



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:28 AM UTC

PDB ID : 4F3Z / pdb_00004f3z
Title : Crystal structure of a swine H1N2 influenza virus hemagglutinin
Authors : Xu, R.; Wilson, I.A.
Deposited on : 2012-05-09
Resolution : 3.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

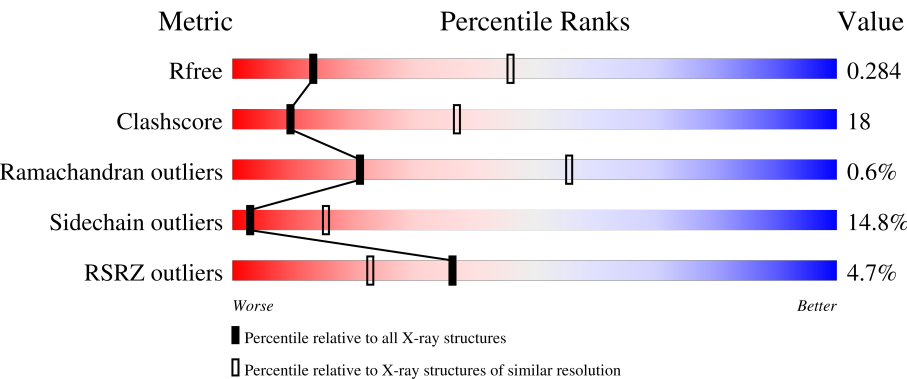
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	329	3% 52%36%8% . .
1	C	329	2% 52%32%7%9%
1	E	329	7% 48%41%9% .
2	B	179	4% 63%30%. .
2	D	179	5% 64%26%6%. .

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Mol	Chain	Length	Quality of chain
2	F	179	6% 53% 37% 6% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11544 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 Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2501	1575	434	479	13			
1	C	301	Total	C	N	O	S	0	0	0
			2364	1492	406	454	12			
1	E	320	Total	C	N	O	S	0	0	0
			2503	1576	434	480	13			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	PRO	-	expression tag	UNP Q8QT89
A	10	GLY	-	expression tag	UNP Q8QT89
A	205	CYS	GLY	engineered mutation	UNP Q8QT89
A	220	CYS	ARG	engineered mutation	UNP Q8QT89
C	9	PRO	-	expression tag	UNP Q8QT89
C	10	GLY	-	expression tag	UNP Q8QT89
C	205	CYS	GLY	engineered mutation	UNP Q8QT89
C	220	CYS	ARG	engineered mutation	UNP Q8QT89
E	9	PRO	-	expression tag	UNP Q8QT89
E	10	GLY	-	expression tag	UNP Q8QT89
E	205	CYS	GLY	engineered mutation	UNP Q8QT89
E	220	CYS	ARG	engineered mutation	UNP Q8QT89

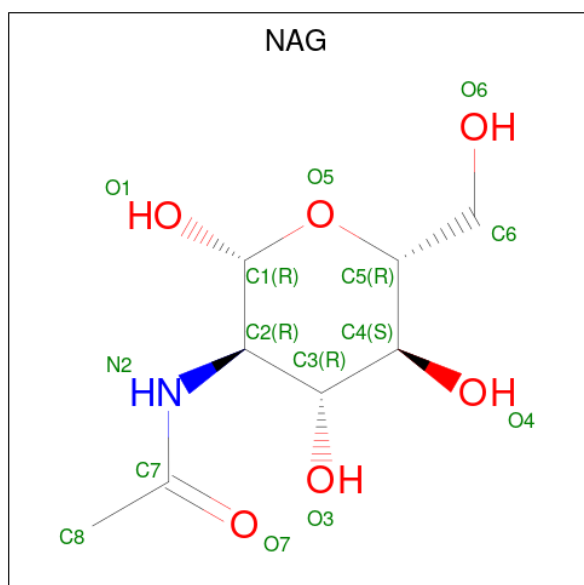
- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	171	Total	C	N	O	S	0	0	0
			1382	869	236	271	6			
2	D	174	Total	C	N	O	S	0	0	0
			1407	884	239	278	6			
2	F	170	Total	C	N	O	S	0	0	0
			1373	864	235	268	6			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	177	SER	-	expression tag	UNP Q8QT89
B	178	GLY	-	expression tag	UNP Q8QT89
B	179	ARG	-	expression tag	UNP Q8QT89
D	177	SER	-	expression tag	UNP Q8QT89
D	178	GLY	-	expression tag	UNP Q8QT89
D	179	ARG	-	expression tag	UNP Q8QT89
F	177	SER	-	expression tag	UNP Q8QT89
F	178	GLY	-	expression tag	UNP Q8QT89
F	179	ARG	-	expression tag	UNP Q8QT89

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

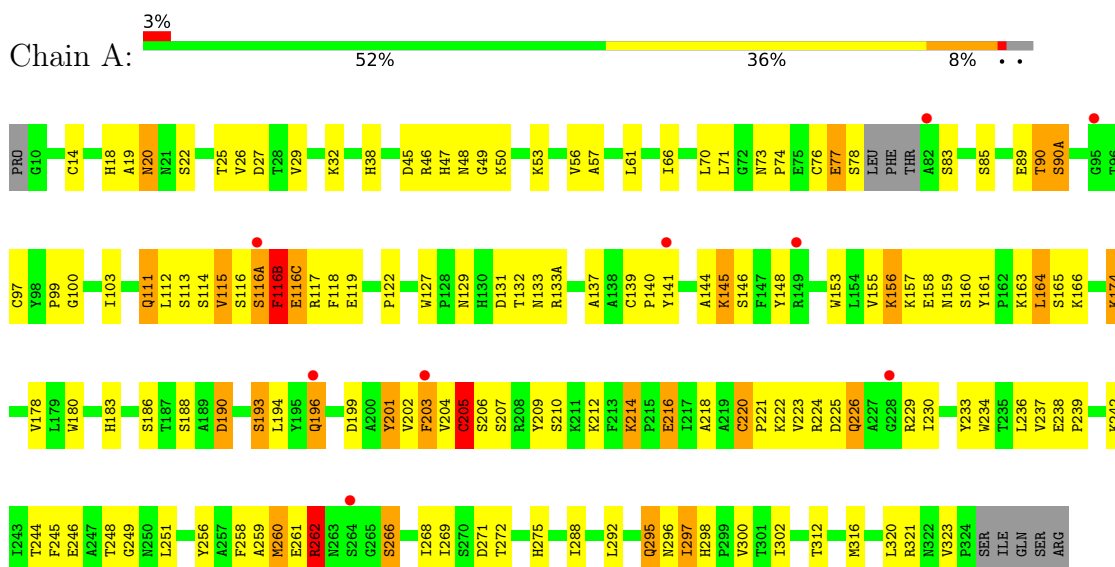


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			14	8	1	5		

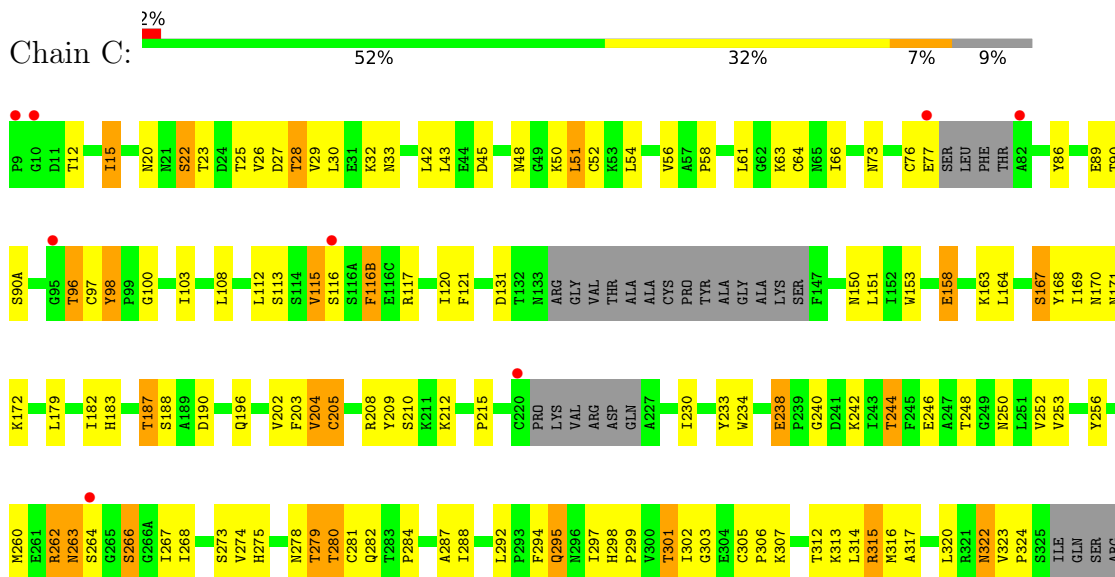
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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.

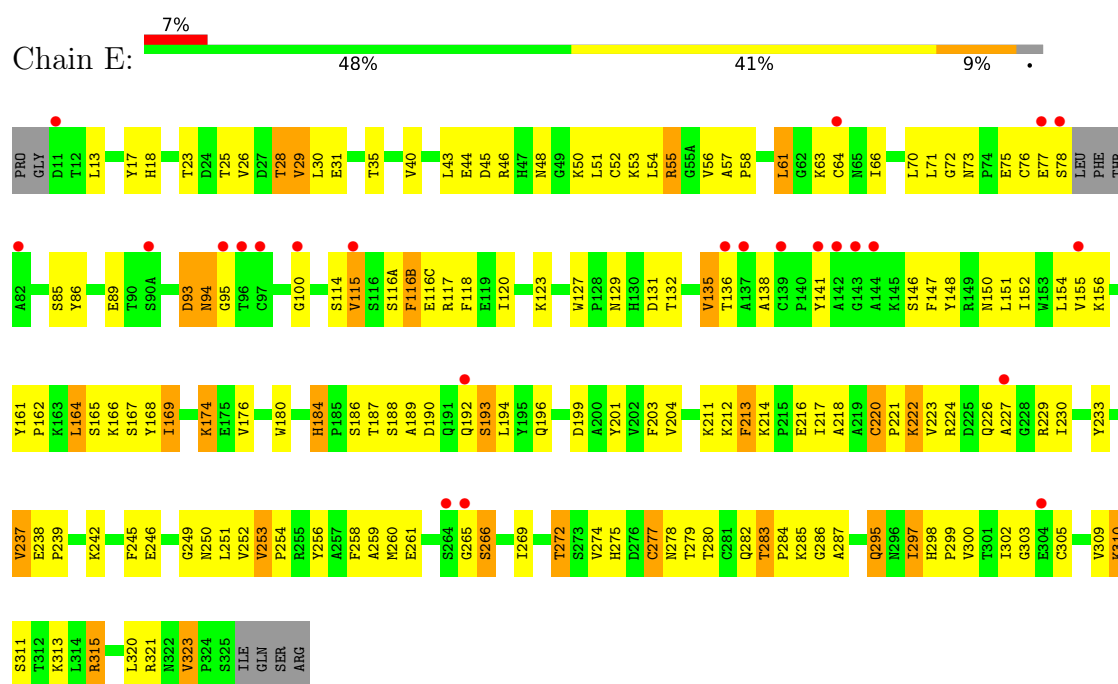
• Molecule 1: Hemagglutinin



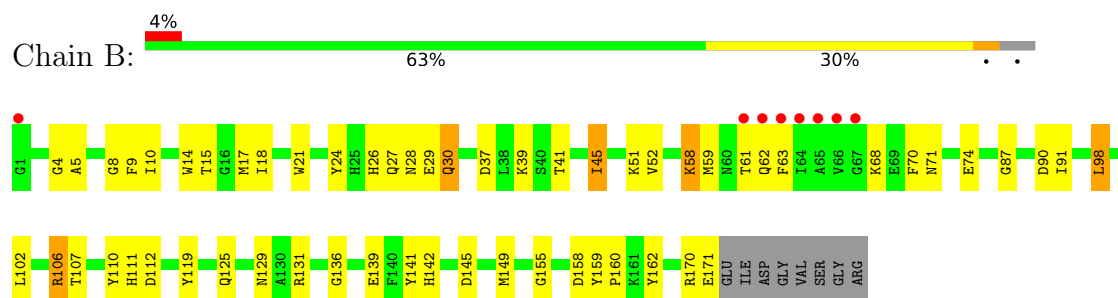
• Molecule 1: Hemagglutinin



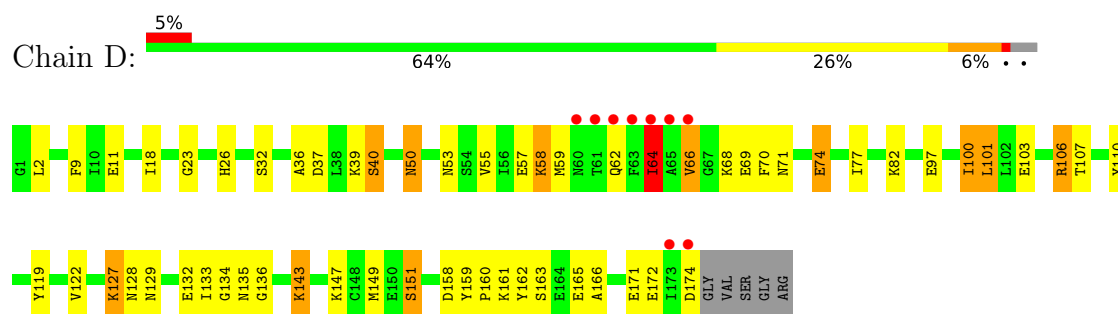
• Molecule 1: Hemagglutinin



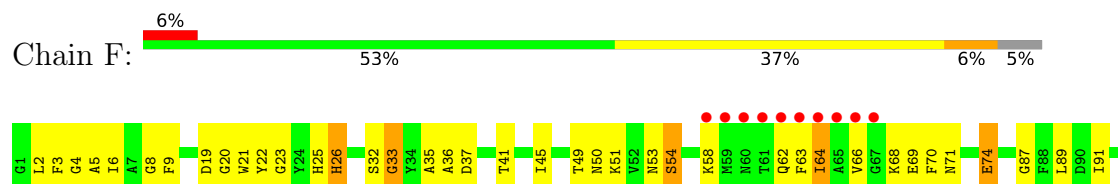
• Molecule 2: Hemagglutinin

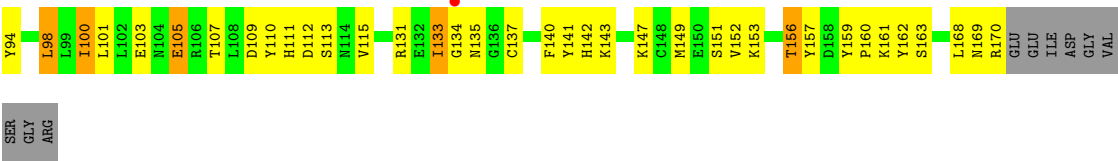


• Molecule 2: Hemagglutinin



• Molecule 2: Hemagglutinin





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.44Å 94.92Å 220.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.27 – 3.20 34.27 – 3.20	Depositor EDS
% Data completeness (in resolution range)	93.2 (34.27-3.20) 92.5 (34.27-3.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.73 (at 3.18Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.234 , 0.293 0.227 , 0.284	Depositor DCC
R_{free} test set	1575 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 33.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.013 for k,h,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	11544	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 2.81% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: NAG

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	0.28	0/2562	0.78	7/3482 (0.2%)
1	C	0.30	0/2421	0.79	6/3288 (0.2%)
1	E	0.29	0/2564	0.81	6/3485 (0.2%)
2	B	0.28	0/1410	0.68	0/1897
2	D	0.30	0/1435	0.69	1/1931 (0.1%)
2	F	0.31	0/1401	0.72	3/1885 (0.2%)
All	All	0.29	0/11793	0.76	23/15968 (0.1%)

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	A	0	2
1	C	0	2
1	E	0	2
2	D	0	1
All	All	0	7

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	98	TYR	CA-C-N	7.36	127.03	119.82
1	C	98	TYR	C-N-CA	7.36	127.03	119.82
1	E	323	VAL	CA-C-N	7.20	126.70	118.85
1	E	323	VAL	C-N-CA	7.20	126.70	118.85
1	E	184	HIS	CA-C-N	6.62	126.58	120.03
1	E	184	HIS	C-N-CA	6.62	126.58	120.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	33	GLY	N-CA-C	6.40	119.86	110.38
1	A	158	GLU	CB-CA-C	-6.36	109.26	116.63
1	E	305	CYS	CA-C-N	6.10	126.06	120.03
1	E	305	CYS	C-N-CA	6.10	126.06	120.03
2	D	127	LYS	CB-CA-C	-5.98	109.70	116.63
2	F	26	HIS	N-CA-C	5.74	117.58	108.79
1	C	121	PHE	CA-C-N	5.69	125.71	119.90
1	C	121	PHE	C-N-CA	5.69	125.71	119.90
1	A	127	TRP	CA-C-N	5.61	124.80	118.97
1	A	127	TRP	C-N-CA	5.61	124.80	118.97
1	C	158	GLU	CB-CA-C	-5.28	110.06	117.23
1	A	292	LEU	CA-C-N	5.17	124.79	119.56
1	A	292	LEU	C-N-CA	5.17	124.79	119.56
1	C	263	ASN	N-CA-C	-5.12	102.44	110.17
1	A	262	ARG	N-CA-C	5.08	116.97	107.99
1	A	20	ASN	N-CA-C	5.05	114.81	108.45
2	F	64	ILE	N-CA-C	5.01	115.49	108.17

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	205	CYS	Peptide
1	A	262	ARG	Peptide
1	C	262	ARG	Peptide
1	C	96	THR	Peptide
2	D	64	ILE	Peptide
1	E	94	ASN	Peptide
1	E	95	GLY	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	2501	0	2429	105	0
1	C	2364	0	2288	85	0
1	E	2503	0	2431	125	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1382	0	1312	41	0
2	D	1407	0	1333	37	0
2	F	1373	0	1306	48	0
3	C	14	0	13	0	0
All	All	11544	0	11112	404	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 18.

All (404) 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:E:189:ALA:O	1:E:193:SER:HB3	1.50	1.11
1:C:96:THR:HA	1:C:97:CYS:HB2	1.37	1.01
1:E:116(A):SER:HB3	1:E:261:GLU:HB3	1.39	1.00
1:E:141:TYR:HB2	1:E:146:SER:HB2	1.46	0.97
1:A:70:LEU:HD11	1:A:112:LEU:HD11	1.53	0.91
1:A:161:TYR:H	1:A:196:GLN:HE21	1.15	0.90
1:A:155:VAL:HG22	1:A:156:LYS:H	1.37	0.90
1:E:155:VAL:HG12	1:E:156:LYS:H	1.36	0.89
1:E:230:ILE:HD11	1:E:252:VAL:HG21	1.52	0.89
1:C:305:CYS:HB2	1:C:306:PRO:CD	2.03	0.87
1:C:170:ASN:HB3	1:C:240:GLY:H	1.42	0.84
1:E:274:VAL:HG22	1:E:275:HIS:H	1.44	0.82
2:F:70:PHE:HB3	2:F:74:GLU:HB2	1.60	0.81
1:E:51:LEU:HB3	1:E:274:VAL:HA	1.62	0.81
2:D:26:HIS:HB2	2:D:149:MET:HE3	1.61	0.80
1:A:205:CYS:HB2	1:E:221:PRO:HD2	1.64	0.80
2:F:26:HIS:O	2:F:32:SER:HA	1.82	0.80
1:E:61:LEU:HD11	1:E:89:GLU:HA	1.65	0.79
1:A:116(C):GLU:HG3	1:A:259:ALA:HB3	1.64	0.78
1:E:169:ILE:HG12	1:E:242:LYS:HG3	1.65	0.78
1:E:116(C):GLU:HG2	1:E:259:ALA:HB3	1.66	0.78
2:D:70:PHE:HB3	2:D:74:GLU:HB3	1.65	0.78
1:E:138:ALA:HB1	1:E:224:ARG:HB3	1.63	0.78
1:E:220:CYS:HB2	1:E:221:PRO:CD	2.14	0.78
1:C:316:MET:HE1	2:D:55:VAL:HG21	1.66	0.77
1:A:47:HIS:HB3	1:A:297:ILE:HD12	1.67	0.75
1:A:316:MET:HE1	2:B:51:LYS:HG2	1.68	0.75
1:A:100:GLY:HA3	1:A:230:ILE:O	1.87	0.75
1:E:220:CYS:HB2	1:E:221:PRO:HD2	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:GLN:HG2	1:A:297:ILE:H	1.50	0.74
1:C:117:ARG:NH1	1:C:150:ASN:HD21	1.87	0.73
1:E:129:ASN:HB3	1:E:162:PRO:HG2	1.72	0.72
1:E:222:LYS:HG3	1:E:227:ALA:HB2	1.71	0.72
1:E:26:VAL:HG12	1:E:315:ARG:HG2	1.72	0.71
1:E:73:ASN:HB3	1:E:76:CYS:SG	2.30	0.71
2:B:28:ASN:HD21	2:B:145:ASP:C	1.98	0.71
1:C:52:CYS:HB2	1:C:279:THR:HG23	1.72	0.71
1:E:28:THR:HG22	1:E:31:GLU:H	1.55	0.71
2:D:39:LYS:N	2:D:39:LYS:HD2	2.04	0.70
1:E:54:LEU:HG	1:E:55:ARG:HG3	1.73	0.70
1:E:51:LEU:HB2	1:E:272:THR:HG23	1.72	0.70
1:A:261:GLU:HG2	1:A:262:ARG:H	1.56	0.70
1:C:100:GLY:HA3	1:C:230:ILE:O	1.93	0.68
1:C:20:ASN:OD1	1:C:22:SER:HB3	1.93	0.68
1:C:96:THR:HA	1:C:97:CYS:CB	2.20	0.68
1:A:164:LEU:O	1:A:246:GLU:HA	1.94	0.67
1:C:204:VAL:HA	1:C:244:THR:O	1.95	0.66
1:E:56:VAL:HG22	1:E:57:ALA:H	1.59	0.66
1:E:155:VAL:HG12	1:E:156:LYS:N	2.08	0.66
1:A:83:SER:O	1:A:116(B):PHE:HB3	1.95	0.66
1:C:305:CYS:HB2	1:C:306:PRO:HD2	1.75	0.66
1:A:99:PRO:HB3	1:A:223:VAL:HB	1.78	0.66
1:A:212:LYS:HD3	1:E:216:GLU:HB3	1.78	0.65
2:F:25:HIS:HA	2:F:33:GLY:O	1.96	0.65
1:E:222:LYS:H	1:E:222:LYS:HD2	1.61	0.65
1:A:131:ASP:HB3	1:A:155:VAL:HG12	1.77	0.65
1:E:100:GLY:HA3	1:E:230:ILE:O	1.96	0.65
2:B:119:TYR:CE1	2:B:136:GLY:HA2	2.32	0.64
2:B:26:HIS:HB2	2:B:149:MET:HE3	1.78	0.64
1:A:206:SER:HA	1:A:242:LYS:O	1.97	0.64
1:A:205:CYS:HA	1:A:209:TYR:O	1.98	0.63
1:A:116(A):SER:O	1:A:116(B):PHE:HB2	1.97	0.63
2:D:171:GLU:HA	2:D:174:ASP:HB3	1.81	0.63
1:E:71:LEU:O	1:E:148:TYR:HB3	1.99	0.63
1:C:52:CYS:HB2	1:C:279:THR:CG2	2.29	0.62
2:F:41:THR:O	2:F:45:ILE:HG12	1.98	0.62
2:F:141:TYR:CE1	2:F:170:ARG:HG3	2.35	0.62
2:B:24:TYR:HE2	2:B:37:ASP:HB2	1.65	0.62
2:B:58:LYS:HG2	2:F:98:LEU:HD23	1.80	0.62
2:D:64:ILE:HD13	2:D:66:VAL:HG22	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:125:GLN:HE22	2:B:155:GLY:HA2	1.66	0.61
2:B:30:GLN:HE21	2:B:145:ASP:HB2	1.66	0.61
1:A:202:VAL:HG11	1:A:251:LEU:HD13	1.82	0.61
2:D:70:PHE:HD1	2:D:74:GLU:HG2	1.66	0.61
1:A:155:VAL:HG22	1:A:156:LYS:N	2.14	0.61
2:D:23:GLY:HA3	2:D:36:ALA:HA	1.82	0.61
1:E:164:LEU:O	1:E:246:GLU:HA	2.00	0.61
2:F:71:ASN:N	2:F:74:GLU:HG3	2.16	0.61
1:C:96:THR:CA	1:C:97:CYS:HB2	2.23	0.60
2:B:142:HIS:CD2	2:B:162:TYR:HB3	2.36	0.60
1:E:283:THR:HG23	1:E:285:LYS:H	1.67	0.60
1:A:26:VAL:HG22	1:A:27:ASP:H	1.65	0.60
1:E:253:VAL:HG22	1:E:254:PRO:HD2	1.83	0.60
1:A:216:GLU:HG2	1:C:212:LYS:HE2	1.84	0.59
1:E:61:LEU:HD13	1:E:64:CYS:O	2.02	0.59
2:D:9:PHE:O	2:D:135:ASN:HA	2.03	0.59
1:A:114:SER:O	1:A:115:VAL:HB	2.02	0.59
2:B:70:PHE:HB3	2:B:74:GLU:HB2	1.83	0.59
1:A:56:VAL:HB	1:A:85:SER:HB3	1.84	0.59
1:E:18:HIS:HB2	2:F:20:GLY:O	2.03	0.59
2:F:71:ASN:H	2:F:74:GLU:HG3	1.68	0.59
1:C:268:ILE:HG23	1:C:284:PRO:HG3	1.86	0.58
1:E:211:LYS:HG3	1:E:212:LYS:N	2.17	0.58
2:F:107:THR:O	2:F:110:TYR:HB3	2.03	0.58
2:D:71:ASN:OD1	2:D:74:GLU:HB2	2.03	0.58
1:C:51:LEU:HA	1:C:282:GLN:NE2	2.17	0.58
1:E:44:GLU:OE1	1:E:46:ARG:HB3	2.04	0.58
1:E:75:GLU:HB2	1:E:94:ASN:HD22	1.68	0.58
2:F:133:ILE:HG23	2:F:137:CYS:O	2.04	0.57
2:D:158:ASP:OD2	2:D:161:LYS:HB2	2.04	0.57
1:C:305:CYS:HB2	1:C:306:PRO:HD3	1.83	0.57
1:A:118:PHE:O	1:A:256:TYR:HA	2.04	0.57
1:A:47:HIS:CE1	1:A:49:GLY:HA2	2.40	0.57
1:A:115:VAL:HG23	1:A:262:ARG:HH12	1.69	0.57
1:E:222:LYS:HB3	1:E:226:GLN:C	2.29	0.57
1:C:117:ARG:HH12	1:C:150:ASN:HD21	1.53	0.56
1:E:192:GLN:O	1:E:196:GLN:HA	2.05	0.56
2:B:71:ASN:H	2:B:74:GLU:HG3	1.70	0.56
1:C:15:ILE:HD11	2:D:122:VAL:HG21	1.86	0.56
1:C:205:CYS:HB3	1:C:209:TYR:O	2.04	0.56
1:A:214:LYS:HZ2	1:A:216:GLU:HG3	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:TYR:CE1	1:A:248:THR:HG23	2.41	0.56
1:A:220:CYS:HB2	1:A:221:PRO:HD2	1.87	0.56
2:B:90:ASP:OD2	2:D:62:GLN:HB2	2.04	0.56
1:C:215:PRO:HB3	1:C:250:ASN:ND2	2.19	0.56
1:E:295:GLN:HG3	1:E:297:ILE:HB	1.87	0.56
1:A:238:GLU:HB3	1:A:239:PRO:HD2	1.86	0.56
2:F:19:ASP:HB3	2:F:36:ALA:HB3	1.88	0.56
1:A:20:ASN:OD1	1:A:22:SER:HB3	2.06	0.56
1:A:204:VAL:HA	1:A:244:THR:O	2.06	0.56
1:E:58:PRO:HB3	1:E:86:TYR:CZ	2.42	0.55
1:E:72:GLY:HA3	1:E:150:ASN:ND2	2.21	0.55
1:E:298:HIS:CE1	1:E:300:VAL:HB	2.41	0.55
2:B:91:ILE:HG21	2:F:91:ILE:HD13	1.87	0.55
2:D:74:GLU:HG3	2:D:77:ILE:HD11	1.89	0.55
1:C:279:THR:HG21	1:C:287:ALA:CB	2.35	0.55
1:E:155:VAL:CG1	1:E:156:LYS:H	2.16	0.55
1:E:28:THR:OG1	2:F:105:GLU:HG2	2.06	0.55
1:A:18:HIS:ND1	1:A:19:ALA:N	2.56	0.54
2:B:41:THR:O	2:B:45:ILE:HG13	2.06	0.54
1:E:131:ASP:HB3	1:E:155:VAL:HB	1.89	0.54
1:E:278:ASN:O	1:E:279:THR:HB	2.07	0.54
2:F:23:GLY:HA2	2:F:37:ASP:H	1.71	0.54
1:A:137:ALA:HA	1:A:145:LYS:HG3	1.89	0.54
1:E:53:LYS:HB2	1:E:277:CYS:O	2.07	0.54
1:E:302:ILE:HD12	1:E:303:GLY:H	1.73	0.54
1:C:295:GLN:HG2	1:C:297:ILE:H	1.73	0.54
2:F:50:ASN:O	2:F:54:SER:HB2	2.09	0.53
1:E:28:THR:HG22	1:E:31:GLU:N	2.23	0.53
1:E:211:LYS:HG2	1:E:213:PHE:CE1	2.43	0.53
1:A:139:CYS:HB2	1:A:146:SER:O	2.08	0.53
1:C:295:GLN:HB3	1:C:306:PRO:HB2	1.90	0.53
1:A:201:TYR:HE1	1:A:248:THR:HG23	1.73	0.53
2:B:87:GLY:O	2:B:91:ILE:HG12	2.08	0.53
1:C:48:ASN:O	1:C:50:LYS:HG3	2.08	0.53
1:C:294:PHE:CE1	2:D:59:MET:HE2	2.43	0.53
1:A:174:LYS:HD2	1:A:261:GLU:HB2	1.90	0.53
1:A:183:HIS:HD2	1:A:229:ARG:O	1.91	0.53
1:C:164:LEU:O	1:C:246:GLU:HA	2.09	0.53
2:D:64:ILE:HD12	2:D:64:ILE:N	2.24	0.53
1:E:222:LYS:H	1:E:222:LYS:CD	2.21	0.52
1:A:50:LYS:HD3	1:A:275:HIS:ND1	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LEU:HD12	1:A:89:GLU:HG2	1.92	0.52
1:A:111:GLN:HG2	1:A:236:LEU:HD11	1.91	0.52
1:A:47:HIS:HE1	1:A:49:GLY:HA2	1.74	0.52
1:C:120:ILE:HG13	1:C:168:TYR:CD1	2.44	0.52
1:E:298:HIS:CG	1:E:299:PRO:HD2	2.45	0.52
1:A:268:ILE:HG13	1:A:302:ILE:HD11	1.90	0.52
1:C:96:THR:OG1	1:C:98:TYR:N	2.43	0.52
1:A:221:PRO:HB3	1:C:242:LYS:HE3	1.91	0.52
1:C:25:THR:HG22	1:C:33:ASN:HA	1.92	0.52
1:C:208:ARG:HH11	1:C:238:GLU:HG2	1.73	0.52
2:D:74:GLU:CG	2:D:77:ILE:HD11	2.39	0.52
2:F:49:THR:O	2:F:53:ASN:HB2	2.10	0.52
1:A:71:LEU:O	1:A:148:TYR:HB3	2.10	0.51
2:B:51:LYS:HE2	1:E:29:VAL:HG21	1.92	0.51
2:F:111:HIS:O	2:F:115:VAL:HG23	2.09	0.51
1:A:262:ARG:HA	1:A:262:ARG:HH11	1.75	0.51
1:A:320:LEU:HD23	1:A:320:LEU:H	1.74	0.51
1:C:115:VAL:HG13	1:C:262:ARG:HH11	1.76	0.51
1:A:145:LYS:HD2	1:A:145:LYS:N	2.25	0.51
2:F:4:GLY:O	2:F:8:GLY:HA3	2.11	0.51
2:B:159:TYR:HB3	2:B:160:PRO:HD3	1.93	0.51
1:E:89:GLU:O	1:E:269:ILE:HA	2.10	0.51
1:E:167:SER:HB2	1:E:242:LYS:HD3	1.91	0.51
1:E:48:ASN:HB3	1:E:287:ALA:HB3	1.93	0.51
1:C:51:LEU:HD12	1:C:282:GLN:NE2	2.25	0.51
2:F:23:GLY:HA3	2:F:36:ALA:HA	1.93	0.51
1:E:155:VAL:HG12	1:E:156:LYS:HG2	1.92	0.50
1:E:274:VAL:HG22	1:E:275:HIS:N	2.20	0.50
1:C:51:LEU:HD12	1:C:282:GLN:HE21	1.75	0.50
1:C:187:THR:HG23	1:C:190:ASP:CG	2.37	0.50
2:D:97:GLU:O	2:D:100:ILE:HG13	2.11	0.50
2:F:21:TRP:CZ3	2:F:45:ILE:HD13	2.47	0.50
1:C:150:ASN:HD22	1:C:256:TYR:HD1	1.58	0.50
1:C:153:TRP:HZ2	1:C:183:HIS:ND1	2.09	0.50
1:E:155:VAL:HG13	1:E:194:LEU:HD12	1.93	0.50
2:B:24:TYR:CE2	2:B:37:ASP:HB2	2.45	0.50
1:E:127:TRP:HE1	1:E:152:ILE:HD12	1.77	0.50
1:E:116(A):SER:HB3	1:E:261:GLU:CB	2.27	0.50
1:E:135:VAL:HG12	1:E:147:PHE:HB2	1.93	0.50
1:A:161:TYR:HB3	1:A:196:GLN:HG2	1.94	0.49
2:B:170:ARG:O	2:B:171:GLU:HG2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:52:CYS:HB2	1:E:282:GLN:HE22	1.77	0.49
1:A:153:TRP:HZ2	1:A:183:HIS:ND1	2.10	0.49
1:A:262:ARG:HA	1:A:262:ARG:NH1	2.27	0.49
2:D:159:TYR:HB3	2:D:160:PRO:HD3	1.93	0.49
1:E:203:PHE:HD1	1:E:212:LYS:HB2	1.76	0.49
2:F:151:SER:HA	2:F:156:THR:HG22	1.94	0.49
1:A:38:HIS:CE1	2:B:21:TRP:HE1	2.30	0.49
1:C:26:VAL:HG12	1:C:315:ARG:HG2	1.93	0.49
1:A:298:HIS:CE1	1:A:300:VAL:HB	2.47	0.49
1:E:187:THR:HG23	1:E:190:ASP:H	1.77	0.49
1:A:203:PHE:CE1	1:A:205:CYS:SG	3.06	0.49
1:E:13:LEU:HD12	2:F:25:HIS:O	2.13	0.49
1:A:97:CYS:N	1:A:148:TYR:OH	2.45	0.49
1:C:298:HIS:CG	1:C:299:PRO:HD2	2.47	0.49
1:C:322:ASN:HD22	1:C:324:PRO:HD3	1.78	0.49
2:F:160:PRO:HA	2:F:163:SER:HB2	1.95	0.49
1:A:116(C):GLU:CG	1:A:259:ALA:HB3	2.36	0.49
1:E:184:HIS:HB2	1:E:229:ARG:HB3	1.95	0.49
2:F:159:TYR:N	2:F:160:PRO:HD2	2.28	0.49
1:A:156:LYS:HD3	1:A:159:ASN:HA	1.94	0.49
1:E:48:ASN:HB3	1:E:287:ALA:H	1.78	0.49
1:C:108:LEU:HB2	1:C:234:TRP:CE2	2.48	0.48
1:E:118:PHE:O	1:E:256:TYR:HA	2.13	0.48
2:D:127:LYS:HB3	2:D:128:ASN:H	1.48	0.48
1:E:57:ALA:HB1	1:E:58:PRO:HD2	1.95	0.48
2:D:161:LYS:HD3	2:D:162:TYR:CZ	2.48	0.48
1:E:265:GLY:O	1:E:266:SER:HB2	2.13	0.48
1:A:160:SER:HA	1:A:196:GLN:NE2	2.29	0.48
1:A:260:MET:HE2	1:A:260:MET:HB3	1.73	0.48
2:B:9:PHE:CE1	2:B:10:ILE:HG13	2.49	0.48
2:D:119:TYR:CE1	2:D:136:GLY:HA2	2.48	0.48
1:C:97:CYS:HB3	1:C:98:TYR:HD2	1.78	0.48
1:A:178:VAL:O	1:A:234:TRP:HA	2.14	0.48
1:C:112:LEU:O	1:C:115:VAL:HG23	2.14	0.48
1:E:52:CYS:HB2	1:E:282:GLN:NE2	2.29	0.47
1:A:161:TYR:H	1:A:196:GLN:NE2	1.97	0.47
1:A:190:ASP:O	1:A:194:LEU:HD12	2.14	0.47
1:E:180:TRP:CE2	1:E:233:TYR:HB2	2.49	0.47
1:A:53:LYS:HE2	1:A:57:ALA:HB2	1.96	0.47
1:C:113:SER:O	1:C:266:SER:HB3	2.13	0.47
2:F:94:TYR:O	2:F:98:LEU:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:GLU:CD	1:A:122:PRO:HA	2.39	0.47
1:C:116:SER:HB3	1:C:263:ASN:HB2	1.96	0.47
1:C:163:LYS:HZ1	1:C:246:GLU:CD	2.22	0.47
2:B:98:LEU:HD23	2:D:58:LYS:HD3	1.96	0.47
1:E:56:VAL:CG1	1:E:85:SER:HB3	2.45	0.47
1:E:161:TYR:CZ	1:E:249:GLY:HA2	2.49	0.47
1:E:282:GLN:HE22	1:E:287:ALA:HB2	1.80	0.47
1:E:117:ARG:HB3	1:E:258:PHE:CD1	2.50	0.47
1:E:251:LEU:HD12	1:E:252:VAL:N	2.29	0.47
1:E:93:ASP:O	1:E:94:ASN:C	2.57	0.47
2:F:140:PHE:HB3	2:F:142:HIS:O	2.15	0.47
1:E:309:VAL:HG12	1:E:311:SER:H	1.80	0.47
1:A:316:MET:SD	2:B:52:VAL:HG22	2.54	0.46
2:B:28:ASN:HD21	2:B:145:ASP:CA	2.28	0.46
1:E:56:VAL:HG13	1:E:85:SER:HB3	1.97	0.46
1:E:77:GLU:O	1:E:78:SER:HB3	2.14	0.46
1:C:279:THR:HG21	1:C:287:ALA:HB1	1.98	0.46
2:F:62:GLN:O	2:F:64:ILE:HG13	2.15	0.46
2:F:98:LEU:HD22	2:F:98:LEU:HA	1.79	0.46
1:C:58:PRO:HB3	1:C:86:TYR:CE2	2.51	0.46
2:D:147:LYS:O	2:D:151:SER:HB3	2.14	0.46
1:E:56:VAL:HG22	1:E:57:ALA:N	2.29	0.46
1:C:116(B):PHE:HE1	1:C:260:MET:HE3	1.81	0.46
1:E:116(B):PHE:HA	1:E:259:ALA:O	2.15	0.46
1:C:168:TYR:O	1:C:242:LYS:HA	2.16	0.46
2:B:51:LYS:HA	1:E:29:VAL:HG23	1.97	0.46
1:E:238:GLU:HB3	1:E:239:PRO:HD2	1.98	0.46
1:C:292:LEU:O	1:C:306:PRO:HG3	2.16	0.45
1:E:120:ILE:HD11	1:E:168:TYR:CB	2.46	0.45
1:E:249:GLY:O	1:E:250:ASN:HB2	2.16	0.45
1:C:61:LEU:HD12	1:C:89:GLU:HG2	1.98	0.45
1:A:116(A):SER:CB	1:A:261:GLU:H	2.29	0.45
1:A:242:LYS:HE3	1:E:221:PRO:HB3	1.97	0.45
1:E:284:PRO:HG2	1:E:298:HIS:CE1	2.52	0.45
1:A:116(B):PHE:CZ	1:A:258:PHE:CD1	3.05	0.45
1:A:222:LYS:HB3	1:A:225:ASP:HA	1.99	0.45
1:E:116(B):PHE:HD1	1:E:260:MET:HE3	1.80	0.45
2:F:152:VAL:HA	2:F:157:TYR:HB2	1.99	0.45
1:A:320:LEU:HB3	2:B:111:HIS:CG	2.52	0.45
1:C:29:VAL:CG1	2:F:51:LYS:HG3	2.47	0.45
1:A:161:TYR:CZ	1:A:249:GLY:HA2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:164:LEU:C	1:C:164:LEU:HD12	2.41	0.45
1:E:222:LYS:HA	1:E:227:ALA:HA	1.98	0.45
2:F:133:ILE:HG12	2:F:137:CYS:HB2	1.99	0.45
1:E:116(B):PHE:CD1	1:E:260:MET:HE3	2.52	0.45
1:A:174:LYS:HB2	1:A:260:MET:O	2.17	0.45
1:E:201:TYR:CD2	1:E:212:LYS:HE2	2.52	0.45
1:E:283:THR:HG22	1:E:286:GLY:H	1.82	0.45
2:B:5:ALA:HB3	2:B:112:ASP:OD1	2.16	0.44
1:A:161:TYR:N	1:A:196:GLN:HE21	1.97	0.44
1:A:113:SER:O	1:A:266:SER:HB3	2.18	0.44
1:A:140:PRO:HA	1:A:144:ALA:O	2.17	0.44
2:B:158:ASP:OD2	2:B:160:PRO:HD2	2.17	0.44
1:E:55:ARG:HH12	2:F:63:PHE:HZ	1.66	0.44
1:C:25:THR:CG2	1:C:33:ASN:HA	2.47	0.44
1:E:310:LYS:HD2	2:F:89:LEU:HG	2.00	0.44
1:A:242:LYS:HE3	1:E:221:PRO:CB	2.48	0.44
2:B:159:TYR:N	2:B:160:PRO:CD	2.80	0.44
1:C:27:ASP:OD1	1:C:32:LYS:HE3	2.18	0.44
1:C:281:CYS:HA	1:C:302:ILE:O	2.18	0.44
2:D:159:TYR:N	2:D:160:PRO:CD	2.81	0.44
1:E:211:LYS:HG2	1:E:213:PHE:CZ	2.52	0.44
1:A:244:THR:OG1	1:E:221:PRO:HD3	2.18	0.44
1:E:156:LYS:HD3	1:E:194:LEU:O	2.18	0.44
2:F:19:ASP:HB3	2:F:36:ALA:CB	2.47	0.44
1:C:170:ASN:CB	1:C:240:GLY:H	2.22	0.44
2:D:2:LEU:HG	2:F:3:PHE:CZ	2.53	0.44
1:E:116(C):GLU:N	1:E:116(C):GLU:OE1	2.50	0.44
1:A:165:SER:HA	1:A:245:PHE:O	2.18	0.43
1:A:242:LYS:NZ	1:E:222:LYS:HE3	2.33	0.43
2:F:5:ALA:HB3	2:F:112:ASP:OD2	2.18	0.43
1:A:156:LYS:HZ3	1:A:193:SER:C	2.27	0.43
1:C:208:ARG:HD2	1:C:238:GLU:CG	2.47	0.43
2:D:129:ASN:HA	2:D:166:ALA:HB1	2.00	0.43
1:E:320:LEU:HD12	1:E:321:ARG:O	2.19	0.43
1:A:180:TRP:CE2	1:A:233:TYR:HB2	2.53	0.43
2:B:107:THR:O	2:B:110:TYR:HB3	2.17	0.43
2:D:26:HIS:O	2:D:32:SER:HB2	2.18	0.43
1:A:224:ARG:O	1:A:226:GLN:HG2	2.18	0.43
1:C:313:LYS:HD3	1:C:314:LEU:N	2.33	0.43
1:E:135:VAL:HG12	1:E:146:SER:C	2.43	0.43
1:E:151:LEU:HA	1:E:253:VAL:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:42:LEU:HA	1:C:292:LEU:HD22	2.01	0.43
1:E:217:ILE:O	1:E:218:ALA:HB2	2.19	0.43
1:A:165:SER:OG	1:A:246:GLU:HG3	2.18	0.43
1:A:218:ALA:HB1	1:C:203:PHE:CZ	2.54	0.43
1:C:263:ASN:OD1	1:C:264:SER:N	2.50	0.43
1:A:73:ASN:HA	1:A:74:PRO:HD3	1.88	0.43
1:A:141:TYR:HB2	1:A:146:SER:CB	2.49	0.43
1:A:261:GLU:HG2	1:A:262:ARG:N	2.28	0.43
2:D:53:ASN:O	2:D:57:GLU:HG3	2.19	0.43
1:E:127:TRP:CD1	1:E:154:LEU:HD21	2.53	0.43
1:E:212:LYS:HD3	1:E:214:LYS:HE2	2.01	0.43
2:F:149:MET:O	2:F:153:LYS:HG3	2.19	0.43
1:A:32:LYS:HD3	2:D:50:ASN:OD1	2.18	0.42
1:A:116(A):SER:HB3	1:A:261:GLU:H	1.83	0.42
1:A:221:PRO:HG3	1:C:242:LYS:O	2.19	0.42
2:B:39:LYS:HB2	2:B:39:LYS:HE3	1.88	0.42
1:C:43:LEU:HD22	1:C:314:LEU:HG	2.00	0.42
2:D:143:LYS:HD2	2:D:143:LYS:HA	1.76	0.42
1:C:282:GLN:O	1:C:301:THR:HA	2.19	0.42
1:E:165:SER:HA	1:E:245:PHE:O	2.20	0.42
1:C:151:LEU:HB3	1:C:252:VAL:HG12	2.00	0.42
1:C:179:LEU:HD23	1:C:234:TRP:HB3	2.01	0.42
1:C:182:ILE:O	1:C:230:ILE:HA	2.20	0.42
1:E:174:LYS:HB3	1:E:260:MET:O	2.19	0.42
2:B:14:TRP:HE3	2:B:17:MET:HE2	1.84	0.42
1:E:44:GLU:HG3	1:E:46:ARG:H	1.84	0.42
1:A:90:THR:O	1:A:90(A):SER:C	2.61	0.42
1:A:166:LYS:HD3	1:A:166:LYS:HA	1.73	0.42
2:D:107:THR:O	2:D:110:TYR:HB3	2.20	0.42
1:A:90(A):SER:HB3	1:A:271:ASP:OD2	2.19	0.42
1:C:58:PRO:HB3	1:C:86:TYR:CZ	2.55	0.42
2:F:87:GLY:O	2:F:91:ILE:HG13	2.19	0.42
2:F:100:ILE:HG22	2:F:101:LEU:N	2.35	0.42
1:E:70:LEU:HD23	1:E:70:LEU:HA	1.82	0.42
1:A:103:ILE:N	1:A:103:ILE:HD12	2.35	0.42
1:A:201:TYR:CD1	1:A:201:TYR:C	2.94	0.42
1:C:66:ILE:HG12	1:C:89:GLU:OE2	2.19	0.42
1:C:76:CYS:O	1:C:77:GLU:CB	2.67	0.42
1:C:280:THR:O	1:C:303:GLY:HA3	2.19	0.42
1:C:50:LYS:HD3	1:C:275:HIS:ND1	2.35	0.42
1:C:97:CYS:HB3	1:C:98:TYR:CD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:103:ILE:N	1:C:103:ILE:HD12	2.35	0.42
2:D:37:ASP:HB3	2:D:40:SER:OG	2.20	0.42
1:C:26:VAL:HG21	1:C:317:ALA:HB2	2.01	0.42
1:A:14:CYS:O	2:B:24:TYR:HA	2.20	0.41
2:B:106:ARG:HG2	2:D:106:ARG:NH2	2.35	0.41
1:C:28:THR:C	2:D:101:LEU:HD13	2.45	0.41
1:C:167:SER:OG	1:C:242:LYS:HD2	2.20	0.41
1:A:77:GLU:O	1:A:78:SER:HB2	2.19	0.41
1:E:166:LYS:HD3	1:E:166:LYS:HA	1.81	0.41
1:A:70:LEU:HD11	1:A:112:LEU:CD1	2.38	0.41
1:C:113:SER:OG	1:C:267:ILE:HG13	2.21	0.41
2:F:168:LEU:HD23	2:F:169:ASN:N	2.36	0.41
1:A:320:LEU:HB3	2:B:111:HIS:CD2	2.56	0.41
2:B:70:PHE:HB3	2:B:74:GLU:CB	2.49	0.41
2:B:98:LEU:HD22	2:B:98:LEU:HA	1.89	0.41
1:E:114:SER:O	1:E:115:VAL:HB	2.19	0.41
1:E:201:TYR:HD2	1:E:212:LYS:HE2	1.86	0.41
1:E:253:VAL:HA	1:E:254:PRO:HD3	1.94	0.41
2:B:4:GLY:O	2:B:8:GLY:HA3	2.21	0.41
1:E:51:LEU:HD22	1:E:272:THR:HG22	2.01	0.41
1:E:320:LEU:HB3	2:F:111:HIS:CD2	2.55	0.41
1:A:204:VAL:O	1:A:210:SER:HA	2.19	0.41
2:F:2:LEU:HD22	2:F:109:ASP:CG	2.46	0.41
2:F:9:PHE:O	2:F:135:ASN:HA	2.20	0.41
2:D:133:ILE:HG13	2:D:135:ASN:HD21	1.86	0.41
1:E:17:TYR:CE1	2:F:6:ILE:HA	2.55	0.41
1:E:174:LYS:H	1:E:174:LYS:HG2	1.67	0.41
1:E:176:VAL:O	1:E:237:VAL:HG23	2.20	0.41
1:A:161:TYR:CE1	1:A:249:GLY:HA2	2.55	0.41
2:B:28:ASN:HD21	2:B:145:ASP:HA	1.86	0.41
1:C:182:ILE:HG13	1:C:233:TYR:CE1	2.56	0.41
2:B:62:GLN:HG2	2:B:63:PHE:H	1.86	0.41
2:B:129:ASN:O	2:B:141:TYR:HB2	2.21	0.41
1:C:203:PHE:CD1	1:C:203:PHE:C	2.99	0.41
2:F:35:ALA:HB3	2:F:153:LYS:NZ	2.35	0.41
1:C:73:ASN:O	1:C:76:CYS:HB2	2.21	0.41
1:E:123:LYS:HA	1:E:152:ILE:CD1	2.50	0.41
1:E:28:THR:HG23	1:E:30:LEU:H	1.86	0.40
1:E:51:LEU:HD12	1:E:51:LEU:HA	1.83	0.40
2:F:133:ILE:HG13	2:F:134:GLY:N	2.36	0.40
1:A:156:LYS:HG3	1:A:194:LEU:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:ASP:HB3	1:A:296:ASN:HD21	1.85	0.40
1:A:77:GLU:O	1:A:78:SER:CB	2.69	0.40
2:F:161:LYS:HD2	2:F:162:TYR:CZ	2.57	0.40
1:C:76:CYS:O	1:C:77:GLU:HG2	2.22	0.40
2:D:134:GLY:O	2:D:135:ASN:CG	2.65	0.40
1:E:115:VAL:HG21	1:E:260: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	316/329 (96%)	293 (93%)	20 (6%)	3 (1%)	14	47
1	C	293/329 (89%)	273 (93%)	18 (6%)	2 (1%)	18	52
1	E	316/329 (96%)	283 (90%)	29 (9%)	4 (1%)	9	40
2	B	169/179 (94%)	164 (97%)	5 (3%)	0	100	100
2	D	172/179 (96%)	156 (91%)	16 (9%)	0	100	100
2	F	168/179 (94%)	156 (93%)	12 (7%)	0	100	100
All	All	1434/1524 (94%)	1325 (92%)	100 (7%)	9 (1%)	21	56

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116(B)	PHE
1	A	266	SER
1	E	116(B)	PHE
1	E	266	SER
1	A	115	VAL
1	C	266	SER

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Mol	Chain	Res	Type
1	E	115	VAL
1	C	116(B)	PHE
1	E	297	ILE

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	280/289 (97%)	231 (82%)	49 (18%)	2	10
1	C	267/289 (92%)	221 (83%)	46 (17%)	2	10
1	E	281/289 (97%)	241 (86%)	40 (14%)	3	17
2	B	146/152 (96%)	131 (90%)	15 (10%)	7	28
2	D	149/152 (98%)	128 (86%)	21 (14%)	3	17
2	F	145/152 (95%)	128 (88%)	17 (12%)	5	23
All	All	1268/1323 (96%)	1080 (85%)	188 (15%)	3	15

All (188) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	THR
1	A	29	VAL
1	A	46	ARG
1	A	48	ASN
1	A	66	ILE
1	A	76	CYS
1	A	77	GLU
1	A	90	THR
1	A	90(A)	SER
1	A	111	GLN
1	A	116	SER
1	A	116(A)	SER
1	A	116(B)	PHE
1	A	116(C)	GLU
1	A	117	ARG

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Mol	Chain	Res	Type
1	A	129	ASN
1	A	132	THR
1	A	133	ASN
1	A	133(A)	ARG
1	A	145	LYS
1	A	156	LYS
1	A	157	LYS
1	A	163	LYS
1	A	164	LEU
1	A	174	LYS
1	A	186	SER
1	A	188	SER
1	A	190	ASP
1	A	193	SER
1	A	196	GLN
1	A	199	ASP
1	A	201	TYR
1	A	203	PHE
1	A	205	CYS
1	A	207	SER
1	A	214	LYS
1	A	216	GLU
1	A	220	CYS
1	A	226	GLN
1	A	237	VAL
1	A	260	MET
1	A	269	ILE
1	A	272	THR
1	A	288	ILE
1	A	295	GLN
1	A	297	ILE
1	A	312	THR
1	A	321	ARG
1	A	323	VAL
2	B	15	THR
2	B	18	ILE
2	B	27	GLN
2	B	29	GLU
2	B	30	GLN
2	B	45	ILE
2	B	58	LYS
2	B	59	MET

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Mol	Chain	Res	Type
2	B	61	THR
2	B	68	LYS
2	B	98	LEU
2	B	102	LEU
2	B	106	ARG
2	B	131	ARG
2	B	139	GLU
1	C	12	THR
1	C	15	ILE
1	C	22	SER
1	C	23	THR
1	C	28	THR
1	C	30	LEU
1	C	45	ASP
1	C	51	LEU
1	C	54	LEU
1	C	56	VAL
1	C	63	LYS
1	C	64	CYS
1	C	90	THR
1	C	90(A)	SER
1	C	115	VAL
1	C	131	ASP
1	C	158	GLU
1	C	167	SER
1	C	169	ILE
1	C	171	ASN
1	C	172	LYS
1	C	187	THR
1	C	188	SER
1	C	196	GLN
1	C	202	VAL
1	C	204	VAL
1	C	205	CYS
1	C	210	SER
1	C	238	GLU
1	C	244	THR
1	C	248	THR
1	C	253	VAL
1	C	273	SER
1	C	274	VAL
1	C	278	ASN

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Mol	Chain	Res	Type
1	C	279	THR
1	C	280	THR
1	C	288	ILE
1	C	295	GLN
1	C	301	THR
1	C	307	LYS
1	C	312	THR
1	C	315	ARG
1	C	320	LEU
1	C	322	ASN
1	C	323	VAL
2	D	11	GLU
2	D	18	ILE
2	D	40	SER
2	D	50	ASN
2	D	58	LYS
2	D	64	ILE
2	D	66	VAL
2	D	68	LYS
2	D	69	GLU
2	D	74	GLU
2	D	82	LYS
2	D	100	ILE
2	D	101	LEU
2	D	103	GLU
2	D	106	ARG
2	D	132	GLU
2	D	143	LYS
2	D	151	SER
2	D	163	SER
2	D	165	GLU
2	D	172	GLU
1	E	23	THR
1	E	25	THR
1	E	28	THR
1	E	29	VAL
1	E	35	THR
1	E	40	VAL
1	E	43	LEU
1	E	45	ASP
1	E	50	LYS
1	E	55	ARG

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Mol	Chain	Res	Type
1	E	61	LEU
1	E	63	LYS
1	E	66	ILE
1	E	93	ASP
1	E	132	THR
1	E	135	VAL
1	E	136	THR
1	E	164	LEU
1	E	169	ILE
1	E	174	LYS
1	E	186	SER
1	E	188	SER
1	E	193	SER
1	E	199	ASP
1	E	204	VAL
1	E	213	PHE
1	E	220	CYS
1	E	222	LYS
1	E	223	VAL
1	E	237	VAL
1	E	253	VAL
1	E	272	THR
1	E	277	CYS
1	E	280	THR
1	E	283	THR
1	E	295	GLN
1	E	310	LYS
1	E	313	LYS
1	E	315	ARG
1	E	323	VAL
2	F	22	TYR
2	F	54	SER
2	F	58	LYS
2	F	66	VAL
2	F	68	LYS
2	F	69	GLU
2	F	74	GLU
2	F	98	LEU
2	F	100	ILE
2	F	103	GLU
2	F	105	GLU
2	F	113	SER

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Mol	Chain	Res	Type
2	F	131	ARG
2	F	133	ILE
2	F	143	LYS
2	F	147	LYS
2	F	156	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	GLN
1	A	129	ASN
1	A	196	GLN
1	A	197	ASN
1	A	296	ASN
2	B	30	GLN
2	B	169	ASN
1	C	150	ASN
1	C	322	ASN
2	D	27	GLN
2	D	125	GLN
1	E	94	ASN
1	E	192	GLN
1	E	282	GLN
2	F	60	ASN
2	F	72	HIS
2	F	114	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 oligosaccharides in this entry.

5.6 Ligand geometry

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	C	401	1	14,14,15	0.53	0	17,19,21	0.79	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	401	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	401	NAG	O5-C5-C6-O6
3	C	401	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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(Å²)	Q<0.9
1	A	320/329 (97%)	0.42	9 (2%)	55	35	22, 48, 89, 124	0
1	C	301/329 (91%)	0.23	8 (2%)	56	36	21, 34, 70, 109	0
1	E	320/329 (97%)	0.74	24 (7%)	20	13	31, 66, 111, 156	0
2	B	171/179 (95%)	0.28	8 (4%)	36	23	22, 41, 78, 145	0
2	D	174/179 (97%)	0.29	9 (5%)	33	21	16, 38, 90, 153	0
2	F	170/179 (94%)	0.44	11 (6%)	25	16	19, 51, 109, 157	0
All	All	1456/1524 (95%)	0.42	69 (4%)	36	23	16, 46, 96, 157	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	65	ALA	6.3
2	F	59	MET	5.0
2	D	65	ALA	4.7
1	E	82	ALA	4.7
2	F	66	VAL	4.4
2	D	64	ILE	4.3
2	F	63	PHE	4.3
1	E	96	THR	4.1
2	F	67	GLY	4.0
2	B	63	PHE	4.0
2	B	64	ILE	3.8
1	E	78	SER	3.7
2	D	66	VAL	3.7
1	E	97	CYS	3.6
1	E	142	ALA	3.5
2	F	61	THR	3.5
2	D	174	ASP	3.4
2	B	67	GLY	3.4
2	B	66	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
2	F	65	ALA	3.2
2	D	61	THR	3.2
1	E	227	ALA	3.2
1	E	141	TYR	3.1
1	E	264	SER	3.1
2	D	63	PHE	3.0
1	A	95	GLY	3.0
1	E	77	GLU	2.8
2	F	62	GLN	2.7
1	A	203	PHE	2.7
1	A	149	ARG	2.6
2	F	60	ASN	2.6
1	C	95	GLY	2.6
1	E	95	GLY	2.5
1	A	228	GLY	2.5
1	C	264	SER	2.5
2	F	134	GLY	2.5
1	A	264	SER	2.5
1	E	139	CYS	2.5
2	F	64	ILE	2.5
2	B	61	THR	2.5
1	E	136	THR	2.4
1	E	143	GLY	2.4
1	C	220	CYS	2.4
1	E	64	CYS	2.4
1	A	82	ALA	2.4
2	D	173	ILE	2.4
1	C	9	PRO	2.4
1	E	155	VAL	2.3
1	C	10	GLY	2.3
1	C	82	ALA	2.3
2	D	60	ASN	2.3
1	C	77	GLU	2.3
1	E	265	GLY	2.3
1	E	100	GLY	2.3
1	E	304	GLU	2.2
2	B	62	GLN	2.2
1	A	116(A)	SER	2.2
1	E	137	ALA	2.2
2	D	62	GLN	2.1
1	C	116	SER	2.1
2	B	1	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	115	VAL	2.1
1	A	141	TYR	2.1
1	E	192	GLN	2.1
2	F	58	LYS	2.0
1	E	90(A)	SER	2.0
1	E	11	ASP	2.0
1	E	144	ALA	2.0
1	A	196	GLN	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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	C	401	14/15	0.89	0.10	23,37,41,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.