



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 22, 2026 – 02:23 PM JST

PDB ID : 9WN0 / pdb_00009wn0
Title : Crystal structure of phospholipase A2
Authors : Zhu, D.
Deposited on : 2025-09-04
Resolution : 2.00 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
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

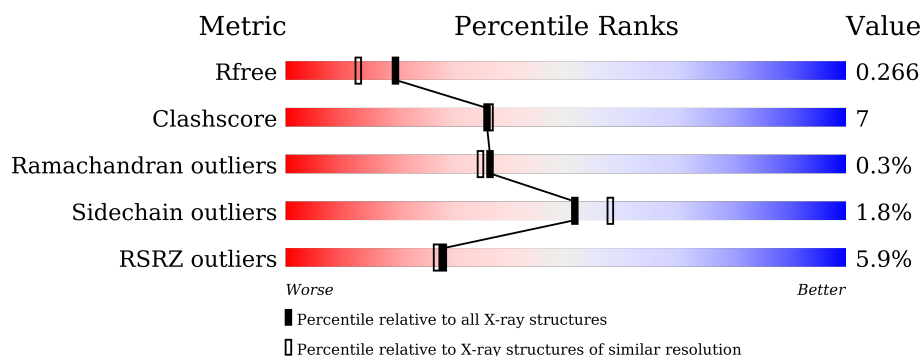
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	333	<div> <div>4%</div> <div> <div></div> <div>85%</div> <div>15%</div> </div> </div>
1	B	333	<div> <div>4%</div> <div> <div></div> <div>86%</div> <div>14%</div> </div> </div>
1	C	333	<div> <div>5%</div> <div> <div></div> <div>87%</div> <div>12%</div> </div> </div>
1	D	333	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>16%</div> </div> </div>
1	E	333	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>17%</div> </div> </div>
1	F	333	<div> <div>8%</div> <div> <div></div> <div>77%</div> <div>22%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	333	<div><div>8%</div><div>83%</div><div>17%</div><div></div></div>
1	H	333	<div><div>11%</div><div>76%</div><div>22%</div><div></div></div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 21314 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 Patatin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2605	1673	435	490	7			
1	B	333	Total	C	N	O	S	0	0	0
			2605	1673	435	490	7			
1	C	333	Total	C	N	O	S	0	0	0
			2605	1673	435	490	7			
1	D	333	Total	C	N	O	S	0	0	0
			2605	1673	435	490	7			
1	E	333	Total	C	N	O	S	0	0	0
			2605	1673	435	490	7			
1	F	333	Total	C	N	O	S	0	0	0
			2605	1673	435	490	7			
1	G	333	Total	C	N	O	S	0	0	0
			2605	1673	435	490	7			
1	H	333	Total	C	N	O	S	0	0	0
			2605	1673	435	490	7			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	75	Total	O	0	0
			75	75		
2	B	80	Total	O	0	0
			80	80		
2	C	58	Total	O	0	0
			58	58		
2	D	68	Total	O	0	0
			68	68		
2	E	59	Total	O	0	0
			59	59		
2	F	44	Total	O	0	0
			44	44		

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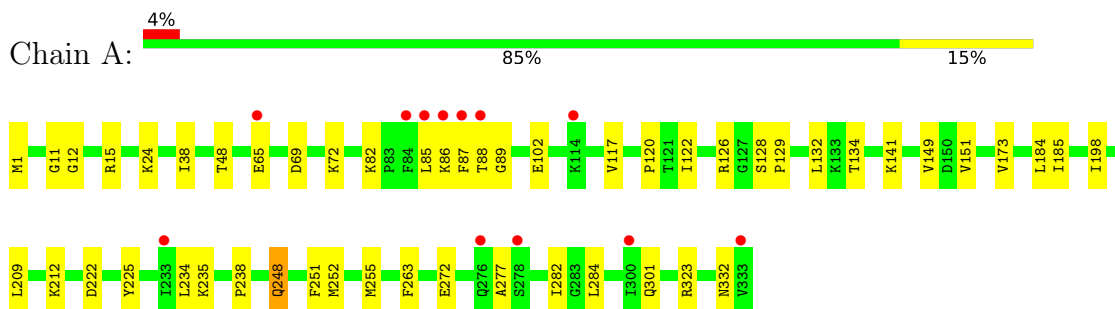
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	59	Total	O	0	0
			59	59		
2	H	31	Total	O	0	0
			31	31		

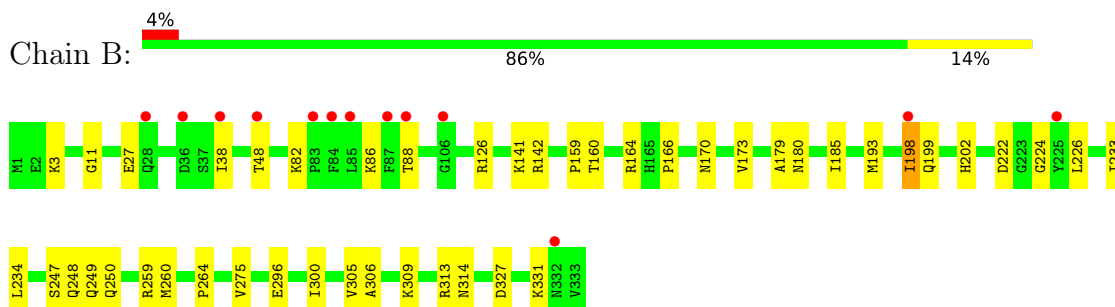
3 Residue-property plots [i](#)

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.

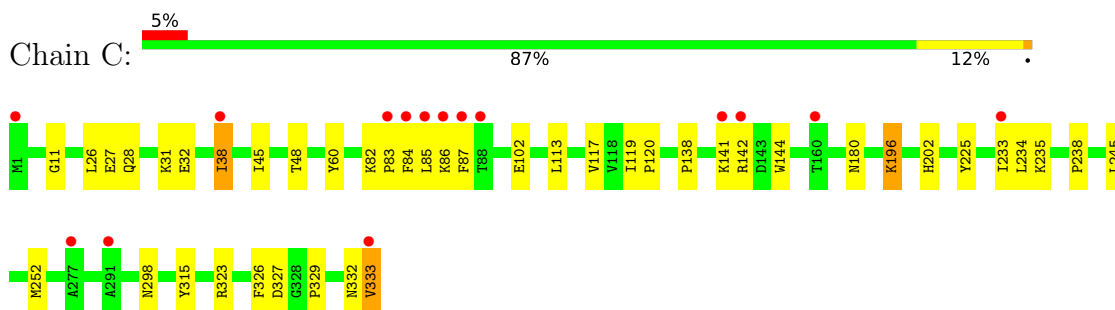
- Molecule 1: Patatin



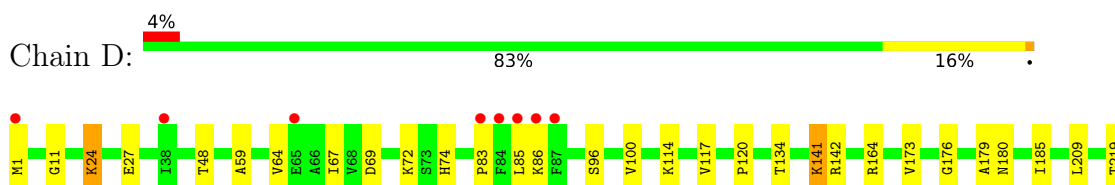
- Molecule 1: Patatin

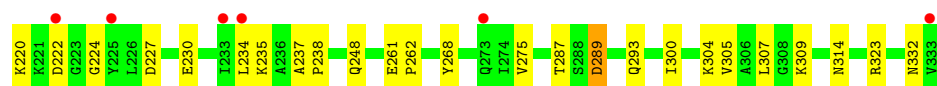


- Molecule 1: Patatin

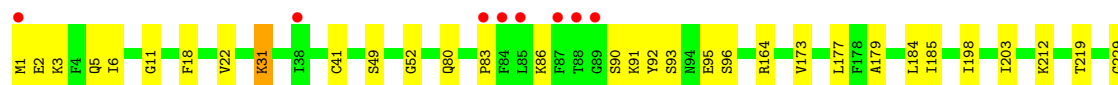
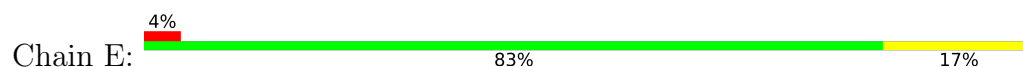


- Molecule 1: Patatin

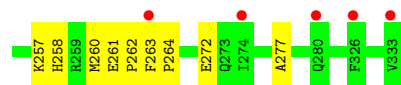
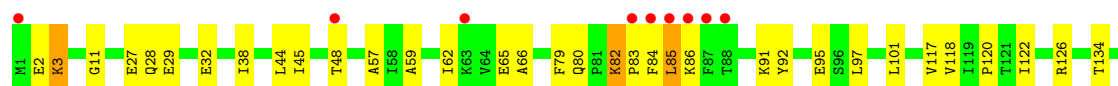
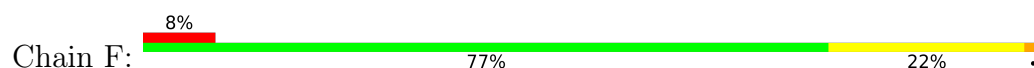




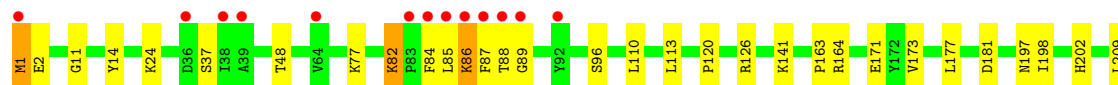
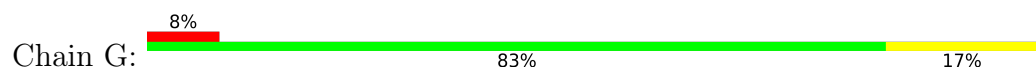
● Molecule 1: Patatin



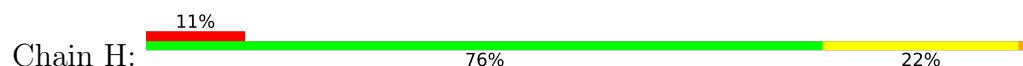
● Molecule 1: Patatin

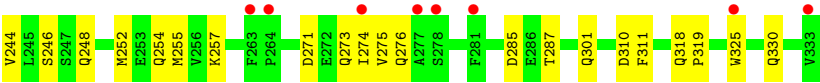


● Molecule 1: Patatin



● Molecule 1: Patatin





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	86.01Å 86.09Å 88.57Å 86.67° 73.37° 86.44°	Depositor
Resolution (Å)	44.64 – 2.00 44.64 – 2.00	Depositor EDS
% Data completeness (in resolution range)	90.6 (44.64-2.00) 90.6 (44.64-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.00Å)	Xtriage
Refinement program	PHENIX (dev_3885: ???)	Depositor
R, R_{free}	0.216 , 0.268 0.220 , 0.266	Depositor DCC
R_{free} test set	7746 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	33.7	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21314	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.42% 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

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	0.51	0/2665	0.73	1/3604 (0.0%)
1	B	0.49	0/2665	0.70	0/3604
1	C	0.49	0/2665	0.74	1/3604 (0.0%)
1	D	0.47	0/2665	0.71	0/3604
1	E	0.53	0/2665	0.78	0/3604
1	F	0.52	0/2665	0.77	2/3604 (0.1%)
1	G	0.46	0/2665	0.73	2/3604 (0.1%)
1	H	0.44	0/2665	0.70	1/3604 (0.0%)
All	All	0.49	0/21320	0.73	7/28832 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	160	THR	CB-CA-C	-8.68	101.24	109.83
1	G	87	PHE	N-CA-C	-5.88	105.36	112.54
1	G	311	PHE	CA-CB-CG	5.43	119.23	113.80
1	H	139	ASN	N-CA-C	-5.21	99.71	110.80
1	A	88	THR	N-CA-C	-5.09	106.76	112.87
1	F	82	LYS	CB-CA-C	5.06	116.83	109.08
1	C	329	PRO	N-CA-C	5.00	120.11	113.86

There are no chirality outliers.

There are no planarity outliers.

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	2605	0	2606	41	1
1	B	2605	0	2606	33	0
1	C	2605	0	2606	30	0
1	D	2605	0	2606	39	1
1	E	2605	0	2606	34	0
1	F	2605	0	2606	55	1
1	G	2605	0	2606	48	1
1	H	2605	0	2606	56	0
2	A	75	0	0	3	0
2	B	80	0	0	2	0
2	C	58	0	0	1	0
2	D	68	0	0	2	0
2	E	59	0	0	1	0
2	F	44	0	0	2	0
2	G	59	0	0	3	0
2	H	31	0	0	0	0
All	All	21314	0	20848	304	2

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 7.

All (304) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:GLN:HE21	1:B:248:GLN:HE21	1.14	0.94
1:G:88:THR:HG21	1:H:226:LEU:HG	1.55	0.89
1:D:268:TYR:OH	1:D:309:LYS:NZ	2.07	0.87
1:A:85:LEU:HD12	1:A:86:LYS:HG2	1.60	0.84
1:E:179:ALA:HB1	1:E:185:ILE:HD11	1.63	0.78
1:D:222:ASP:OD1	1:D:227:ASP:HB2	1.83	0.78
1:H:179:ALA:HB1	1:H:185:ILE:HD11	1.66	0.76
1:A:248:GLN:HE21	1:B:248:GLN:NE2	1.82	0.75
1:A:184:LEU:HD23	1:A:252:MET:HE2	1.68	0.75
1:C:28:GLN:NE2	1:C:32:GLU:OE2	2.18	0.75
1:F:141:LYS:O	1:F:141:LYS:HG2	1.89	0.73
1:G:88:THR:HG22	1:H:225:TYR:HB2	1.71	0.73
1:C:48:THR:HG22	1:C:120:PRO:HG2	1.71	0.72
1:A:248:GLN:NE2	1:B:248:GLN:HE21	1.87	0.72
1:D:219:THR:CG2	1:D:230:GLU:HA	2.20	0.71
1:D:179:ALA:HB1	1:D:185:ILE:HD11	1.72	0.71
1:D:69:ASP:HA	1:D:72:LYS:HE2	1.74	0.70
1:D:219:THR:HG21	1:D:230:GLU:HA	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:27:GLU:HB2	1:F:38:ILE:HD11	1.72	0.69
1:H:137:ASN:ND2	1:H:193:MET:O	2.25	0.69
1:F:254:GLN:HA	1:F:257:LYS:HE3	1.75	0.69
1:B:198:ILE:HD11	1:B:259:ARG:CZ	2.23	0.68
1:B:27:GLU:HB2	1:B:38:ILE:HD11	1.77	0.66
1:G:163:PRO:O	2:G:402:HOH:O	2.15	0.64
1:A:69:ASP:HA	1:A:72:LYS:HG2	1.79	0.63
1:H:181:ASP:HB3	1:H:252:MET:HE2	1.80	0.63
1:D:180:ASN:H	1:D:248:GLN:HE21	1.45	0.62
1:B:179:ALA:HB1	1:B:185:ILE:HD11	1.81	0.62
1:A:48:THR:HG22	1:A:120:PRO:HG2	1.81	0.62
1:A:89:GLY:HA2	1:B:222:ASP:OD2	2.00	0.62
1:G:14:TYR:OH	2:G:401:HOH:O	2.15	0.62
1:C:84:PHE:O	1:C:85:LEU:C	2.43	0.61
1:F:179:ALA:HA	1:F:252:MET:HE1	1.81	0.61
1:E:1:MET:SD	1:E:1:MET:N	2.73	0.61
1:E:5:GLN:NE2	1:E:41:CYS:O	2.34	0.61
1:F:48:THR:OG1	1:F:120:PRO:HG2	2.01	0.61
1:B:296:GLU:O	1:B:300:ILE:HG13	2.00	0.61
1:E:184:LEU:HD23	1:E:252:MET:HE2	1.82	0.61
1:F:82:LYS:HB3	1:F:83:PRO:HD2	1.83	0.61
1:D:180:ASN:N	1:D:248:GLN:HE21	1.98	0.60
1:G:82:LYS:HD2	1:G:86:LYS:HZ2	1.68	0.59
1:C:138:PRO:HG3	1:C:326:PHE:CD2	2.37	0.59
1:E:198:ILE:HD11	1:E:259:ARG:HH21	1.68	0.59
1:G:164:ARG:HG2	1:G:173:VAL:HG22	1.85	0.59
1:C:86:LYS:O	1:C:87:PHE:C	2.45	0.59
1:F:197:ASN:ND2	2:F:401:HOH:O	2.29	0.59
1:C:82:LYS:HB2	1:C:83:PRO:HD2	1.84	0.58
1:E:219:THR:HG21	1:E:230:GLU:HG2	1.86	0.58
1:F:137:ASN:HB3	1:F:140:LEU:HG	1.86	0.57
1:G:86:LYS:HE2	1:G:89:GLY:HA3	1.86	0.57
1:C:84:PHE:O	1:C:87:PHE:N	2.37	0.57
1:E:80:GLN:OE1	1:E:96:SER:OG	2.16	0.56
1:D:234:LEU:CD1	1:F:141:LYS:NZ	2.68	0.56
1:G:216:ASN:O	1:G:219:THR:HG22	2.05	0.56
1:H:220:LYS:HE3	1:H:226:LEU:HB3	1.88	0.56
1:F:80:GLN:OE1	1:F:95:GLU:HB3	2.05	0.56
1:F:141:LYS:O	1:F:141:LYS:CG	2.54	0.55
1:D:85:LEU:HG	1:D:86:LYS:HG3	1.87	0.55
1:G:275:VAL:O	1:G:276:GLN:C	2.49	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LEU:CD2	1:B:233:ILE:HD11	2.37	0.55
1:G:86:LYS:CE	1:G:89:GLY:HA3	2.37	0.55
1:H:236:ALA:O	1:H:240:ILE:HG13	2.06	0.55
1:H:318:GLN:HB3	1:H:319:PRO:HD2	1.89	0.55
1:C:27:GLU:HA	1:C:38:ILE:HG12	1.88	0.55
1:D:176:GLY:O	1:D:248:GLN:NE2	2.39	0.55
1:F:3:LYS:HG2	1:F:202:HIS:CD2	2.42	0.55
1:G:209:LEU:HD13	1:G:271:ASP:HB2	1.88	0.55
1:G:248:GLN:OE1	1:H:248:GLN:NE2	2.37	0.55
1:H:144:TRP:CZ2	1:H:145:LYS:HE2	2.42	0.54
1:E:18:PHE:O	1:E:22:VAL:HG23	2.08	0.54
1:C:27:GLU:O	1:C:31:LYS:HG3	2.08	0.54
1:C:298:ASN:ND2	2:C:403:HOH:O	2.27	0.54
1:D:74:HIS:HD2	1:D:287:THR:O	1.91	0.53
1:B:275:VAL:HG12	1:B:305:VAL:HG21	1.90	0.53
1:D:234:LEU:HD13	1:F:141:LYS:NZ	2.24	0.53
1:F:198:ILE:HG23	1:F:263:PHE:CE2	2.43	0.53
1:G:248:GLN:O	1:G:252:MET:HG2	2.08	0.53
1:B:193:MET:HE2	1:B:193:MET:HA	1.89	0.53
1:F:229:GLY:HA3	1:F:232:SER:O	2.08	0.53
1:G:126:ARG:O	1:H:212:LYS:NZ	2.41	0.53
1:G:89:GLY:N	1:H:224:GLY:HA3	2.23	0.53
1:G:181:ASP:HB3	1:G:252:MET:HB2	1.90	0.53
1:B:82:LYS:NZ	1:B:88:THR:O	2.41	0.53
1:G:197:ASN:OD1	1:G:198:ILE:N	2.42	0.53
1:D:64:VAL:HA	1:D:67:ILE:HD12	1.90	0.52
1:C:26:LEU:C	1:C:38:ILE:HD11	2.33	0.52
1:E:234:LEU:CD2	1:F:233:ILE:HD11	2.39	0.52
1:C:141:LYS:HB2	1:C:327:ASP:OD2	2.09	0.52
1:B:226:LEU:O	2:B:401:HOH:O	2.19	0.52
1:C:138:PRO:HG3	1:C:326:PHE:CE2	2.44	0.52
1:E:310:ASP:OD1	1:E:311:PHE:N	2.42	0.52
1:F:272:GLU:HB3	1:F:277:ALA:HB3	1.91	0.52
1:A:141:LYS:NZ	2:A:408:HOH:O	2.43	0.52
1:A:222:ASP:OD2	2:A:401:HOH:O	2.19	0.52
1:D:117:VAL:HG22	1:D:134:THR:HG23	1.92	0.52
1:A:120:PRO:HG3	1:A:185:ILE:HD13	1.91	0.51
1:D:219:THR:HG23	1:D:230:GLU:HA	1.92	0.51
1:H:93:SER:OG	1:H:95:GLU:HG2	2.10	0.51
1:B:27:GLU:HA	1:B:38:ILE:HG12	1.91	0.51
1:A:87:PHE:C	1:A:89:GLY:H	2.17	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LEU:HD23	1:B:233:ILE:HD11	1.92	0.51
1:D:59:ALA:O	1:D:114:LYS:HG2	2.11	0.51
1:F:235:LYS:HA	1:F:238:PRO:HD2	1.92	0.51
1:A:15:ARG:HD3	1:A:282:ILE:O	2.11	0.51
1:F:253:GLU:HG2	1:F:257:LYS:HE2	1.93	0.51
1:H:72:LYS:HB3	1:H:77:LYS:NZ	2.26	0.51
1:E:177:LEU:HG	1:F:244:VAL:HG22	1.94	0.50
1:A:87:PHE:C	1:A:89:GLY:N	2.67	0.50
1:G:84:PHE:CG	1:G:85:LEU:N	2.80	0.50
1:A:235:LYS:C	1:A:238:PRO:HD2	2.36	0.50
1:A:248:GLN:O	1:A:252:MET:HG3	2.11	0.50
1:E:271:ASP:O	1:E:274:ILE:HG12	2.12	0.50
1:G:202:HIS:ND1	1:G:266:GLN:HG2	2.27	0.50
1:A:122:ILE:HB	1:A:173:VAL:HB	1.93	0.50
1:C:85:LEU:O	1:C:86:LYS:HB2	2.12	0.50
1:C:235:LYS:C	1:C:238:PRO:HD2	2.37	0.50
1:C:323:ARG:O	1:C:332:ASN:ND2	2.35	0.49
1:E:248:GLN:NE2	1:F:248:GLN:OE1	2.44	0.49
1:G:237:ALA:HB3	1:G:238:PRO:HD3	1.93	0.49
1:A:323:ARG:O	1:A:332:ASN:ND2	2.44	0.49
1:C:180:ASN:ND2	1:C:245:LEU:O	2.36	0.49
1:G:219:THR:HG23	1:G:220:LYS:O	2.12	0.49
1:G:268:TYR:CD2	1:G:312:ILE:HD11	2.47	0.49
1:E:83:PRO:HB2	1:E:86:LYS:HB2	1.93	0.49
1:E:318:GLN:HB2	1:E:321:ARG:HH12	1.77	0.49
1:H:254:GLN:HA	1:H:257:LYS:HE3	1.93	0.49
1:E:318:GLN:OE1	1:E:321:ARG:NH2	2.44	0.49
1:F:261:GLU:N	1:F:262:PRO:CD	2.75	0.49
1:G:232:SER:HB2	1:G:235:LYS:HG2	1.95	0.49
1:C:180:ASN:H	1:C:252:MET:HE1	1.77	0.49
1:F:254:GLN:HG2	1:F:258:HIS:CE1	2.47	0.49
1:D:164:ARG:HG2	1:D:173:VAL:HG22	1.95	0.49
1:H:150:ASP:HB3	1:H:167:ILE:HG12	1.93	0.49
1:E:237:ALA:HB3	1:E:238:PRO:HD3	1.95	0.49
1:H:83:PRO:HD2	1:H:84:PHE:CE1	2.48	0.49
1:H:275:VAL:HG11	1:H:301:GLN:HB3	1.93	0.49
1:H:4:PHE:HB2	1:H:196:LYS:HE2	1.95	0.48
1:F:48:THR:HG21	1:F:176:GLY:HA2	1.95	0.48
1:F:260:MET:C	1:F:262:PRO:HD2	2.38	0.48
1:H:310:ASP:OD1	1:H:311:PHE:N	2.46	0.48
1:D:120:PRO:HG3	1:D:185:ILE:HD13	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:300:ILE:O	1:D:304:LYS:HG2	2.13	0.48
1:E:164:ARG:HG2	1:E:173:VAL:HG22	1.95	0.48
1:F:44:LEU:HD11	1:F:118:VAL:HG23	1.94	0.48
1:G:110:LEU:O	1:G:113:LEU:HG	2.14	0.48
1:H:65:GLU:HG3	1:H:66:ALA:N	2.29	0.48
1:A:1:MET:SD	1:A:1:MET:N	2.77	0.48
1:F:249:GLN:HG3	1:F:250:GLN:N	2.28	0.48
1:H:82:LYS:HD3	1:H:87:PHE:HB3	1.94	0.48
1:A:89:GLY:HA2	1:B:224:GLY:HA3	1.96	0.47
1:H:85:LEU:HG	1:H:85:LEU:O	2.15	0.47
1:E:49:SER:O	1:E:52:GLY:N	2.47	0.47
1:D:180:ASN:H	1:D:248:GLN:NE2	2.11	0.47
1:D:289:ASP:O	1:D:293:GLN:HG3	2.15	0.47
1:F:178:PHE:C	1:F:248:GLN:NE2	2.73	0.47
1:F:179:ALA:HB1	1:F:185:ILE:HD11	1.96	0.47
1:F:235:LYS:C	1:F:238:PRO:HD2	2.40	0.47
1:E:243:LEU:HD11	1:F:122:ILE:HG21	1.96	0.47
1:G:271:ASP:O	1:G:274:ILE:HG12	2.14	0.47
1:G:272:GLU:HA	1:G:275:VAL:HG22	1.97	0.47
1:H:3:LYS:HG2	1:H:202:HIS:CE1	2.49	0.47
1:H:45:ILE:HD12	1:H:59:ALA:HB2	1.97	0.47
1:A:272:GLU:HB3	1:A:277:ALA:HB3	1.96	0.47
1:G:77:LYS:HA	1:G:96:SER:OG	2.14	0.47
1:H:209:LEU:HD11	1:H:273:GLN:HG2	1.96	0.47
1:C:326:PHE:CE2	1:C:333:VAL:HG11	2.50	0.47
1:H:79:PHE:HB3	1:H:92:TYR:CD1	2.50	0.47
1:E:2:GLU:HG3	1:E:3:LYS:N	2.30	0.46
1:F:181:ASP:OD2	1:F:183:SER:HB2	2.15	0.46
1:F:159:PRO:O	1:F:160:THR:OG1	2.28	0.46
1:G:86:LYS:HB3	1:G:88:THR:O	2.15	0.46
1:A:198:ILE:HG23	1:A:263:PHE:CE2	2.50	0.46
1:H:84:PHE:HB2	1:H:87:PHE:HB2	1.97	0.46
1:A:251:PHE:CE2	1:A:255:MET:HE3	2.51	0.46
1:F:28:GLN:O	1:F:32:GLU:HG3	2.16	0.46
1:D:24:LYS:HA	1:D:24:LYS:HD2	1.68	0.46
1:E:2:GLU:HG3	1:E:3:LYS:H	1.81	0.46
1:G:327:ASP:HA	1:G:331:LYS:HD2	1.97	0.46
1:E:275:VAL:O	1:E:276:GLN:C	2.58	0.46
1:G:86:LYS:HE2	1:G:86:LYS:HB2	1.60	0.46
1:G:171:GLU:OE2	1:H:221:LYS:NZ	2.49	0.46
1:G:202:HIS:CE1	1:G:266:GLN:HG2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:234:LEU:HD12	1:C:234:LEU:N	2.31	0.45
1:D:83:PRO:HG2	1:D:86:LYS:HB2	1.99	0.45
1:A:225:TYR:OH	1:B:234:LEU:O	2.29	0.45
1:F:237:ALA:HB3	1:F:238:PRO:HD3	1.98	0.45
1:A:129:PRO:HG3	1:B:247:SER:HA	1.98	0.45
1:A:132:LEU:HD12	1:A:151:VAL:HG13	1.98	0.45
1:E:293:GLN:OE1	2:E:401:HOH:O	2.21	0.45
1:D:275:VAL:HG12	1:D:305:VAL:HG21	1.99	0.45
1:F:65:GLU:HG3	1:F:66:ALA:N	2.32	0.45
1:A:117:VAL:HG22	1:A:134:THR:HG23	1.98	0.45
1:H:285:ASP:O	1:H:287:THR:HG23	2.16	0.45
1:A:212:LYS:NZ	1:H:276:GLN:O	2.42	0.45
1:B:141:LYS:O	1:B:142:ARG:HD3	2.17	0.45
1:B:264:PRO:HG3	1:F:263:PHE:CZ	2.52	0.45
1:H:84:PHE:N	1:H:84:PHE:CD1	2.85	0.45
1:H:235:LYS:C	1:H:238:PRO:HD2	2.41	0.45
1:C:333:VAL:HG13	1:H:319:PRO:HD3	1.99	0.45
1:D:237:ALA:HB3	1:D:238:PRO:HD3	1.98	0.45
1:G:177:LEU:HG	1:H:244:VAL:HG22	1.99	0.45
1:H:209:LEU:C	1:H:211:SER:H	2.23	0.45
1:H:276:GLN:O	1:H:276:GLN:HG2	2.17	0.45
1:A:12:GLY:C	1:A:284:LEU:HD23	2.42	0.44
1:G:235:LYS:C	1:G:238:PRO:HD2	2.42	0.44
1:G:234:LEU:O	1:H:225:TYR:OH	2.32	0.44
1:D:74:HIS:CD2	1:D:287:THR:O	2.71	0.44
1:G:220:LYS:C	1:G:221:LYS:HG3	2.43	0.44
1:B:159:PRO:O	1:B:160:THR:OG1	2.32	0.44
1:D:323:ARG:O	1:D:332:ASN:ND2	2.45	0.44
1:D:209:LEU:HD23	1:D:209:LEU:HA	1.82	0.44
1:E:229:GLY:HA3	1:E:232:SER:O	2.17	0.44
1:B:260:MET:HE3	1:B:260:MET:HB3	1.86	0.44
1:B:249:GLN:HG3	1:B:250:GLN:N	2.33	0.43
1:F:227:ASP:OD1	1:F:227:ASP:C	2.60	0.43
1:G:37:SER:HB3	2:G:450:HOH:O	2.17	0.43
1:A:102:GLU:HG2	1:A:149:VAL:HG21	2.00	0.43
1:C:326:PHE:CD2	1:C:333:VAL:HG11	2.53	0.43
1:F:97:LEU:O	1:F:101:LEU:HG	2.17	0.43
1:D:224:GLY:O	1:D:227:ASP:HB3	2.17	0.43
1:D:141:LYS:O	1:D:142:ARG:HD3	2.18	0.43
1:G:84:PHE:CE2	1:G:86:LYS:HG3	2.54	0.43
1:E:93:SER:OG	1:E:95:GLU:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:LYS:HE2	1:A:65:GLU:HG2	2.00	0.43
1:A:212:LYS:HE2	1:B:126:ARG:O	2.19	0.43
1:B:314:ASN:ND2	2:B:407:HOH:O	2.36	0.43
1:G:215:LEU:HB3	1:H:125:THR:HA	2.01	0.43
1:G:229:GLY:HA3	1:G:232:SER:O	2.19	0.43
1:A:301:GLN:HG3	1:G:126:ARG:NE	2.33	0.43
1:C:196:LYS:HD2	1:C:196:LYS:HA	1.69	0.43
1:H:85:LEU:HD23	1:H:85:LEU:H	1.83	0.43
1:B:48:THR:HG21	1:B:180:ASN:HA	2.01	0.43
1:D:234:LEU:HD12	1:D:234:LEU:H	1.83	0.43
1:D:261:GLU:N	1:D:262:PRO:CD	2.81	0.43
1:F:45:ILE:HD12	1:F:59:ALA:HB2	2.01	0.43
1:H:212:LYS:HA	1:H:246:SER:OG	2.17	0.42
1:D:27:GLU:OE1	2:D:401:HOH:O	2.21	0.42
1:B:166:PRO:HA	1:B:170:ASN:O	2.19	0.42
1:B:249:GLN:HE21	1:B:250:GLN:HG3	1.85	0.42
1:E:290:ALA:O	1:E:294:VAL:HG23	2.18	0.42
1:G:1:MET:HE2	1:G:1:MET:HB2	1.68	0.42
1:H:84:PHE:CD1	1:H:87:PHE:CG	3.08	0.42
1:H:85:LEU:O	1:H:86:LYS:C	2.62	0.42
1:A:126:ARG:NH2	1:A:128:SER:OG	2.51	0.42
1:C:45:ILE:O	1:C:117:VAL:HA	2.20	0.42
1:C:119:ILE:HA	1:C:120:PRO:HD3	1.83	0.42
1:H:142:ARG:HD2	1:H:145:LYS:HZ1	1.85	0.42
1:E:212:LYS:HE2	1:F:126:ARG:O	2.20	0.42
1:A:126:ARG:NH2	2:A:420:HOH:O	2.53	0.42
1:E:31:LYS:HB3	1:E:31:LYS:HE3	1.58	0.42
1:F:85:LEU:HD21	1:F:91:LYS:HE3	2.02	0.42
1:H:7:LEU:HB2	1:H:42:PHE:CE2	2.55	0.42
1:F:262:PRO:HB2	1:F:263:PHE:CD2	2.55	0.42
1:F:82:LYS:O	1:F:84:PHE:N	2.53	0.42
1:H:84:PHE:HB2	1:H:87:PHE:CB	2.50	0.42
1:H:120:PRO:HG3	1:H:185:ILE:HD13	2.02	0.42
1:E:6:ILE:HD12	1:E:203:ILE:HG12	2.02	0.41
1:E:300:ILE:O	1:E:304:LYS:HG2	2.20	0.41
1:G:84:PHE:HD2	1:G:86:LYS:HD3	1.84	0.41
1:H:50:VAL:O	1:H:54:VAL:HG23	2.20	0.41
1:H:138:PRO:O	1:H:139:ASN:C	2.61	0.41
1:H:142:ARG:HD2	1:H:145:LYS:NZ	2.35	0.41
1:A:69:ASP:O	1:A:72:LYS:HG2	2.20	0.41
1:D:234:LEU:HD13	1:F:141:LYS:HZ1	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:TYR:HB2	1:C:233:ILE:CD1	2.50	0.41
1:D:307:LEU:HD23	1:D:307:LEU:HA	1.90	0.41
1:D:314:ASN:ND2	2:D:408:HOH:O	2.43	0.41
1:G:181:ASP:CB	1:G:252:MET:HB2	2.50	0.41
1:G:227:ASP:OD1	1:G:227:ASP:C	2.63	0.41
1:H:84:PHE:HB2	1:H:87:PHE:CD1	2.56	0.41
1:E:83:PRO:HB2	1:E:86:LYS:CB	2.50	0.41
1:H:325:TRP:NE1	1:H:330:GLN:OE1	2.53	0.41
1:D:96:SER:O	1:D:100:VAL:HG23	2.21	0.41
1:F:149:VAL:HG23	2:F:427:HOH:O	2.20	0.41
1:F:263:PHE:HA	1:F:264:PRO:HD3	1.79	0.41
1:H:184:LEU:HD22	1:H:255:MET:HE3	2.03	0.41
1:B:164:ARG:HG2	1:B:173:VAL:HG22	2.03	0.41
1:F:117:VAL:HG22	1:F:134:THR:HG23	2.03	0.41
1:F:216:ASN:CG	1:F:219:THR:HG23	2.46	0.41
1:H:271:ASP:O	1:H:274:ILE:HG12	2.20	0.41
1:H:92:TYR:OH	1:H:285:ASP:OD2	2.21	0.41
1:H:122:ILE:O	1:H:172:TYR:HA	2.21	0.41
1:B:3:LYS:HD3	1:B:202:HIS:CE1	2.56	0.41
1:B:327:ASP:HA	1:B:331:LYS:HD2	2.03	0.41
1:E:91:LYS:HD3	1:E:92:TYR:CE2	2.56	0.41
1:F:48:THR:HG21	1:F:176:GLY:CA	2.50	0.41
1:G:48:THR:HG22	1:G:120:PRO:HG2	2.02	0.41
1:B:313:ARG:NH2	1:E:326:PHE:HZ	2.19	0.41
1:C:142:ARG:C	1:C:144:TRP:N	2.79	0.41
1:C:180:ASN:N	1:C:252:MET:HE1	2.35	0.41
1:D:235:LYS:C	1:D:238:PRO:HD2	2.46	0.41
1:F:120:PRO:HG3	1:F:185:ILE:HD13	2.03	0.41
1:G:289:ASP:O	1:G:293:GLN:HG3	2.19	0.41
1:A:38:ILE:HD13	1:A:38:ILE:HA	1.96	0.40
1:B:306:ALA:O	1:B:309:LYS:HB2	2.21	0.40
1:F:57:ALA:HB1	1:F:62:ILE:HB	2.03	0.40
1:A:235:LYS:O	1:A:238:PRO:HD2	2.21	0.40
1:F:179:ALA:CB	1:F:185:ILE:HD11	2.51	0.40
1:G:24:LYS:HD3	1:G:24:LYS:C	2.47	0.40
1:F:180:ASN:H	1:F:248:GLN:HE21	1.69	0.40
1:A:128:SER:HA	1:A:129:PRO:HD3	1.97	0.40
1:C:202:HIS:HB3	1:C:315:TYR:OH	2.21	0.40
1:F:79:PHE:HB3	1:F:92:TYR:CD1	2.56	0.40
1:C:60:TYR:HD1	1:C:113:LEU:HD23	1.86	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:141:LYS:NZ	1:F:29:GLU:OE1[1_545]	2.13	0.07
1:A:85:LEU:CD2	1:G:276:GLN:NE2[1_565]	2.15	0.05

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	331/333 (99%)	318 (96%)	12 (4%)	1 (0%)	36	35
1	B	331/333 (99%)	320 (97%)	10 (3%)	1 (0%)	36	35
1	C	331/333 (99%)	314 (95%)	16 (5%)	1 (0%)	36	35
1	D	331/333 (99%)	321 (97%)	9 (3%)	1 (0%)	36	35
1	E	331/333 (99%)	319 (96%)	11 (3%)	1 (0%)	36	35
1	F	331/333 (99%)	321 (97%)	9 (3%)	1 (0%)	36	35
1	G	331/333 (99%)	316 (96%)	14 (4%)	1 (0%)	36	35
1	H	331/333 (99%)	315 (95%)	15 (4%)	1 (0%)	36	35
All	All	2648/2664 (99%)	2544 (96%)	96 (4%)	8 (0%)	36	35

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	11	GLY
1	B	11	GLY
1	C	11	GLY
1	D	11	GLY
1	E	11	GLY
1	F	11	GLY
1	G	11	GLY
1	A	11	GLY

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	283/283 (100%)	280 (99%)	3 (1%)	65	73
1	B	283/283 (100%)	280 (99%)	3 (1%)	65	73
1	C	283/283 (100%)	279 (99%)	4 (1%)	59	66
1	D	283/283 (100%)	277 (98%)	6 (2%)	47	52
1	E	283/283 (100%)	280 (99%)	3 (1%)	65	73
1	F	283/283 (100%)	276 (98%)	7 (2%)	42	45
1	G	283/283 (100%)	278 (98%)	5 (2%)	51	58
1	H	283/283 (100%)	273 (96%)	10 (4%)	32	32
All	All	2264/2264 (100%)	2223 (98%)	41 (2%)	51	58

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	LYS
1	A	209	LEU
1	A	248	GLN
1	B	86	LYS
1	B	198	ILE
1	B	199	GLN
1	C	38	ILE
1	C	102	GLU
1	C	196	LYS
1	C	333	VAL
1	D	1	MET
1	D	24	LYS
1	D	48	THR
1	D	141	LYS
1	D	220	LYS
1	D	289	ASP
1	E	31	LYS
1	E	90	SER
1	E	309	LYS

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Mol	Chain	Res	Type
1	F	2	GLU
1	F	3	LYS
1	F	85	LEU
1	F	86	LYS
1	F	142	ARG
1	F	212	LYS
1	F	233	ILE
1	G	1	MET
1	G	2	GLU
1	G	82	LYS
1	G	86	LYS
1	G	141	LYS
1	H	1	MET
1	H	3	LYS
1	H	31	LYS
1	H	48	THR
1	H	77	LYS
1	H	84	PHE
1	H	85	LEU
1	H	86	LYS
1	H	88	THR
1	H	221	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	250	GLN
1	A	273	GLN
1	B	214	GLN
1	B	248	GLN
1	C	214	GLN
1	C	248	GLN
1	C	250	GLN
1	C	254	GLN
1	C	273	GLN
1	C	330	GLN
1	D	35	HIS
1	D	214	GLN
1	D	248	GLN
1	D	254	GLN
1	E	28	GLN
1	E	214	GLN

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Mol	Chain	Res	Type
1	F	214	GLN
1	F	301	GLN
1	G	28	GLN
1	G	214	GLN
1	H	28	GLN
1	H	202	HIS
1	H	266	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
1	A	333/333 (100%)	0.62	12 (3%)	46	45	28, 39, 56, 83	0
1	B	333/333 (100%)	0.57	13 (3%)	43	42	26, 38, 53, 93	0
1	C	333/333 (100%)	0.67	15 (4%)	38	37	29, 41, 58, 88	0
1	D	333/333 (100%)	0.62	14 (4%)	40	39	29, 40, 55, 96	0
1	E	333/333 (100%)	0.75	13 (3%)	43	42	30, 43, 62, 84	0
1	F	333/333 (100%)	0.85	28 (8%)	17	15	31, 43, 64, 87	0
1	G	333/333 (100%)	0.87	25 (7%)	20	19	32, 44, 63, 91	0
1	H	333/333 (100%)	1.03	37 (11%)	10	9	34, 47, 67, 96	0
All	All	2664/2664 (100%)	0.75	157 (5%)	28	27	26, 42, 62, 96	0

All (157) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	87	PHE	7.1
1	F	85	LEU	6.1
1	H	85	LEU	5.4
1	F	83	PRO	5.3
1	A	84	PHE	5.0
1	H	234	LEU	5.0
1	F	84	PHE	4.9
1	C	84	PHE	4.8
1	B	84	PHE	4.7
1	C	1	MET	4.5
1	D	84	PHE	4.2
1	C	141	LYS	4.1
1	F	141	LYS	4.1
1	C	85	LEU	3.9
1	H	86	LYS	3.9
1	D	87	PHE	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	85	LEU	3.9
1	F	1	MET	3.8
1	H	84	PHE	3.8
1	A	87	PHE	3.8
1	G	85	LEU	3.7
1	E	85	LEU	3.7
1	C	87	PHE	3.6
1	E	84	PHE	3.6
1	B	83	PRO	3.6
1	H	87	PHE	3.5
1	G	84	PHE	3.4
1	H	88	THR	3.3
1	H	231	GLY	3.3
1	G	233	ILE	3.3
1	C	86	LYS	3.3
1	E	277	ALA	3.3
1	H	226	LEU	3.2
1	D	333	VAL	3.2
1	F	233	ILE	3.2
1	F	263	PHE	3.2
1	F	138	PRO	3.2
1	A	233	ILE	3.1
1	B	85	LEU	3.1
1	D	83	PRO	3.1
1	E	38	ILE	3.1
1	H	218	SER	3.1
1	A	88	THR	3.1
1	A	333	VAL	3.1
1	H	333	VAL	3.0
1	B	198	ILE	3.0
1	F	140	LEU	3.0
1	H	83	PRO	3.0
1	G	333	VAL	2.9
1	H	198	ILE	2.9
1	C	83	PRO	2.9
1	H	1	MET	2.9
1	H	277	ALA	2.9
1	H	146	LEU	2.8
1	F	217	PRO	2.8
1	G	64	VAL	2.8
1	H	225	TYR	2.8
1	D	85	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	226	LEU	2.8
1	B	332	ASN	2.8
1	F	280	GLN	2.8
1	G	1	MET	2.8
1	H	217	PRO	2.7
1	H	140	LEU	2.7
1	B	38	ILE	2.7
1	F	86	LYS	2.7
1	C	142	ARG	2.7
1	G	311	PHE	2.7
1	G	83	PRO	2.6
1	H	138	PRO	2.6
1	G	88	THR	2.6
1	E	278	SER	2.6
1	F	232	SER	2.6
1	C	233	ILE	2.6
1	F	225	TYR	2.6
1	D	1	MET	2.6
1	C	291	ALA	2.6
1	D	234	LEU	2.6
1	D	65	GLU	2.5
1	G	278	SER	2.5
1	H	2	GLU	2.5
1	E	87	PHE	2.5
1	F	87	PHE	2.5
1	G	89	GLY	2.5
1	H	148	ILE	2.5
1	H	233	ILE	2.5
1	H	79	PHE	2.5
1	F	274	ILE	2.5
1	B	87	PHE	2.4
1	B	28	GLN	2.4
1	F	198	ILE	2.4
1	E	1	MET	2.4
1	G	307	LEU	2.4
1	A	86	LYS	2.4
1	H	169	PRO	2.4
1	A	276	GLN	2.4
1	H	89	GLY	2.4
1	E	83	PRO	2.4
1	H	325	TRP	2.4
1	H	215	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	114	LYS	2.3
1	G	86	LYS	2.3
1	H	100	VAL	2.3
1	E	89	GLY	2.3
1	H	223	GLY	2.3
1	B	48	THR	2.3
1	B	36	ASP	2.3
1	F	63	LYS	2.3
1	G	304	LYS	2.3
1	H	278	SER	2.3
1	H	281	PHE	2.3
1	F	333	VAL	2.3
1	C	88	THR	2.3
1	C	277	ALA	2.3
1	A	300	ILE	2.2
1	G	274	ILE	2.2
1	F	190	ALA	2.2
1	B	225	TYR	2.2
1	F	88	THR	2.2
1	D	222	ASP	2.2
1	H	224	GLY	2.2
1	F	234	LEU	2.2
1	F	48	THR	2.2
1	H	263	PHE	2.2
1	G	226	LEU	2.2
1	G	92	TYR	2.1
1	C	160	THR	2.1
1	G	219	THR	2.1
1	D	233	ILE	2.1
1	H	194	PHE	2.1
1	G	218	SER	2.1
1	B	88	THR	2.1
1	F	160	THR	2.1
1	H	142	ARG	2.1
1	C	333	VAL	2.1
1	A	278	SER	2.1
1	D	273	GLN	2.1
1	G	276	GLN	2.1
1	E	88	THR	2.1
1	H	264	PRO	2.1
1	D	38	ILE	2.1
1	H	274	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	333	VAL	2.1
1	F	326	PHE	2.1
1	G	305	VAL	2.1
1	A	65	GLU	2.0
1	E	320	SER	2.1
1	G	39	ALA	2.1
1	F	227	ASP	2.0
1	G	36	ASP	2.0
1	D	86	LYS	2.0
1	C	38	ILE	2.0
1	E	300	ILE	2.0
1	G	38	ILE	2.0
1	F	218	SER	2.0
1	B	106	GLY	2.0
1	D	225	TYR	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.