



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 05:02 am BST

PDB ID : 4ZRK
Title : Merlin-FERM and Lats1 complex
Authors : Lin, Z.; Li, Y.; Wei, Z.; Zhang, M.
Deposited on : 2015-05-12
Resolution : 2.32 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

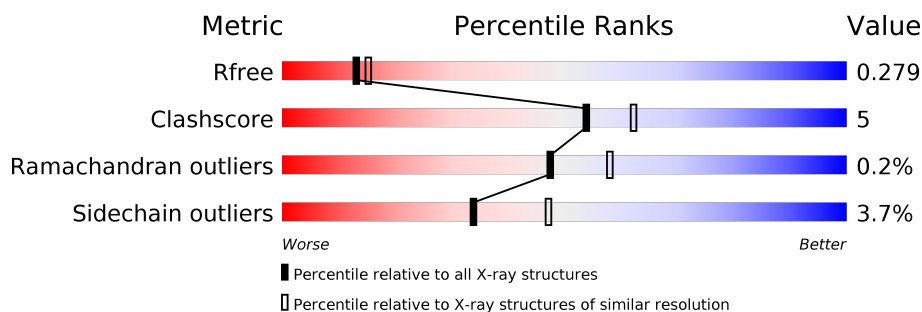
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.32 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}	130704	5974 (2.34-2.30)
Clashscore	141614	6604 (2.34-2.30)
Ramachandran outliers	138981	6523 (2.34-2.30)
Sidechain outliers	138945	6523 (2.34-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	324	<div> <div>76%</div> <div>13%</div> <div>•</div> <div>10%</div> </div>
1	B	324	<div> <div>74%</div> <div>16%</div> <div>•</div> <div>10%</div> </div>
1	C	324	<div> <div>75%</div> <div>15%</div> <div>•</div> <div>9%</div> </div>
1	D	324	<div> <div>71%</div> <div>19%</div> <div>•</div> <div>10%</div> </div>
2	E	32	<div> <div>59%</div> <div>6%</div> <div>34%</div> </div>
2	F	32	<div> <div>53%</div> <div>13%</div> <div>34%</div> </div>
2	G	32	<div> <div>59%</div> <div>9%</div> <div>31%</div> </div>

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Mol	Chain	Length	Quality of chain
2	H	32	 66% 34%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10303 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 Merlin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	1	0
			2395	1553	401	428	13			
1	B	293	Total	C	N	O	S	0	1	0
			2382	1545	394	431	12			
1	C	295	Total	C	N	O	S	0	0	0
			2423	1571	406	433	13			
1	D	293	Total	C	N	O	S	0	0	0
			2393	1555	396	429	13			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P46662
A	-2	PRO	-	expression tag	UNP P46662
A	-1	GLY	-	expression tag	UNP P46662
A	0	SER	-	expression tag	UNP P46662
B	-3	GLY	-	expression tag	UNP P46662
B	-2	PRO	-	expression tag	UNP P46662
B	-1	GLY	-	expression tag	UNP P46662
B	0	SER	-	expression tag	UNP P46662
C	-3	GLY	-	expression tag	UNP P46662
C	-2	PRO	-	expression tag	UNP P46662
C	-1	GLY	-	expression tag	UNP P46662
C	0	SER	-	expression tag	UNP P46662
D	-3	GLY	-	expression tag	UNP P46662
D	-2	PRO	-	expression tag	UNP P46662
D	-1	GLY	-	expression tag	UNP P46662
D	0	SER	-	expression tag	UNP P46662

- Molecule 2 is a protein called Serine/threonine-protein kinase LATS1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	21	Total	C	N	O	0	0	0
			170	111	32	27			
2	F	21	Total	C	N	O	0	0	0
			166	108	31	27			
2	G	22	Total	C	N	O	0	0	0
			175	114	33	28			
2	H	21	Total	C	N	O	0	0	0
			166	108	31	27			

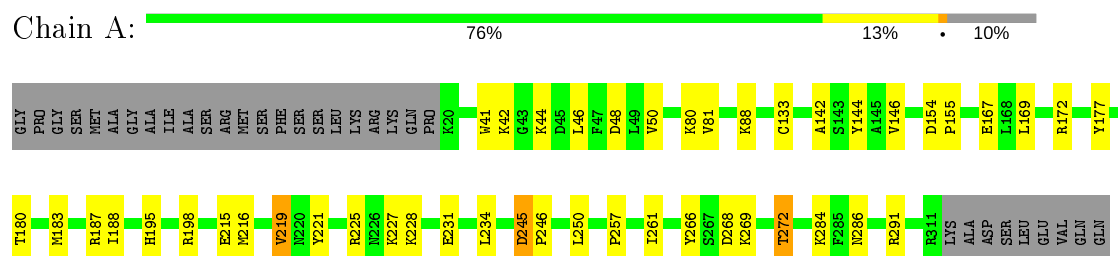
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	6	Total	O	0	0
			6	6		
3	C	7	Total	O	0	0
			7	7		
3	D	8	Total	O	0	0
			8	8		
3	G	1	Total	O	0	0
			1	1		

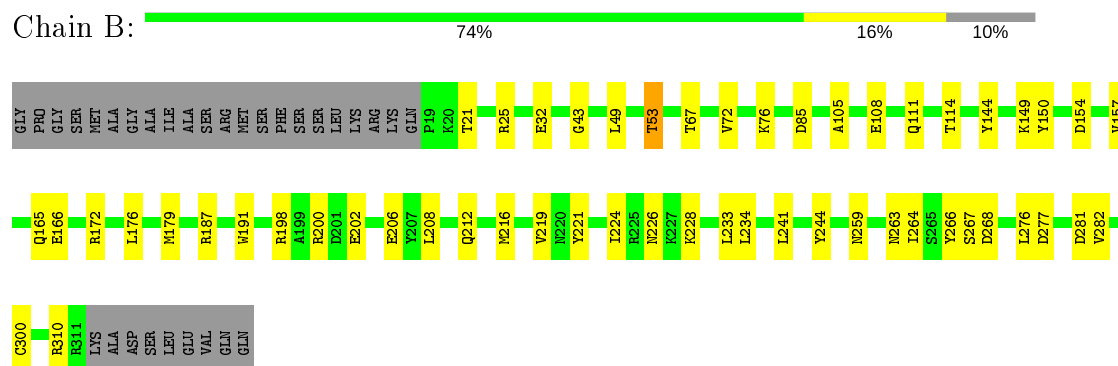
3 Residue-property plots [i](#)

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

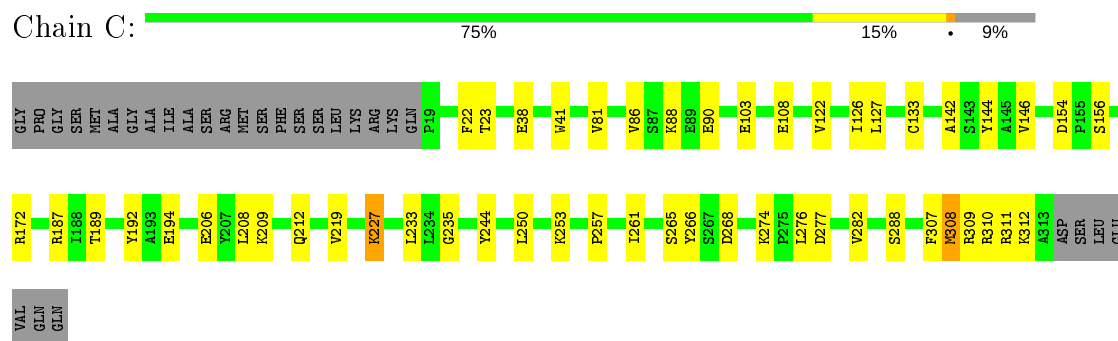
• Molecule 1: Merlin



• Molecule 1: Merlin

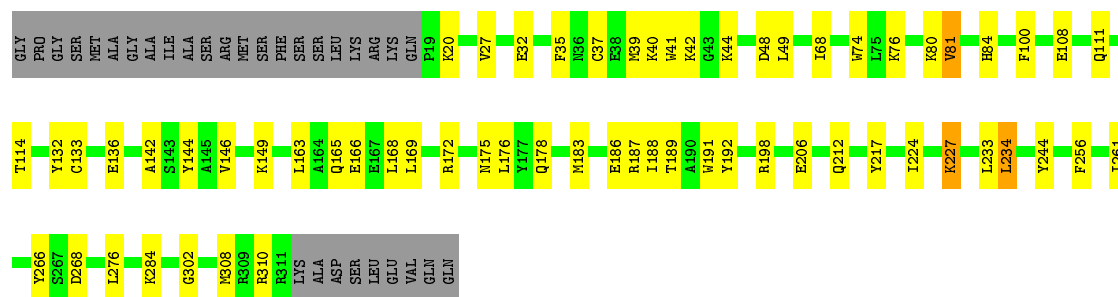


• Molecule 1: Merlin



• Molecule 1: Merlin





- Molecule 2: Serine/threonine-protein kinase LATS1



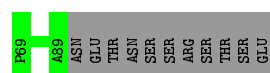
- Molecule 2: Serine/threonine-protein kinase LATS1



- Molecule 2: Serine/threonine-protein kinase LATS1



- Molecule 2: Serine/threonine-protein kinase LATS1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.05Å 110.77Å 168.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.78 – 2.32 42.78 – 2.32	Depositor EDS
% Data completeness (in resolution range)	97.9 (42.78-2.32) 97.9 (42.78-2.32)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.34 (at 2.32Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.246 , 0.279 0.246 , 0.279	Depositor DCC
R_{free} test set	4477 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	56.4	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.287 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10303	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.82% 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	0.47	0/2457	0.60	0/3324
1	B	0.49	0/2445	0.61	0/3315
1	C	0.50	1/2483 (0.0%)	0.62	0/3356
1	D	0.47	1/2453 (0.0%)	0.63	1/3320 (0.0%)
2	E	0.44	0/175	0.60	0/234
2	F	0.41	0/171	0.69	0/230
2	G	0.40	0/180	0.62	0/241
2	H	0.42	0/171	0.58	0/230
All	All	0.48	2/10535 (0.0%)	0.61	1/14250 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	133	CYS	CB-SG	-8.16	1.68	1.82
1	D	133	CYS	CB-SG	-5.38	1.73	1.81

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	234	LEU	CA-CB-CG	5.75	128.53	115.30

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	2395	0	2353	23	0
1	B	2382	0	2309	29	0
1	C	2423	0	2394	25	0
1	D	2393	0	2354	30	0
2	E	170	0	174	1	0
2	F	166	0	163	1	0
2	G	175	0	176	4	0
2	H	166	0	163	0	0
3	A	11	0	0	0	0
3	B	6	0	0	1	0
3	C	7	0	0	0	0
3	D	8	0	0	0	0
3	G	1	0	0	0	0
All	All	10303	0	10086	108	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 5.

All (108) 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:180:THR:H	1:A:183:MET:HE3	1.31	0.93
1:C:227:LYS:H	1:C:227:LYS:HD2	1.47	0.79
1:A:44:LYS:NZ	1:A:48:ASP:OD2	2.22	0.71
1:D:224:ILE:HD13	1:D:234:LEU:HB2	1.75	0.67
1:C:23:THR:OG1	1:C:90:GLU:OE2	2.13	0.67
1:D:187:ARG:NH2	1:D:191:TRP:HE1	1.92	0.66
1:D:42:LYS:HA	1:D:80:LYS:HA	1.79	0.65
1:A:169:LEU:HD21	1:A:188:ILE:HD11	1.78	0.64
1:D:149:LYS:NZ	1:D:166:GLU:OE1	2.26	0.64
2:G:70:LYS:HD2	2:G:70:LYS:H	1.63	0.62
1:D:27:VAL:HG22	1:D:32:GLU:HG2	1.82	0.61
1:B:49:LEU:O	1:B:53:THR:HG23	2.00	0.61
1:A:272:THR:HB	1:A:284:LYS:HG2	1.82	0.59
1:C:208:LEU:O	1:C:212:GLN:HB2	2.03	0.57
1:D:20:LYS:HD3	1:D:39:MET:HG3	1.86	0.57
1:B:233:LEU:HB2	1:B:244:TYR:HB2	1.86	0.57
1:D:44:LYS:NZ	1:D:48:ASP:OD2	2.37	0.56
1:B:276:LEU:O	1:B:277:ASP:HB3	2.07	0.55
1:C:308:MET:HE3	1:C:308:MET:HA	1.88	0.54
1:C:142:ALA:O	1:C:146:VAL:HG13	2.08	0.54
1:B:187:ARG:NH2	1:B:191:TRP:HE1	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:132:TYR:CE1	1:D:172:ARG:NH2	2.76	0.54
1:B:150:TYR:OH	1:B:166:GLU:OE2	2.17	0.53
1:D:266:TYR:OH	1:D:268:ASP:O	2.15	0.53
1:C:227:LYS:CD	1:C:227:LYS:H	2.12	0.53
1:B:198:ARG:NH1	1:B:202:GLU:OE2	2.42	0.52
1:B:105:ALA:N	1:B:212:GLN:OE1	2.27	0.52
1:B:165:GLN:HA	1:B:165:GLN:OE1	2.08	0.51
1:A:266:TYR:OH	1:A:268:ASP:O	2.17	0.51
1:B:208:LEU:O	1:B:212:GLN:HB2	2.11	0.50
1:D:175:ASN:O	1:D:176:LEU:HB2	2.11	0.50
1:A:228:LYS:O	1:B:228:LYS:NZ	2.36	0.50
1:B:263:ASN:OD1	1:B:264:ILE:N	2.45	0.49
1:D:169:LEU:HD21	1:D:188:ILE:HD11	1.93	0.49
1:D:227:LYS:HD2	1:D:284:LYS:HD2	1.93	0.49
1:B:67:THR:HG22	1:B:72:VAL:HG12	1.94	0.49
1:A:215:GLU:HB2	1:A:250:LEU:HD12	1.94	0.49
1:B:266:TYR:OH	1:B:268:ASP:O	2.28	0.49
1:C:309:ARG:HA	1:C:312:LYS:HE3	1.93	0.49
1:D:108:GLU:OE1	1:D:310:ARG:HD3	2.13	0.49
1:D:144:TYR:CZ	1:D:206:GLU:HG2	2.48	0.48
1:B:154:ASP:OD2	1:B:157:VAL:HG23	2.13	0.48
1:C:154:ASP:OD2	1:C:156:SER:OG	2.32	0.48
1:C:266:TYR:OH	1:C:268:ASP:O	2.26	0.47
1:B:144:TYR:CZ	1:B:206:GLU:HG2	2.49	0.47
1:B:216:MET:O	1:B:219:VAL:HG12	2.15	0.47
1:C:189:THR:HA	1:C:192:TYR:HB3	1.97	0.47
1:C:103:GLU:HG3	1:C:310:ARG:NH2	2.30	0.47
1:C:206:GLU:HG3	2:G:74:HIS:CE1	2.50	0.47
1:A:269:LYS:O	1:A:286:ASN:ND2	2.47	0.47
1:B:108:GLU:OE1	1:B:310:ARG:HD3	2.14	0.46
1:A:219:VAL:HG22	1:A:221:TYR:CE1	2.51	0.46
1:D:100:PHE:CZ	1:D:302:GLY:HA3	2.49	0.46
1:A:133:CYS:O	1:A:172[B]:ARG:NH2	2.48	0.46
1:B:224:ILE:HD13	1:B:234:LEU:HB2	1.98	0.46
1:A:227:LYS:HA	1:A:227:LYS:HD3	1.68	0.46
1:C:144:TYR:CZ	1:C:206:GLU:HG2	2.51	0.46
1:B:111:GLN:HB2	1:B:114:THR:OG1	2.17	0.45
1:D:111:GLN:HB2	1:D:114:THR:OG1	2.16	0.45
1:C:122:VAL:O	1:C:126:ILE:HG12	2.16	0.45
1:A:46:LEU:O	1:A:50:VAL:HG23	2.16	0.45
1:A:154:ASP:HA	1:A:155:PRO:HD2	1.84	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:256:PHE:HB3	1:D:261:ILE:HD11	1.99	0.45
1:B:25:ARG:HH21	1:B:32:GLU:CD	2.20	0.45
1:B:85:ASP:OD1	1:B:85:ASP:N	2.47	0.44
1:B:241:LEU:HD21	1:B:300:CYS:SG	2.57	0.44
1:A:142:ALA:O	1:A:146:VAL:HG13	2.17	0.44
1:D:142:ALA:O	1:D:146:VAL:HG13	2.17	0.44
1:D:74:TRP:O	1:D:76:LYS:NZ	2.51	0.44
1:D:37:CYS:SG	1:D:41:TRP:HB2	2.57	0.44
1:C:206:GLU:HG3	2:G:74:HIS:HE1	1.82	0.44
1:D:163:LEU:HB3	1:D:168:LEU:HD11	1.99	0.44
1:C:276:LEU:O	1:C:277:ASP:HB3	2.18	0.43
2:F:86:LEU:HB3	2:F:87:PRO:HD3	2.00	0.43
1:D:233:LEU:HB2	1:D:244:TYR:HB2	2.00	0.43
1:B:219:VAL:HG13	1:B:221:TYR:CE2	2.53	0.43
1:A:257:PRO:O	1:A:261:ILE:HG13	2.18	0.43
1:D:266:TYR:CZ	1:D:268:ASP:O	2.72	0.43
1:C:108:GLU:OE1	1:C:310:ARG:HD3	2.18	0.43
1:B:43:GLY:HA3	1:B:76:LYS:O	2.19	0.43
1:D:183:MET:O	1:D:186:GLU:HB3	2.18	0.43
1:C:307:PHE:O	1:C:311:ARG:HG2	2.19	0.43
1:A:216:MET:HG3	1:A:250:LEU:HD13	2.01	0.42
1:D:189:THR:O	1:D:192:TYR:HB3	2.19	0.42
1:A:42:LYS:HA	1:A:80:LYS:HA	2.01	0.42
1:B:172:ARG:O	1:B:176:LEU:HD13	2.19	0.42
1:B:149:LYS:NZ	1:B:166:GLU:OE1	2.24	0.42
1:A:144:TYR:CE1	1:A:195:HIS:CE1	3.08	0.42
1:A:41:TRP:O	1:A:81:VAL:HG23	2.20	0.42
1:B:187:ARG:HA	1:B:187:ARG:HD2	1.81	0.42
1:C:257:PRO:O	1:C:261:ILE:HG13	2.20	0.41
1:A:225:ARG:HG2	1:A:231:GLU:HA	2.02	0.41
1:C:38:GLU:HB2	1:C:41:TRP:CD1	2.54	0.41
1:D:40:LYS:O	1:D:80:LYS:HE2	2.20	0.41
1:C:253:LYS:HE3	1:C:253:LYS:HB3	1.70	0.41
1:D:136:GLU:N	1:D:136:GLU:OE1	2.41	0.41
1:A:144:TYR:CZ	1:A:195:HIS:CE1	3.08	0.41
1:C:194:GLU:OE1	2:G:70:LYS:HD3	2.21	0.41
1:B:226:ASN:HB2	3:B:403:HOH:O	2.19	0.41
1:C:22:PHE:HZ	1:C:81:VAL:CG2	2.32	0.41
1:D:81:VAL:HA	1:D:84:HIS:HD1	1.86	0.41
1:D:212:GLN:HA	1:D:217:TYR:CG	2.56	0.41
1:C:233:LEU:HB2	1:C:244:TYR:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:ASN:OD1	1:B:259:ASN:N	2.50	0.40
1:C:235:GLY:HA3	1:C:244:TYR:HE1	1.86	0.40
1:A:177:TYR:HD2	2:E:88:PHE:HB2	1.86	0.40
1:A:245:ASP:HA	1:A:246:PRO:HD3	1.89	0.40
1:D:35:PHE:CE1	1:D:49:LEU:HG	2.56	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	291/324 (90%)	288 (99%)	2 (1%)	1 (0%)	41	50
1	B	292/324 (90%)	286 (98%)	5 (2%)	1 (0%)	41	50
1	C	293/324 (90%)	288 (98%)	5 (2%)	0	100	100
1	D	291/324 (90%)	284 (98%)	7 (2%)	0	100	100
2	E	19/32 (59%)	18 (95%)	1 (5%)	0	100	100
2	F	19/32 (59%)	17 (90%)	2 (10%)	0	100	100
2	G	20/32 (62%)	19 (95%)	1 (5%)	0	100	100
2	H	19/32 (59%)	18 (95%)	1 (5%)	0	100	100
All	All	1244/1424 (87%)	1218 (98%)	24 (2%)	2 (0%)	47	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	281	ASP
1	A	167	GLU

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	253/291 (87%)	245 (97%)	8 (3%)	39	53
1	B	250/291 (86%)	244 (98%)	6 (2%)	49	65
1	C	257/291 (88%)	243 (95%)	14 (5%)	22	30
1	D	254/291 (87%)	246 (97%)	8 (3%)	40	55
2	E	18/29 (62%)	17 (94%)	1 (6%)	21	29
2	F	17/29 (59%)	15 (88%)	2 (12%)	5	5
2	G	18/29 (62%)	17 (94%)	1 (6%)	21	29
2	H	17/29 (59%)	17 (100%)	0	100	100
All	All	1084/1280 (85%)	1044 (96%)	40 (4%)	34	47

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	LYS
1	A	187	ARG
1	A	198	ARG
1	A	219	VAL
1	A	234	LEU
1	A	245	ASP
1	A	272	THR
1	A	291	ARG
1	B	21	THR
1	B	53	THR
1	B	179	MET
1	B	200	ARG
1	B	267	SER
1	B	282	VAL
1	C	86	VAL
1	C	88	LYS
1	C	127	LEU
1	C	172	ARG
1	C	187	ARG

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Mol	Chain	Res	Type
1	C	209	LYS
1	C	219	VAL
1	C	227	LYS
1	C	250	LEU
1	C	265	SER
1	C	274	LYS
1	C	282	VAL
1	C	288	SER
1	C	308	MET
1	D	68	ILE
1	D	81	VAL
1	D	165	GLN
1	D	178	GLN
1	D	198	ARG
1	D	227	LYS
1	D	276	LEU
1	D	308	MET
2	E	86	LEU
2	F	69	PRO
2	F	79	GLN
2	G	86	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	G	79	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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 ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.