1.91	0.53
1:C:167:PRO:HG2	1:C:170:ARG:CG	2.38	0.53
1:A:55:LEU:HD22	1:A:55:LEU:N	2.23	0.53
1:F:172:ASP:HA	1:F:177:LYS:HG3	1.90	0.53
1:E:167:PRO:HG2	1:E:170:ARG:CG	2.38	0.53
1:B:163:ASN:ND2	1:B:165:LYS:H	2.07	0.53
1:D:175:PHE:HA	1:D:179:GLU:HB2	1.91	0.53
1:G:139:PHE:H	3:G:479:NCA:C6	2.22	0.53
1:A:167:PRO:HG2	1:A:170:ARG:CG	2.38	0.53
1:D:172:ASP:HA	1:D:177:LYS:HG3	1.91	0.53
1:C:18:TYR:HA	1:C:22:ARG:HG2	1.91	0.53
1:C:196:HIS:CE1	1:C:202:ILE:HG23	2.44	0.53
1:A:129:LYS:HD2	1:A:129:LYS:N	2.20	0.52
1:F:129:LYS:HD2	1:F:129:LYS:N	2.22	0.52
1:H:103:GLY:O	1:H:107:ASN:HB2	2.10	0.52
1:B:129:LYS:HD2	1:B:129:LYS:N	2.22	0.52
1:E:163:ASN:C	1:E:163:ASN:ND2	2.60	0.52
1:A:163:ASN:ND2	1:A:165:LYS:H	2.07	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:128:PHE:CD2	1:F:136:ARG:HD3	2.43	0.52
1:A:31:ARG:HH21	1:C:191:LYS:NZ	2.06	0.52
1:G:72:ASN:HA	1:G:154:GLU:O	2.10	0.52
1:B:103:GLY:O	1:B:107:ASN:HB2	2.10	0.52
1:E:145:SER:OG	1:E:178:TYR:HB3	2.08	0.52
1:H:97:LEU:HD13	1:H:147:TYR:HB2	1.92	0.52
1:E:101:LEU:HG	1:E:105:MET:CE	2.39	0.52
1:F:101:LEU:HG	1:F:105:MET:CE	2.40	0.52
1:C:175:PHE:HA	1:C:179:GLU:HB2	1.92	0.52
1:E:55:LEU:HD22	1:E:55:LEU:H	1.75	0.52
1:H:196:HIS:CE1	1:H:202:ILE:HG23	2.44	0.52
1:F:18:TYR:HA	1:F:22:ARG:CG	2.40	0.52
1:B:101:LEU:HG	1:B:105:MET:CE	2.40	0.52
1:E:163:ASN:ND2	1:E:165:LYS:H	2.07	0.51
1:A:175:PHE:HA	1:A:179:GLU:HB2	1.93	0.51
1:G:172:ASP:HA	1:G:177:LYS:HG3	1.91	0.51
1:G:196:HIS:CE1	1:G:202:ILE:HG23	2.45	0.51
1:A:196:HIS:CE1	1:A:202:ILE:HG23	2.46	0.51
1:B:5:ARG:HG2	1:C:185:ASN:HB2	1.91	0.51
1:H:55:LEU:H	1:H:55:LEU:HD22	1.74	0.51
1:G:101:LEU:HG	1:G:105:MET:CE	2.40	0.51
2:E:1179:N:H1'	3:E:359:NCA:H2	1.92	0.51
1:E:27:LEU:HB3	1:E:67:GLN:NE2	2.25	0.51
1:A:62:PHE:CZ	1:A:143:ALA:HB2	2.45	0.51
1:D:129:LYS:HB3	1:G:217:LEU:CD2	2.40	0.51
1:G:101:LEU:HG	1:G:105:MET:HE1	1.93	0.51
1:A:55:LEU:HD22	1:A:55:LEU:H	1.74	0.51
1:G:129:LYS:HD2	1:G:129:LYS:N	2.23	0.50
1:B:123:LYS:HD3	1:C:149:HIS:HE1	1.75	0.50
1:H:163:ASN:ND2	1:H:165:LYS:H	2.08	0.50
1:G:180:LEU:HB3	1:G:181:PRO:HD3	1.94	0.50
1:D:133:VAL:HG11	1:G:177:LYS:HD2	1.93	0.50
1:E:196:HIS:HE1	1:E:204:GLU:O	1.94	0.50
1:C:62:PHE:CZ	1:C:143:ALA:HB2	2.46	0.50
1:D:62:PHE:CZ	1:D:143:ALA:HB2	2.46	0.50
1:F:55:LEU:HD22	1:F:55:LEU:H	1.77	0.50
1:B:128:PHE:CD2	1:B:136:ARG:HD3	2.47	0.50
1:B:136:ARG:O	3:B:429:NCA:H5	2.12	0.50
1:A:69:LEU:N	1:A:69:LEU:HD13	2.26	0.50
1:G:62:PHE:CZ	1:G:143:ALA:HB2	2.47	0.50
1:E:103:GLY:O	1:E:107:ASN:HB2	2.12	0.50
1:H:101:LEU:HG	1:H:105:MET:CE	2.41	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:55:LEU:N	1:C:55:LEU:HD22	2.26	0.49
1:C:85:HIS:NE2	1:C:98:GLU:OE1	2.42	0.49
1:E:97:LEU:HD13	1:E:147:TYR:HB2	1.93	0.49
1:F:101:LEU:HG	1:F:105:MET:HE1	1.94	0.49
1:H:72:ASN:OD1	1:H:154:GLU:N	2.25	0.49
1:D:129:LYS:CG	1:G:171:PRO:HG3	2.41	0.49
1:E:106:LEU:N	1:E:106:LEU:HD22	2.27	0.49
1:B:69:LEU:HD22	1:B:150:SER:HB2	1.95	0.49
1:G:163:ASN:ND2	1:G:165:LYS:H	2.11	0.49
1:E:137:GLU:OE1	1:E:177:LYS:NZ	2.44	0.49
1:H:196:HIS:HE1	1:H:204:GLU:O	1.95	0.49
1:D:18:TYR:HA	1:D:22:ARG:CG	2.41	0.49
1:A:129:LYS:N	1:A:129:LYS:CD	2.70	0.49
1:D:129:LYS:CB	1:G:217:LEU:CD2	2.91	0.49
1:D:163:ASN:ND2	1:D:165:LYS:H	2.10	0.49
1:B:196:HIS:HE1	1:B:204:GLU:O	1.96	0.49
1:A:18:TYR:HA	1:A:22:ARG:CG	2.43	0.49
1:H:18:TYR:HA	1:H:22:ARG:CG	2.42	0.49
1:C:69:LEU:HD22	1:C:150:SER:HB2	1.91	0.49
1:B:5:ARG:CD	1:C:185:ASN:O	2.58	0.49
1:A:101:LEU:HG	1:A:105:MET:HE1	1.94	0.49
1:G:175:PHE:HA	1:G:179:GLU:HB2	1.94	0.49
1:C:55:LEU:H	1:C:55:LEU:HD22	1.77	0.48
1:H:106:LEU:HD22	1:H:106:LEU:N	2.28	0.48
1:A:1:ILE:HG22	1:A:125:CYS:O	2.12	0.48
1:H:1:ILE:HG22	1:H:125:CYS:O	2.13	0.48
1:B:124:VAL:HG22	1:B:125:CYS:N	2.27	0.48
1:A:101:LEU:HG	1:A:105:MET:CE	2.43	0.48
1:G:18:TYR:HA	1:G:22:ARG:CG	2.44	0.48
1:B:55:LEU:CD2	1:B:55:LEU:H	2.26	0.48
1:E:98:GLU:OE2	3:E:359:NCA:N7	2.46	0.48
1:F:55:LEU:HD22	1:F:55:LEU:N	2.26	0.48
1:F:180:LEU:HB3	1:F:181:PRO:HD3	1.95	0.48
1:C:103:GLY:O	1:C:107:ASN:HB2	2.14	0.48
1:D:196:HIS:HE1	1:D:204:GLU:O	1.97	0.48
1:A:27:LEU:HD21	1:A:95:ILE:HD13	1.96	0.48
2:G:1179:N:H1'	3:G:379:NCA:C6	2.43	0.48
1:C:1:ILE:HG22	1:C:125:CYS:O	2.13	0.48
1:E:165:LYS:HG2	1:G:202:ILE:HG13	1.95	0.48
1:F:67:GLN:O	1:F:69:LEU:HD13	2.13	0.48
1:C:129:LYS:CD	1:C:129:LYS:N	2.72	0.48
1:H:163:ASN:C	1:H:163:ASN:ND2	2.67	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:106:LEU:HD22	1:C:106:LEU:N	2.29	0.48
1:D:101:LEU:HG	1:D:105:MET:HE1	1.95	0.47
1:D:139:PHE:HB3	3:D:449:NCA:C6	2.44	0.47
1:F:196:HIS:CE1	1:F:202:ILE:HG23	2.49	0.47
1:D:1:ILE:HG22	1:D:125:CYS:O	2.14	0.47
1:F:18:TYR:HA	1:F:22:ARG:HG2	1.96	0.47
1:C:250:LEU:CD1	1:D:232:ARG:HD2	2.45	0.47
1:F:196:HIS:HE1	1:F:204:GLU:O	1.97	0.47
1:B:18:TYR:HA	1:B:22:ARG:CG	2.43	0.47
1:G:152:GLU:CD	1:G:186:LYS:HD3	2.34	0.47
1:B:18:TYR:HA	1:B:22:ARG:HG2	1.95	0.47
1:B:175:PHE:HA	1:B:179:GLU:HB2	1.96	0.47
1:A:250:LEU:CD1	1:B:232:ARG:HD2	2.45	0.47
1:C:69:LEU:HD22	1:C:150:SER:HB3	1.96	0.47
1:D:27:LEU:HD23	1:D:70:PRO:HD3	1.96	0.47
1:D:129:LYS:N	1:D:129:LYS:HD2	2.28	0.47
1:A:103:GLY:O	1:A:107:ASN:HB2	2.15	0.47
1:C:163:ASN:C	1:C:163:ASN:ND2	2.67	0.47
1:G:103:GLY:O	1:G:107:ASN:HB2	2.14	0.47
1:D:106:LEU:N	1:D:106:LEU:HD22	2.30	0.47
1:G:1:ILE:HG22	1:G:125:CYS:O	2.15	0.47
1:A:180:LEU:HB3	1:A:181:PRO:HD3	1.95	0.47
1:A:106:LEU:N	1:A:106:LEU:HD22	2.29	0.47
1:E:101:LEU:HG	1:E:105:MET:HE2	1.96	0.47
1:C:180:LEU:HB3	1:C:181:PRO:HD3	1.95	0.47
1:E:180:LEU:HB3	1:E:181:PRO:HD3	1.96	0.47
1:G:128:PHE:CD2	1:G:136:ARG:HD2	2.50	0.47
1:H:124:VAL:HG22	1:H:125:CYS:N	2.30	0.47
1:D:139:PHE:HB3	3:D:449:NCA:C5	2.45	0.47
1:E:62:PHE:CZ	1:E:143:ALA:HB2	2.50	0.47
1:F:62:PHE:CZ	1:F:143:ALA:HB2	2.50	0.47
1:G:236:PHE:CZ	1:H:248:CYS:HB3	2.50	0.47
1:F:103:GLY:O	1:F:107:ASN:HB2	2.15	0.47
2:G:1179:N:H1'	3:G:379:NCA:H6	1.97	0.47
1:H:69:LEU:CD2	1:H:150:SER:HB2	2.45	0.46
1:C:101:LEU:HG	1:C:105:MET:CE	2.45	0.46
1:A:144:SER:HB3	1:A:179:GLU:HG3	1.97	0.46
1:A:196:HIS:HE1	1:A:204:GLU:O	1.98	0.46
1:E:18:TYR:HA	1:E:22:ARG:CG	2.44	0.46
1:H:55:LEU:CD2	1:H:55:LEU:H	2.28	0.46
1:E:50:PRO:HB2	1:E:126:PRO:HG3	1.97	0.46
1:E:110:VAL:HG13	3:E:459:NCA:O7	2.16	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:103:GLY:O	1:D:107:ASN:HB2	2.15	0.46
1:A:163:ASN:ND2	1:A:163:ASN:C	2.66	0.46
1:C:221:LYS:O	1:C:222:HIS:HB2	2.16	0.46
1:G:55:LEU:H	1:G:55:LEU:HD22	1.81	0.46
1:E:163:ASN:HD22	1:E:164:PRO:N	2.14	0.46
1:F:163:ASN:ND2	1:F:165:LYS:H	2.13	0.46
1:A:55:LEU:CD2	1:A:55:LEU:H	2.29	0.46
1:H:72:ASN:HA	1:H:154:GLU:O	2.16	0.46
1:D:128:PHE:CD2	1:D:136:ARG:HD3	2.51	0.46
1:D:124:VAL:HG22	1:D:125:CYS:N	2.29	0.46
1:D:167:PRO:HG2	1:D:170:ARG:CG	2.45	0.46
1:F:152:GLU:HG2	1:F:186:LYS:HB3	1.98	0.46
1:F:69:LEU:HA	1:F:70:PRO:HD2	1.70	0.46
1:G:244:ASN:O	1:G:245:ALA:C	2.54	0.46
1:E:246:ARG:HD3	1:E:246:ARG:O	2.16	0.46
1:H:62:PHE:CZ	1:H:143:ALA:HB2	2.51	0.46
1:C:50:PRO:HB2	1:C:126:PRO:HG3	1.98	0.46
1:B:167:PRO:HG2	1:B:170:ARG:CG	2.46	0.45
1:C:140:TRP:HD1	3:C:439:NCA:C5	2.28	0.45
1:C:139:PHE:HB3	3:C:439:NCA:N1	2.31	0.45
1:D:90:THR:O	1:D:90:THR:HG22	2.16	0.45
1:A:69:LEU:HD22	1:A:150:SER:HB3	1.98	0.45
1:A:31:ARG:NH2	1:C:191:LYS:HZ2	2.13	0.45
1:B:65:ALA:CB	1:B:102:PRO:HG3	2.46	0.45
1:C:196:HIS:HE1	1:C:204:GLU:O	2.00	0.45
1:A:217:LEU:O	1:A:220:ALA:HB3	2.16	0.45
1:G:196:HIS:HE1	1:G:204:GLU:O	1.99	0.45
1:G:18:TYR:HA	1:G:22:ARG:HG2	1.98	0.45
1:D:107:ASN:O	1:D:108:SER:HB2	2.16	0.45
1:B:98:GLU:H	1:B:98:GLU:CD	2.20	0.45
1:B:70:PRO:HB2	1:B:73:LYS:CG	2.47	0.45
1:H:69:LEU:HD22	1:H:150:SER:CB	2.47	0.45
1:D:55:LEU:H	1:D:55:LEU:CD2	2.29	0.45
1:C:124:VAL:HG22	1:C:125:CYS:N	2.31	0.45
1:H:163:ASN:HA	1:H:164:PRO:HD2	1.85	0.45
1:F:167:PRO:HG2	1:F:170:ARG:CG	2.45	0.45
1:H:246:ARG:HD3	1:H:246:ARG:O	2.17	0.45
1:E:68:GLN:HE21	1:E:68:GLN:HB2	1.50	0.45
1:E:140:TRP:HD1	3:E:459:NCA:H5	1.80	0.45
1:D:177:LYS:HG2	1:G:133:VAL:HG13	1.99	0.45
1:D:18:TYR:HA	1:D:22:ARG:HG2	1.97	0.45
1:A:18:TYR:HA	1:A:22:ARG:HG2	1.99	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:124:VAL:HG22	1:A:125:CYS:N	2.32	0.45
1:E:18:TYR:HA	1:E:22:ARG:HG2	1.98	0.45
1:G:124:VAL:HG22	1:G:125:CYS:N	2.32	0.45
1:F:69:LEU:HD22	1:F:150:SER:CB	2.39	0.45
1:B:27:LEU:HD21	1:B:95:ILE:HD13	1.98	0.45
1:D:180:LEU:HB3	1:D:181:PRO:HD3	1.98	0.45
1:F:14:ARG:NH2	1:F:107:ASN:O	2.49	0.45
1:B:246:ARG:HD3	1:B:246:ARG:O	2.16	0.45
1:F:220:ALA:C	1:F:222:HIS:H	2.21	0.45
1:E:163:ASN:HA	1:E:164:PRO:HD2	1.85	0.44
1:C:62:PHE:CE1	1:C:143:ALA:HA	2.52	0.44
1:H:27:LEU:HD21	1:H:95:ILE:HD13	1.99	0.44
1:B:70:PRO:HB2	1:B:73:LYS:HG2	2.00	0.44
1:H:180:LEU:HB3	1:H:181:PRO:HD3	1.99	0.44
1:B:180:LEU:HB3	1:B:181:PRO:HD3	1.98	0.44
1:C:71:LYS:HG2	1:C:72:ASN:HD21	1.77	0.44
1:E:55:LEU:H	1:E:55:LEU:CD2	2.29	0.44
1:E:55:LEU:HD13	1:E:55:LEU:HA	1.84	0.44
1:H:18:TYR:HA	1:H:22:ARG:HG2	1.99	0.44
1:B:22:ARG:HA	1:B:22:ARG:HD3	1.83	0.44
1:G:167:PRO:HG2	1:G:170:ARG:CG	2.45	0.44
1:A:74:VAL:HG13	1:A:147:TYR:CZ	2.52	0.44
1:H:69:LEU:HD22	1:H:150:SER:HB2	1.99	0.44
1:B:163:ASN:ND2	1:B:163:ASN:C	2.67	0.44
1:D:133:VAL:CG1	1:G:177:LYS:CG	2.96	0.44
1:C:163:ASN:HD22	1:C:165:LYS:H	1.66	0.44
1:F:163:ASN:ND2	1:F:163:ASN:C	2.70	0.44
1:F:106:LEU:HD22	1:F:106:LEU:N	2.32	0.44
1:C:69:LEU:HD21	1:C:150:SER:HB2	1.96	0.44
1:A:166:VAL:HG13	1:A:170:ARG:HD3	2.00	0.44
1:D:163:ASN:HA	1:D:164:PRO:HD2	1.83	0.44
1:G:55:LEU:HD22	1:G:55:LEU:N	2.32	0.43
1:B:90:THR:HG22	1:B:90:THR:O	2.18	0.43
1:C:101:LEU:HG	1:C:105:MET:HE1	1.99	0.43
1:A:220:ALA:C	1:A:222:HIS:H	2.22	0.43
1:A:14:ARG:NH2	1:A:107:ASN:O	2.51	0.43
1:C:55:LEU:CD2	1:C:55:LEU:H	2.32	0.43
1:A:50:PRO:HB2	1:A:126:PRO:HG3	2.01	0.43
2:G:1179:N:C1'	3:G:379:NCA:H6	2.47	0.43
1:G:106:LEU:HD22	1:G:106:LEU:N	2.33	0.43
1:A:71:LYS:O	1:A:72:ASN:CB	2.63	0.43
1:F:55:LEU:H	1:F:55:LEU:CD2	2.32	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:73:LYS:HE2	1:F:93:LYS:O	2.18	0.43
1:H:14:ARG:NH2	1:H:107:ASN:O	2.51	0.43
1:E:124:VAL:HG22	1:E:125:CYS:N	2.33	0.43
1:C:144:SER:HB3	1:C:179:GLU:HG3	2.01	0.43
1:C:90:THR:O	1:C:90:THR:HG22	2.18	0.43
1:C:217:LEU:O	1:C:220:ALA:HB3	2.19	0.43
1:G:246:ARG:O	1:G:246:ARG:HD3	2.19	0.43
1:F:120:PHE:N	1:F:120:PHE:CD1	2.87	0.43
1:B:26:ILE:N	1:B:26:ILE:CD1	2.81	0.43
1:E:14:ARG:NH2	1:E:107:ASN:O	2.49	0.42
1:B:163:ASN:HA	1:B:164:PRO:HD2	1.82	0.42
1:E:204:GLU:HB3	1:E:211:LEU:HD22	2.01	0.42
1:G:62:PHE:CE1	1:G:143:ALA:HA	2.54	0.42
1:B:244:ASN:O	1:B:245:ALA:C	2.57	0.42
1:B:217:LEU:O	1:B:220:ALA:HB3	2.19	0.42
1:A:62:PHE:CE1	1:A:143:ALA:HA	2.54	0.42
1:E:72:ASN:HA	1:E:154:GLU:O	2.18	0.42
1:E:11:PHE:C	1:E:11:PHE:CD1	2.93	0.42
1:G:90:THR:HG22	1:G:90:THR:O	2.19	0.42
1:F:246:ARG:HD3	1:F:246:ARG:O	2.20	0.42
1:E:65:ALA:CB	1:E:102:PRO:HG3	2.49	0.42
1:D:246:ARG:HD3	1:D:246:ARG:O	2.19	0.42
1:F:62:PHE:CE1	1:F:143:ALA:HA	2.54	0.42
1:F:1:ILE:HG22	1:F:125:CYS:O	2.20	0.42
1:D:67:GLN:O	1:D:69:LEU:HD13	2.19	0.42
1:A:113:GLY:HA2	1:A:120:PHE:HA	2.00	0.42
1:H:221:LYS:O	1:H:222:HIS:HB2	2.19	0.42
1:E:128:PHE:HB3	1:E:129:LYS:CD	2.41	0.42
1:D:65:ALA:CB	1:D:102:PRO:HG3	2.49	0.42
1:H:101:LEU:HG	1:H:105:MET:HE2	2.01	0.42
1:B:1:ILE:HG22	1:B:125:CYS:O	2.19	0.42
1:C:189:ARG:HG2	1:C:190:VAL:N	2.34	0.42
1:H:90:THR:O	1:H:90:THR:HG22	2.19	0.42
1:F:124:VAL:HG22	1:F:125:CYS:N	2.35	0.42
1:B:106:LEU:N	1:B:106:LEU:HD22	2.33	0.42
1:F:244:ASN:O	1:F:245:ALA:C	2.58	0.42
1:A:246:ARG:O	1:A:246:ARG:HD3	2.20	0.42
1:H:55:LEU:HD13	1:H:55:LEU:HA	1.90	0.42
1:E:220:ALA:C	1:E:222:HIS:H	2.22	0.42
1:D:244:ASN:O	1:D:245:ALA:C	2.58	0.42
1:C:246:ARG:HD3	1:C:246:ARG:O	2.20	0.41
1:H:98:GLU:CD	1:H:98:GLU:H	2.23	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:65:ALA:CB	1:A:102:PRO:HG3	2.48	0.41
1:A:174:PHE:HB2	2:A:1179:N:O5'	2.20	0.41
1:A:55:LEU:HD13	1:A:55:LEU:HA	1.87	0.41
1:H:220:ALA:C	1:H:222:HIS:H	2.24	0.41
1:H:50:PRO:HB2	1:H:126:PRO:HG3	2.02	0.41
1:F:87:TYR:CE2	1:F:94:TYR:HE2	2.39	0.41
1:G:152:GLU:OE2	1:G:186:LYS:HD3	2.21	0.41
1:G:202:ILE:HG21	1:G:205:LYS:HE3	2.02	0.41
1:H:113:GLY:HA2	1:H:120:PHE:HA	2.03	0.41
1:F:128:PHE:HB3	1:F:129:LYS:CD	2.43	0.41
1:H:166:VAL:HG13	1:H:170:ARG:HD3	2.03	0.41
1:G:163:ASN:C	1:G:163:ASN:ND2	2.72	0.41
1:F:189:ARG:HG2	1:F:190:VAL:N	2.35	0.41
1:D:220:ALA:C	1:D:222:HIS:H	2.22	0.41
1:F:50:PRO:HB2	1:F:126:PRO:HG3	2.01	0.41
1:G:40:ASP:HB3	1:G:61:PHE:HA	2.01	0.41
1:F:90:THR:HG22	1:F:90:THR:O	2.21	0.41
1:E:90:THR:O	1:E:90:THR:HG22	2.20	0.41
1:G:220:ALA:C	1:G:222:HIS:H	2.23	0.41
1:D:69:LEU:HD22	1:D:150:SER:CB	2.40	0.41
1:C:174:PHE:HB2	2:C:1179:N:O5'	2.20	0.41
1:G:27:LEU:HA	1:G:28:PRO:HD3	1.91	0.41
1:A:128:PHE:HB3	1:A:129:LYS:CD	2.42	0.41
1:H:65:ALA:CB	1:H:102:PRO:HG3	2.50	0.41
1:G:251:ALA:HB3	1:H:249:ARG:HB2	2.02	0.41
1:B:50:PRO:HB2	1:B:126:PRO:HG3	2.01	0.41
1:A:72:ASN:HA	1:A:154:GLU:O	2.21	0.41
1:F:67:GLN:OE1	1:F:100:THR:HB	2.21	0.41
1:H:163:ASN:HD22	1:H:165:LYS:H	1.68	0.41
1:G:107:ASN:O	1:G:108:SER:HB2	2.20	0.41
1:E:40:ASP:HB3	1:E:61:PHE:HA	2.03	0.41
1:E:113:GLY:HA2	1:E:120:PHE:HA	2.03	0.41
1:G:113:GLY:HA2	1:G:120:PHE:HA	2.02	0.41
1:A:90:THR:O	1:A:90:THR:HG22	2.20	0.41
1:H:139:PHE:HB3	3:H:489:NCA:C6	2.51	0.41
1:E:166:VAL:HG13	1:E:170:ARG:HD3	2.03	0.41
1:A:67:GLN:O	1:A:69:LEU:HD13	2.20	0.41
1:A:163:ASN:HD22	1:A:165:LYS:H	1.69	0.41
1:C:22:ARG:HA	1:C:22:ARG:HD3	1.78	0.41
1:F:18:TYR:O	1:F:29:ARG:NH1	2.54	0.41
1:C:250:LEU:HD12	1:D:232:ARG:HD2	2.03	0.41
1:H:49:ASN:O	1:H:50:PRO:C	2.59	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:189:ARG:HG2	1:E:190:VAL:N	2.36	0.41
1:A:17:ASP:CG	1:B:14:ARG:HD2	2.39	0.41
1:E:1:ILE:HG23	1:E:125:CYS:HB2	2.03	0.41
1:A:189:ARG:HG2	1:A:190:VAL:N	2.36	0.41
1:F:163:ASN:HA	1:F:164:PRO:HD2	1.81	0.40
1:D:120:PHE:CD1	1:D:120:PHE:N	2.89	0.40
1:B:14:ARG:NH2	1:B:107:ASN:O	2.52	0.40
1:G:87:TYR:CE2	1:G:94:TYR:HE2	2.39	0.40
1:B:163:ASN:HD22	1:B:165:LYS:H	1.68	0.40
1:B:27:LEU:HA	1:B:28:PRO:HD3	1.86	0.40
1:B:113:GLY:HA2	1:B:120:PHE:HA	2.02	0.40
1:H:87:TYR:CE2	1:H:94:TYR:HE2	2.39	0.40
2:A:1179:N:H1'	3:A:319:NCA:H6	2.03	0.40
1:D:27:LEU:HA	1:D:28:PRO:HD3	1.89	0.40
1:B:220:ALA:C	1:B:222:HIS:H	2.23	0.40
1:C:73:LYS:HD3	1:C:93:LYS:O	2.21	0.40
1:E:248:CYS:HB3	1:F:236:PHE:CZ	2.56	0.40
1:A:244:ASN:O	1:A:245:ALA:C	2.59	0.40
1:C:166:VAL:HG13	1:C:170:ARG:HD3	2.04	0.40
1:A:251:ALA:HA	1:B:251:ALA:CA	2.47	0.40
1:D:163:ASN:C	1:D:163:ASN:ND2	2.69	0.40
1:H:22:ARG:HD3	1:H:22:ARG:HA	1.78	0.40
1:E:217:LEU:O	1:E:220:ALA:HB3	2.21	0.40
1:B:40:ASP:HB3	1:B:61:PHE:HA	2.03	0.40
1:G:232:ARG:HH22	1:H:244:ASN:HA	1.86	0.40
1:G:248:CYS:HB3	1:H:236:PHE:CZ	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:122:GLU:OE1	1:H:31:ARG:NH2[1_545]	1.99	0.21

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	249/258 (96%)	233 (94%)	15 (6%)	1 (0%)	43	61
1	B	249/258 (96%)	235 (94%)	14 (6%)	0	100	100
1	C	249/258 (96%)	232 (93%)	15 (6%)	2 (1%)	27	39
1	D	249/258 (96%)	230 (92%)	18 (7%)	1 (0%)	43	61
1	E	249/258 (96%)	234 (94%)	14 (6%)	1 (0%)	43	61
1	F	249/258 (96%)	234 (94%)	13 (5%)	2 (1%)	27	39
1	G	249/258 (96%)	234 (94%)	13 (5%)	2 (1%)	27	39
1	H	249/258 (96%)	232 (93%)	16 (6%)	1 (0%)	43	61
All	All	1992/2064 (96%)	1864 (94%)	118 (6%)	10 (0%)	38	53

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	128	PHE
1	E	128	PHE
1	F	128	PHE
1	G	128	PHE
1	H	128	PHE
1	A	128	PHE
1	C	128	PHE
1	G	80	VAL
1	C	80	VAL
1	F	80	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%)	208 (94%)	12 (6%)	30	46
1	B	220/226 (97%)	209 (95%)	11 (5%)	34	51
1	C	220/226 (97%)	208 (94%)	12 (6%)	30	46
1	D	220/226 (97%)	210 (96%)	10 (4%)	38	57
1	E	220/226 (97%)	209 (95%)	11 (5%)	34	51

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	220/226 (97%)	209 (95%)	11 (5%)	34	51
1	G	220/226 (97%)	212 (96%)	8 (4%)	47	68
1	H	220/226 (97%)	210 (96%)	10 (4%)	38	57
All	All	1760/1808 (97%)	1675 (95%)	85 (5%)	35	53

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLU
1	A	16	LYS
1	A	30	VAL
1	A	32	SER
1	A	69	LEU
1	A	133	VAL
1	A	134	GLN
1	A	155	VAL
1	A	163	ASN
1	A	211	LEU
1	A	242	ASN
1	A	246	ARG
1	B	8	GLU
1	B	30	VAL
1	B	32	SER
1	B	69	LEU
1	B	133	VAL
1	B	134	GLN
1	B	155	VAL
1	B	163	ASN
1	B	211	LEU
1	B	242	ASN
1	B	246	ARG
1	C	8	GLU
1	C	16	LYS
1	C	30	VAL
1	C	32	SER
1	C	69	LEU
1	C	133	VAL
1	C	134	GLN
1	C	155	VAL
1	C	163	ASN
1	C	211	LEU

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Mol	Chain	Res	Type
1	C	242	ASN
1	C	246	ARG
1	D	8	GLU
1	D	30	VAL
1	D	32	SER
1	D	69	LEU
1	D	133	VAL
1	D	134	GLN
1	D	163	ASN
1	D	211	LEU
1	D	242	ASN
1	D	246	ARG
1	E	8	GLU
1	E	30	VAL
1	E	32	SER
1	E	68	GLN
1	E	69	LEU
1	E	133	VAL
1	E	134	GLN
1	E	163	ASN
1	E	211	LEU
1	E	242	ASN
1	E	246	ARG
1	F	8	GLU
1	F	30	VAL
1	F	32	SER
1	F	69	LEU
1	F	133	VAL
1	F	134	GLN
1	F	155	VAL
1	F	163	ASN
1	F	211	LEU
1	F	242	ASN
1	F	246	ARG
1	G	8	GLU
1	G	30	VAL
1	G	133	VAL
1	G	134	GLN
1	G	163	ASN
1	G	211	LEU
1	G	242	ASN
1	G	246	ARG

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Mol	Chain	Res	Type
1	H	8	GLU
1	H	30	VAL
1	H	32	SER
1	H	69	LEU
1	H	133	VAL
1	H	134	GLN
1	H	163	ASN
1	H	211	LEU
1	H	242	ASN
1	H	246	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (52) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	68	GLN
1	A	89	ASN
1	A	134	GLN
1	A	163	ASN
1	A	196	HIS
1	A	242	ASN
1	B	66	GLN
1	B	89	ASN
1	B	134	GLN
1	B	163	ASN
1	B	196	HIS
1	B	242	ASN
1	C	66	GLN
1	C	89	ASN
1	C	134	GLN
1	C	149	HIS
1	C	163	ASN
1	C	196	HIS
1	C	242	ASN
1	D	66	GLN
1	D	89	ASN
1	D	134	GLN
1	D	163	ASN
1	D	196	HIS
1	D	242	ASN
1	E	66	GLN
1	E	68	GLN

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Mol	Chain	Res	Type
1	E	89	ASN
1	E	134	GLN
1	E	163	ASN
1	E	196	HIS
1	E	242	ASN
1	F	66	GLN
1	F	89	ASN
1	F	134	GLN
1	F	163	ASN
1	F	196	HIS
1	F	242	ASN
1	G	66	GLN
1	G	68	GLN
1	G	89	ASN
1	G	134	GLN
1	G	163	ASN
1	G	196	HIS
1	G	242	ASN
1	H	66	GLN
1	H	89	ASN
1	H	134	GLN
1	H	163	ASN
1	H	196	HIS
1	H	242	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 ⓘ

24 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	N	A	1179	1	13,13,13	3.34	6 (46%)	19,19,19	2.29	9 (47%)
3	NCA	A	319	-	9,9,9	2.90	5 (55%)	11,11,11	1.87	4 (36%)
3	NCA	A	419	-	9,9,9	2.17	3 (33%)	11,11,11	1.79	3 (27%)
2	N	B	1179	1	13,13,13	3.57	5 (38%)	19,19,19	2.47	9 (47%)
3	NCA	B	329	-	9,9,9	2.52	5 (55%)	11,11,11	1.75	4 (36%)
3	NCA	B	429	-	9,9,9	2.71	4 (44%)	11,11,11	1.81	5 (45%)
2	N	C	1179	1	13,13,13	3.45	6 (46%)	19,19,19	2.34	10 (52%)
3	NCA	C	339	-	9,9,9	2.65	3 (33%)	11,11,11	1.83	4 (36%)
3	NCA	C	439	-	9,9,9	2.63	3 (33%)	11,11,11	1.94	4 (36%)
2	N	D	1179	1	13,13,13	3.31	5 (38%)	19,19,19	2.31	9 (47%)
3	NCA	D	349	-	9,9,9	2.62	4 (44%)	11,11,11	1.75	4 (36%)
3	NCA	D	449	-	9,9,9	2.73	5 (55%)	11,11,11	1.69	3 (27%)
2	N	E	1179	1	13,13,13	3.46	6 (46%)	19,19,19	2.35	9 (47%)
3	NCA	E	359	-	9,9,9	2.40	4 (44%)	11,11,11	1.60	3 (27%)
3	NCA	E	459	-	9,9,9	2.34	4 (44%)	11,11,11	1.84	3 (27%)
2	N	F	1179	1	13,13,13	3.45	6 (46%)	19,19,19	2.40	9 (47%)
3	NCA	F	369	-	9,9,9	2.78	4 (44%)	11,11,11	1.77	4 (36%)
3	NCA	F	469	-	9,9,9	3.07	7 (77%)	11,11,11	1.84	4 (36%)
2	N	G	1179	1	13,13,13	3.14	5 (38%)	19,19,19	2.37	9 (47%)
3	NCA	G	379	-	9,9,9	2.81	5 (55%)	11,11,11	2.00	3 (27%)
3	NCA	G	479	-	9,9,9	3.35	7 (77%)	11,11,11	2.01	4 (36%)
2	N	H	1179	1	13,13,13	3.26	6 (46%)	19,19,19	2.31	10 (52%)
3	NCA	H	389	-	9,9,9	3.21	5 (55%)	11,11,11	1.94	5 (45%)
3	NCA	H	489	-	9,9,9	2.11	2 (22%)	11,11,11	1.43	1 (9%)

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	N	A	1179	1	1/1/4/4	0/6/19/19	0/1/1/1
3	NCA	A	319	-	-	0/4/4/4	0/1/1/1
3	NCA	A	419	-	-	0/4/4/4	0/1/1/1
2	N	B	1179	1	1/1/4/4	0/6/19/19	0/1/1/1
3	NCA	B	329	-	-	0/4/4/4	0/1/1/1
3	NCA	B	429	-	-	0/4/4/4	0/1/1/1
2	N	C	1179	1	1/1/4/4	0/6/19/19	0/1/1/1
3	NCA	C	339	-	-	0/4/4/4	0/1/1/1
3	NCA	C	439	-	-	0/4/4/4	0/1/1/1
2	N	D	1179	1	1/1/4/4	0/6/19/19	0/1/1/1
3	NCA	D	349	-	-	0/4/4/4	0/1/1/1
3	NCA	D	449	-	-	0/4/4/4	0/1/1/1
2	N	E	1179	1	1/1/4/4	0/6/19/19	0/1/1/1
3	NCA	E	359	-	-	0/4/4/4	0/1/1/1
3	NCA	E	459	-	-	0/4/4/4	0/1/1/1
2	N	F	1179	1	1/1/4/4	0/6/19/19	0/1/1/1
3	NCA	F	369	-	-	0/4/4/4	0/1/1/1
3	NCA	F	469	-	-	0/4/4/4	0/1/1/1
2	N	G	1179	1	1/1/4/4	0/6/19/19	0/1/1/1
3	NCA	G	379	-	-	0/4/4/4	0/1/1/1
3	NCA	G	479	-	-	0/4/4/4	0/1/1/1
2	N	H	1179	1	1/1/4/4	0/6/19/19	0/1/1/1
3	NCA	H	389	-	-	0/4/4/4	0/1/1/1
3	NCA	H	489	-	-	0/4/4/4	0/1/1/1

All (115) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1179	N	O4'-C4'	6.68	1.56	1.44
2	C	1179	N	C1'-C2'	6.58	1.64	1.51
2	F	1179	N	O4'-C4'	6.55	1.56	1.44
2	C	1179	N	O4'-C4'	6.51	1.56	1.44
2	B	1179	N	C2'-C3'	6.34	1.62	1.53
2	E	1179	N	C1'-C2'	6.30	1.63	1.51
2	G	1179	N	O4'-C4'	6.30	1.55	1.44
2	A	1179	N	C1'-C2'	6.30	1.63	1.51
2	B	1179	N	C1'-C2'	6.27	1.63	1.51
3	G	479	NCA	C2-N1	6.25	1.48	1.34
2	D	1179	N	O4'-C4'	6.25	1.55	1.44
2	E	1179	N	O4'-C4'	6.21	1.55	1.44
2	H	1179	N	O4'-C4'	6.19	1.55	1.44
2	A	1179	N	O4'-C4'	6.15	1.55	1.44
3	H	389	NCA	C3-C7	6.13	1.61	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	1179	N	C1'-C2'	6.05	1.63	1.51
2	D	1179	N	C1'-C2'	6.02	1.63	1.51
3	F	469	NCA	C2-N1	5.82	1.47	1.34
2	F	1179	N	C2'-C3'	5.80	1.62	1.53
2	F	1179	N	C1'-C2'	5.79	1.62	1.51
2	A	1179	N	P-O5'	5.70	1.79	1.60
2	C	1179	N	P-O5'	5.66	1.79	1.60
2	E	1179	N	P-O5'	5.58	1.79	1.60
2	E	1179	N	C2'-C3'	5.49	1.61	1.53
2	H	1179	N	P-O5'	5.46	1.79	1.60
2	G	1179	N	C1'-C2'	5.40	1.62	1.51
2	F	1179	N	P-O5'	5.24	1.78	1.60
3	F	369	NCA	C3-C7	5.23	1.59	1.50
2	D	1179	N	P-O5'	5.18	1.78	1.60
2	G	1179	N	P-O5'	5.15	1.77	1.60
3	C	339	NCA	C3-C7	5.13	1.59	1.50
3	D	449	NCA	C2-N1	5.10	1.45	1.34
2	B	1179	N	P-O5'	5.09	1.77	1.60
3	A	319	NCA	C3-C7	5.09	1.59	1.50
2	D	1179	N	C2'-C3'	5.06	1.61	1.53
3	D	349	NCA	C2-N1	5.00	1.45	1.34
3	E	359	NCA	C2-N1	4.78	1.45	1.34
2	C	1179	N	C2'-C3'	4.75	1.60	1.53
3	B	329	NCA	C2-N1	4.68	1.44	1.34
3	G	379	NCA	C3-C7	4.64	1.58	1.50
3	B	429	NCA	C2-N1	4.61	1.44	1.34
3	H	389	NCA	C2-N1	4.59	1.44	1.34
2	A	1179	N	C2'-C3'	4.56	1.60	1.53
3	H	489	NCA	C2-N1	4.56	1.44	1.34
3	C	439	NCA	C2-N1	4.56	1.44	1.34
2	G	1179	N	C2'-C3'	4.55	1.60	1.53
3	A	319	NCA	C4-C3	4.46	1.47	1.39
3	C	439	NCA	C3-C7	4.45	1.58	1.50
3	E	459	NCA	C2-N1	4.38	1.44	1.34
2	H	1179	N	C2'-C3'	4.32	1.59	1.53
3	G	379	NCA	C4-C3	4.31	1.46	1.39
3	G	379	NCA	C2-N1	4.22	1.43	1.34
3	A	319	NCA	C2-N1	4.15	1.43	1.34
3	A	419	NCA	C2-N1	4.13	1.43	1.34
3	H	389	NCA	C4-C3	4.03	1.46	1.39
3	G	479	NCA	C4-C3	4.00	1.46	1.39
3	F	369	NCA	C2-N1	3.98	1.43	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	339	NCA	C2-N1	3.86	1.43	1.34
3	F	369	NCA	C4-C3	3.80	1.45	1.39
3	B	329	NCA	C4-C3	3.80	1.45	1.39
3	C	339	NCA	C4-C3	3.79	1.45	1.39
3	B	429	NCA	C4-C3	3.76	1.45	1.39
3	A	419	NCA	C3-C7	3.74	1.57	1.50
3	D	349	NCA	C4-C3	3.71	1.45	1.39
3	D	449	NCA	C3-C7	3.70	1.56	1.50
3	C	439	NCA	C4-C3	3.62	1.45	1.39
3	D	449	NCA	C2-C3	3.57	1.45	1.39
3	G	479	NCA	C3-C7	3.51	1.56	1.50
3	F	469	NCA	C2-C3	3.46	1.44	1.39
3	B	429	NCA	C3-C7	3.30	1.56	1.50
3	D	349	NCA	C3-C7	3.29	1.56	1.50
3	G	479	NCA	C2-C3	3.24	1.44	1.39
3	F	469	NCA	C4-C3	3.14	1.44	1.39
3	E	459	NCA	C3-C7	3.11	1.55	1.50
3	B	429	NCA	C7-N7	-3.06	1.26	1.33
3	H	389	NCA	C2-C3	3.03	1.44	1.39
3	E	359	NCA	C4-C3	2.98	1.44	1.39
3	E	459	NCA	C4-C3	2.92	1.44	1.39
2	F	1179	N	C3'-C4'	2.74	1.60	1.53
3	F	369	NCA	C2-C3	2.74	1.43	1.39
3	E	359	NCA	C3-C7	2.74	1.55	1.50
2	H	1179	N	C3'-C4'	2.70	1.60	1.53
3	F	469	NCA	C3-C7	2.67	1.55	1.50
3	F	469	NCA	C7-N7	-2.66	1.27	1.33
2	E	1179	N	C3'-C4'	2.65	1.60	1.53
3	G	479	NCA	C7-N7	-2.62	1.27	1.33
2	C	1179	N	C3'-C4'	2.62	1.60	1.53
3	D	449	NCA	C4-C3	2.61	1.43	1.39
3	G	479	NCA	C6-N1	2.59	1.41	1.33
3	A	419	NCA	C2-C3	2.53	1.43	1.39
3	F	469	NCA	O7-C7	-2.52	1.18	1.24
3	G	379	NCA	C2-C3	2.50	1.43	1.39
2	D	1179	N	C3'-C4'	2.50	1.59	1.53
3	B	329	NCA	C2-C3	2.46	1.43	1.39
3	B	329	NCA	C3-C7	2.44	1.54	1.50
2	B	1179	N	C3'-C4'	2.40	1.59	1.53
3	F	469	NCA	C6-N1	2.39	1.41	1.33
3	H	489	NCA	C2-C3	2.37	1.43	1.39
2	A	1179	N	C3'-C4'	2.36	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	319	NCA	C6-N1	2.34	1.41	1.33
3	A	319	NCA	C2-C3	2.32	1.42	1.39
3	E	359	NCA	C2-C3	2.25	1.42	1.39
3	G	379	NCA	C6-N1	2.24	1.40	1.33
3	D	349	NCA	C2-C3	2.23	1.42	1.39
3	G	479	NCA	O7-C7	-2.23	1.18	1.24
2	A	1179	N	O5'-C5'	2.23	1.54	1.44
3	H	389	NCA	C6-N1	2.21	1.40	1.33
3	B	329	NCA	C6-N1	2.14	1.40	1.33
3	E	459	NCA	C6-N1	2.12	1.40	1.33
3	D	449	NCA	C6-N1	2.09	1.40	1.33
2	F	1179	N	O5'-C5'	2.07	1.53	1.44
2	G	1179	N	C3'-C4'	2.07	1.58	1.53
2	H	1179	N	O5'-C5'	2.06	1.53	1.44
2	C	1179	N	O5'-C5'	2.06	1.53	1.44
2	E	1179	N	O5'-C5'	2.04	1.53	1.44

All (132) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1179	N	OP2-P-O5'	-4.83	93.31	106.65
2	G	1179	N	OP2-P-O5'	-4.58	94.00	106.65
2	F	1179	N	OP2-P-O5'	-4.49	94.27	106.65
2	D	1179	N	OP2-P-O5'	-4.44	94.38	106.65
2	G	1179	N	O4'-C1'-C2'	-4.40	97.05	106.14
2	B	1179	N	O4'-C1'-C2'	-4.40	97.06	106.14
2	F	1179	N	O4'-C1'-C2'	-4.33	97.19	106.14
2	E	1179	N	O4'-C1'-C2'	-4.27	97.33	106.14
2	C	1179	N	O4'-C1'-C2'	-4.26	97.34	106.14
2	H	1179	N	O4'-C1'-C2'	-4.26	97.35	106.14
2	C	1179	N	OP2-P-O5'	-4.18	95.11	106.65
2	D	1179	N	O4'-C1'-C2'	-4.16	97.56	106.14
2	A	1179	N	OP2-P-O5'	-4.09	95.37	106.65
2	A	1179	N	O4'-C1'-C2'	-4.08	97.72	106.14
2	E	1179	N	OP2-P-O5'	-4.05	95.46	106.65
2	H	1179	N	OP2-P-O5'	-3.75	96.29	106.65
2	B	1179	N	C1'-O4'-C4'	3.73	117.37	108.19
3	G	479	NCA	C3-C7-N7	3.69	121.97	117.77
3	G	379	NCA	C3-C7-N7	3.63	121.90	117.77
3	F	469	NCA	C3-C7-N7	3.56	121.83	117.77
2	F	1179	N	C1'-O4'-C4'	3.54	116.90	108.19
2	E	1179	N	C1'-O4'-C4'	3.50	116.80	108.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	459	NCA	C3-C7-N7	3.48	121.73	117.77
2	D	1179	N	C1'-O4'-C4'	3.41	116.59	108.19
2	G	1179	N	C1'-O4'-C4'	3.40	116.57	108.19
2	H	1179	N	C1'-O4'-C4'	3.40	116.57	108.19
2	C	1179	N	C1'-O4'-C4'	3.40	116.56	108.19
2	E	1179	N	O5'-C5'-C4'	3.35	121.22	108.94
2	C	1179	N	O5'-C5'-C4'	3.31	121.08	108.94
2	F	1179	N	O5'-C5'-C4'	3.26	120.91	108.94
2	B	1179	N	O5'-C5'-C4'	3.26	120.89	108.94
2	A	1179	N	C1'-O4'-C4'	3.26	116.21	108.19
2	H	1179	N	O5'-C5'-C4'	3.25	120.85	108.94
3	G	379	NCA	C5-C4-C3	3.20	124.47	120.32
3	A	319	NCA	C3-C7-N7	3.19	121.41	117.77
3	C	439	NCA	C3-C7-N7	3.19	121.40	117.77
2	A	1179	N	O5'-C5'-C4'	3.18	120.61	108.94
3	G	479	NCA	C5-C4-C3	3.18	124.44	120.32
3	A	419	NCA	C3-C7-N7	3.17	121.38	117.77
2	G	1179	N	O5'-C5'-C4'	3.16	120.54	108.94
2	D	1179	N	O5'-C5'-C4'	3.14	120.45	108.94
2	H	1179	N	O4'-C4'-C5'	-3.09	102.49	109.57
3	H	389	NCA	C5-C4-C3	3.06	124.30	120.32
3	C	339	NCA	C5-C4-C3	3.06	124.29	120.32
3	H	389	NCA	C4-C3-C2	-3.03	113.92	117.64
3	C	439	NCA	C4-C3-C2	-3.03	113.92	117.64
3	C	439	NCA	C5-C4-C3	2.97	124.18	120.32
3	A	319	NCA	C5-C4-C3	2.96	124.17	120.32
2	E	1179	N	O4'-C4'-C5'	-2.95	102.82	109.57
3	A	419	NCA	C4-C3-C2	-2.93	114.05	117.64
3	G	479	NCA	C4-C3-C2	-2.92	114.06	117.64
2	G	1179	N	OP2-P-OP1	2.91	119.97	110.44
2	F	1179	N	OP2-P-OP1	2.91	119.94	110.44
2	D	1179	N	OP2-P-OP1	2.89	119.89	110.44
3	C	339	NCA	C4-C3-C2	-2.89	114.09	117.64
2	F	1179	N	O4'-C4'-C5'	-2.88	102.97	109.57
2	C	1179	N	OP2-P-OP1	2.87	119.81	110.44
2	B	1179	N	OP2-P-OP1	2.86	119.79	110.44
3	D	449	NCA	C5-C4-C3	2.86	124.03	120.32
2	B	1179	N	O4'-C4'-C5'	-2.84	103.06	109.57
3	D	449	NCA	C4-C3-C2	-2.84	114.16	117.64
3	F	369	NCA	C5-C4-C3	2.82	123.99	120.32
3	B	329	NCA	C3-C7-N7	2.81	120.97	117.77
2	A	1179	N	O4'-C4'-C5'	-2.80	103.16	109.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	379	NCA	C4-C3-C2	-2.80	114.21	117.64
3	D	349	NCA	C4-C3-C2	-2.78	114.22	117.64
2	C	1179	N	O4'-C4'-C5'	-2.78	103.20	109.57
2	B	1179	N	O2'-C2'-C3'	2.77	116.67	111.31
2	A	1179	N	OP2-P-OP1	2.77	119.50	110.44
3	H	389	NCA	C3-C7-N7	2.77	120.92	117.77
3	D	349	NCA	C5-C4-C3	2.76	123.91	120.32
3	B	429	NCA	C4-C3-C2	-2.76	114.25	117.64
3	F	369	NCA	C3-C7-N7	2.76	120.91	117.77
2	B	1179	N	O5'-P-OP1	2.75	114.78	106.71
3	E	459	NCA	C5-C4-C3	2.75	123.89	120.32
2	E	1179	N	OP2-P-OP1	2.71	119.29	110.44
2	G	1179	N	O2'-C2'-C3'	2.71	116.55	111.31
3	F	369	NCA	C4-C3-C2	-2.70	114.32	117.64
3	A	319	NCA	C4-C3-C2	-2.69	114.33	117.64
3	B	329	NCA	C5-C4-C3	2.69	123.81	120.32
3	E	359	NCA	C3-C7-N7	2.69	120.83	117.77
2	F	1179	N	O2'-C2'-C3'	2.67	116.47	111.31
3	B	429	NCA	C3-C7-N7	2.65	120.79	117.77
3	F	469	NCA	C5-C4-C3	2.65	123.76	120.32
2	D	1179	N	O4'-C4'-C5'	-2.63	103.55	109.57
3	A	419	NCA	C5-C4-C3	2.63	123.73	120.32
3	B	429	NCA	C5-C4-C3	2.62	123.72	120.32
3	C	339	NCA	C3-C7-N7	2.61	120.74	117.77
2	H	1179	N	OP2-P-OP1	2.60	118.94	110.44
2	E	1179	N	O2'-C2'-C3'	2.58	116.30	111.31
3	B	329	NCA	C4-C3-C2	-2.58	114.48	117.64
3	D	349	NCA	C3-C7-N7	2.49	120.61	117.77
2	D	1179	N	O2'-C2'-C3'	2.47	116.09	111.31
2	A	1179	N	O2'-C2'-C3'	2.47	116.08	111.31
2	A	1179	N	O5'-P-OP1	2.45	113.90	106.71
2	D	1179	N	O5'-P-OP1	2.45	113.88	106.71
2	C	1179	N	O5'-P-OP1	2.44	113.86	106.71
3	F	469	NCA	C4-C3-C2	-2.44	114.65	117.64
2	E	1179	N	O5'-P-OP1	2.42	113.81	106.71
2	H	1179	N	O5'-P-OP1	2.42	113.81	106.71
3	D	449	NCA	C3-C7-N7	2.42	120.52	117.77
3	G	479	NCA	O7-C7-N7	-2.41	119.10	122.59
3	E	359	NCA	C4-C3-C2	-2.41	114.68	117.64
3	B	429	NCA	C6-N1-C2	2.41	121.17	116.85
3	H	389	NCA	O7-C7-N7	-2.38	119.15	122.59
3	F	469	NCA	O7-C7-N7	-2.37	119.17	122.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	359	NCA	C5-C4-C3	2.36	123.38	120.32
3	C	339	NCA	C6-N1-C2	2.34	121.05	116.85
3	D	349	NCA	C6-N1-C2	2.33	121.03	116.85
2	F	1179	N	O5'-P-OP1	2.31	113.49	106.71
3	B	429	NCA	O7-C7-N7	-2.30	119.26	122.59
2	G	1179	N	O5'-P-OP1	2.30	113.46	106.71
2	G	1179	N	O4'-C4'-C5'	-2.30	104.30	109.57
3	C	439	NCA	C6-N1-C2	2.28	120.94	116.85
2	H	1179	N	O2'-C2'-C3'	2.27	115.70	111.31
3	H	389	NCA	C6-N1-C2	2.23	120.86	116.85
3	E	459	NCA	C4-C3-C2	-2.23	114.90	117.64
2	H	1179	N	C1'-C2'-C3'	2.23	105.17	101.65
2	C	1179	N	O2'-C2'-C3'	2.21	115.59	111.31
2	H	1179	N	C5'-C4'-C3'	2.14	123.77	115.21
3	B	329	NCA	C6-N1-C2	2.13	120.67	116.85
2	F	1179	N	C5'-C4'-C3'	2.09	123.57	115.21
3	F	369	NCA	C6-N1-C2	2.06	120.56	116.85
2	G	1179	N	C5'-C4'-C3'	2.06	123.47	115.21
3	H	489	NCA	C5-C4-C3	2.06	122.99	120.32
2	E	1179	N	C5'-C4'-C3'	2.06	123.45	115.21
3	A	319	NCA	C6-N1-C2	2.05	120.53	116.85
2	C	1179	N	C1'-C2'-C3'	2.05	104.89	101.65
2	C	1179	N	C5'-C4'-C3'	2.03	123.36	115.21
2	D	1179	N	C5'-C4'-C3'	2.03	123.35	115.21
2	B	1179	N	P-O5'-C5'	2.03	124.05	118.19
2	A	1179	N	C5'-C4'-C3'	2.03	123.33	115.21

All (8) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	1179	N	C3'
2	E	1179	N	C3'
2	B	1179	N	C3'
2	C	1179	N	C3'
2	H	1179	N	C3'
2	F	1179	N	C3'
2	A	1179	N	C3'
2	G	1179	N	C3'

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 ⓘ

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	251/258 (97%)	0.08	5 (1%) 62 59	21, 41, 63, 72	0
1	B	251/258 (97%)	-0.03	3 (1%) 75 75	15, 35, 55, 68	0
1	C	251/258 (97%)	0.04	4 (1%) 68 67	19, 39, 60, 73	0
1	D	251/258 (97%)	-0.09	2 (0%) 83 82	14, 34, 53, 75	0
1	E	251/258 (97%)	0.13	6 (2%) 56 54	23, 46, 63, 71	0
1	F	251/258 (97%)	-0.09	1 (0%) 90 90	14, 34, 54, 75	0
1	G	251/258 (97%)	-0.13	0 100 100	12, 35, 55, 78	0
1	H	251/258 (97%)	0.23	7 (2%) 50 48	26, 49, 63, 74	0
All	All	2008/2064 (97%)	0.02	28 (1%) 72 71	12, 39, 61, 78	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	128	PHE	7.1
1	A	128	PHE	5.7
1	D	251	ALA	5.1
1	B	251	ALA	3.7
1	B	1	ILE	3.3
1	D	250	LEU	2.9
1	H	117	ASN	2.8
1	E	134	GLN	2.6
1	C	115	ARG	2.6
1	H	128	PHE	2.6
1	H	127	ASP	2.5
1	E	122	GLU	2.4
1	H	134	GLN	2.4
1	E	126	PRO	2.4
1	F	115	ARG	2.4
1	H	232	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	47	PHE	2.3
1	B	250	LEU	2.3
1	A	135	ALA	2.3
1	C	222	HIS	2.2
1	E	31	ARG	2.2
1	H	31	ARG	2.2
1	A	250	LEU	2.2
1	A	126	PRO	2.2
1	E	127	ASP	2.1
1	A	132	PRO	2.1
1	E	46	SER	2.1
1	H	123	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	NCA	A	319	9/9	0.33	11.98	59,61,62,62	0
3	NCA	F	369	9/9	0.27	9.67	49,52,52,53	0
3	NCA	D	449	9/9	0.25	8.85	41,44,46,47	0
3	NCA	F	469	9/9	0.32	8.79	48,49,53,55	0
3	NCA	B	429	9/9	0.28	8.26	43,45,49,49	0
2	N	A	1179	13/13	0.31	7.97	69,71,85,86	0
3	NCA	D	349	9/9	0.35	7.25	67,67,68,68	0
3	NCA	G	379	9/9	0.26	5.99	61,62,62,63	0
3	NCA	H	489	9/9	0.29	5.58	48,49,54,54	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NCA	H	389	9/9	0.30	5.17	57,58,60,62	0
2	N	F	1179	13/13	0.23	5.01	59,61,68,68	0
2	N	H	1179	13/13	0.22	4.46	69,71,74,74	0
3	NCA	B	329	9/9	0.24	4.37	35,39,40,41	0
2	N	G	1179	13/13	0.22	4.11	59,60,66,66	0
3	NCA	A	419	9/9	0.42	3.71	58,59,60,61	0
3	NCA	C	339	9/9	0.24	3.65	53,54,54,55	0
3	NCA	E	459	9/9	0.25	3.21	52,53,55,55	0
3	NCA	G	479	9/9	0.26	2.88	38,40,44,44	0
3	NCA	C	439	9/9	0.24	2.82	52,54,55,56	0
2	N	B	1179	13/13	0.25	2.82	58,61,66,67	0
2	N	C	1179	13/13	0.23	2.64	71,74,82,84	0
3	NCA	E	359	9/9	0.22	2.51	54,54,55,55	0
2	N	E	1179	13/13	0.20	2.42	69,72,79,80	0
2	N	D	1179	13/13	0.20	1.75	53,54,68,68	0

6.5 Other polymers ⓘ

There are no such residues in this entry.