



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 09:22 AM GMT

PDB ID : 1RD4
Title : An allosteric inhibitor of LFA-1 bound to its I-domain
Authors : Crump, M.P.; Ceska, T.A.; Spyropoulos, L.; Henry, A.; Archibald, S.C.;
Alexander, R.; Taylor, R.J.; Findlow, S.C.; O'Connell, J.; Robinson, M.K.;
Shock, A.
Deposited on : 2003-11-05
Resolution : 2.40 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

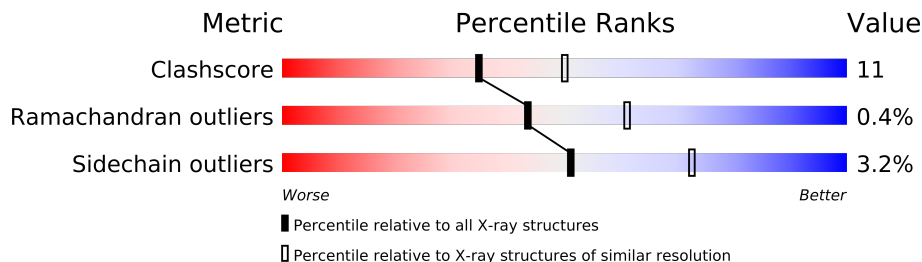
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	191	
1	B	191	
1	C	191	
1	D	191	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6052 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

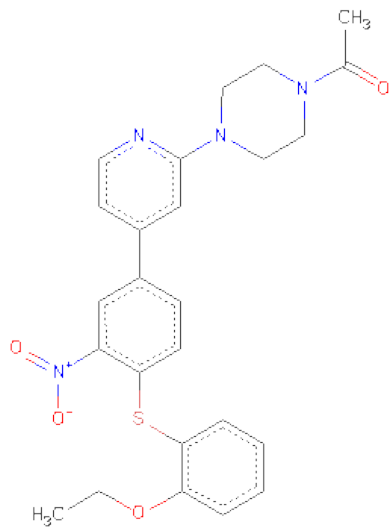
- Molecule 1 is a protein called Integrin alpha-L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	184	Total	C	N	O	S	0	0	0
			1479	956	236	283	4			
1	B	184	Total	C	N	O	S	0	0	0
			1479	956	236	283	4			
1	C	184	Total	C	N	O	S	0	0	0
			1479	956	236	283	4			
1	D	184	Total	C	N	O	S	0	0	0
			1479	956	236	283	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	121	GLY	-	CLONING ARTIFACT	UNP P20701
A	122	ALA	-	CLONING ARTIFACT	UNP P20701
A	123	MET	-	CLONING ARTIFACT	UNP P20701
A	124	SER	-	CLONING ARTIFACT	UNP P20701
B	121	GLY	-	CLONING ARTIFACT	UNP P20701
B	122	ALA	-	CLONING ARTIFACT	UNP P20701
B	123	MET	-	CLONING ARTIFACT	UNP P20701
B	124	SER	-	CLONING ARTIFACT	UNP P20701
C	121	GLY	-	CLONING ARTIFACT	UNP P20701
C	122	ALA	-	CLONING ARTIFACT	UNP P20701
C	123	MET	-	CLONING ARTIFACT	UNP P20701
C	124	SER	-	CLONING ARTIFACT	UNP P20701
D	121	GLY	-	CLONING ARTIFACT	UNP P20701
D	122	ALA	-	CLONING ARTIFACT	UNP P20701
D	123	MET	-	CLONING ARTIFACT	UNP P20701
D	124	SER	-	CLONING ARTIFACT	UNP P20701

- Molecule 2 is 1-ACETYL-4-(4-{4-[(2-ETHOXYPHENYL)THIO]-3-NITROPHENYL}PYRIDIN-2-YL)PIPERAZINE (three-letter code: L08) (formula: C₂₅H₂₆N₄O₄S).



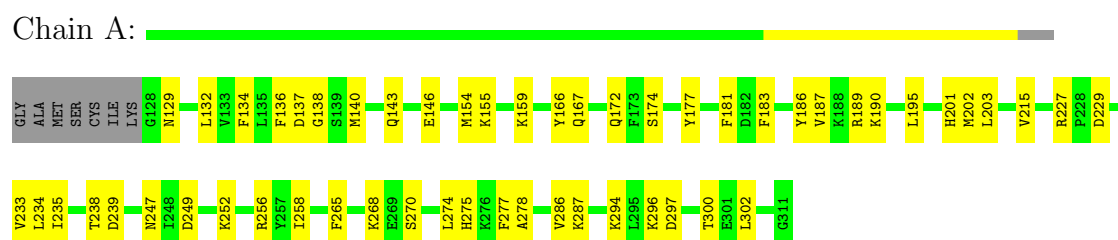
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	B	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	C	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	D	1	Total	C	N	O	S	0	0
			34	25	4	4	1		

3 Residue-property plots

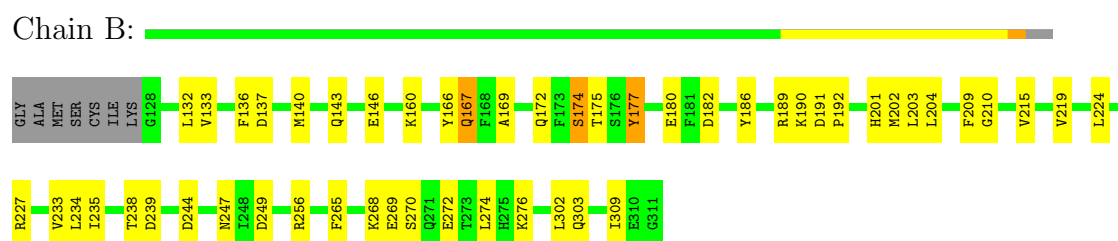
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

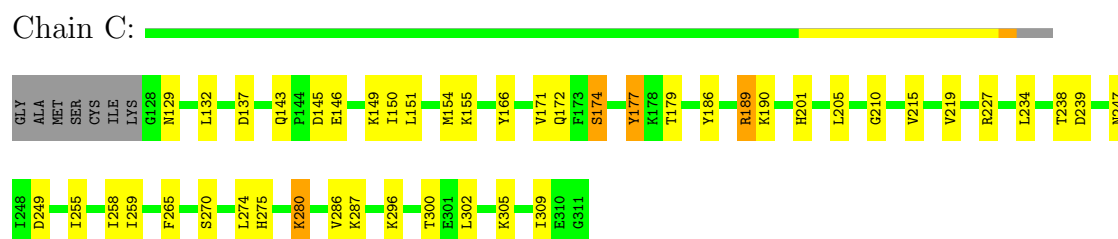
• Molecule 1: Integrin alpha-L



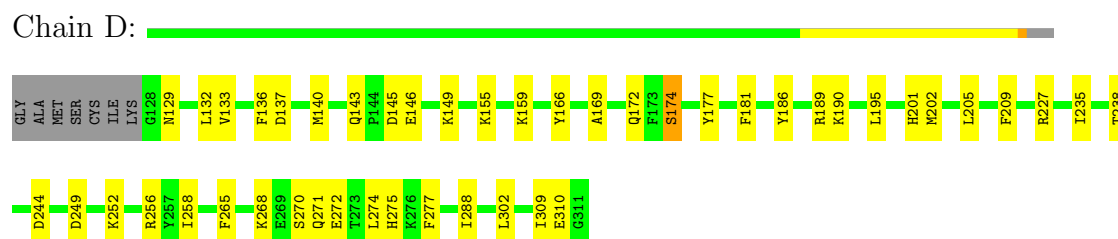
• Molecule 1: Integrin alpha-L



• Molecule 1: Integrin alpha-L



• Molecule 1: Integrin alpha-L



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.96Å 64.41Å 66.21Å 74.21° 90.00° 87.26°	Depositor
Resolution (Å)	40.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (40.00-2.40)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.236 , 0.290	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6052	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: L08

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	0.45	0/1507	0.60	0/2026
1	B	0.45	0/1507	0.58	0/2026
1	C	0.45	0/1507	0.60	0/2026
1	D	0.45	0/1507	0.59	0/2026
All	All	0.45	0/6028	0.59	0/8104

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1479	0	1487	35	0
1	B	1479	0	1487	34	0
1	C	1479	0	1487	35	0
1	D	1479	0	1487	26	0
2	A	34	0	26	2	0
2	B	34	0	26	3	0
2	C	34	0	26	5	0
2	D	34	0	26	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6052	0	6052	130	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (130) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:132:LEU:HD11	1:A:235:ILE:HD11	1.46	0.96
1:C:309:ILE:HD13	1:D:309:ILE:HD13	1.48	0.95
1:B:132:LEU:HD21	1:B:235:ILE:HD12	1.64	0.78
1:D:268:LYS:O	1:D:272:GLU:HG2	1.83	0.78
1:B:132:LEU:HD11	1:B:235:ILE:HD11	1.64	0.77
1:B:247:ASN:HD22	1:B:249:ASP:H	1.37	0.72
1:A:155:LYS:O	1:A:159:LYS:HG3	1.90	0.72
1:A:132:LEU:HD21	1:A:235:ILE:HD12	1.73	0.70
1:B:160:LYS:HE2	1:B:303:GLN:HE21	1.57	0.70
1:A:167:GLN:HG3	1:A:227:ARG:NH2	2.07	0.69
1:C:305:LYS:HD3	1:D:310:GLU:HB3	1.75	0.68
1:A:137:ASP:HB2	1:A:238:THR:HA	1.76	0.68
1:B:137:ASP:HB2	1:B:238:THR:HA	1.75	0.68
1:A:249:ASP:O	1:A:252:LYS:HG3	1.94	0.67
1:C:166:TYR:CE2	2:C:2328:L08:H35	2.28	0.67
1:D:132:LEU:HD21	1:D:235:ILE:HD12	1.77	0.66
1:C:137:ASP:HB2	1:C:238:THR:HA	1.79	0.65
1:A:132:LEU:HD11	1:A:235:ILE:CD1	2.26	0.64
1:D:137:ASP:HB2	1:D:238:THR:HA	1.79	0.64
1:D:132:LEU:HD11	1:D:235:ILE:HD11	1.79	0.63
1:B:174:SER:HB3	1:B:202:MET:HE1	1.83	0.60
1:C:258:ILE:HG22	1:C:286:VAL:HG22	1.84	0.60
1:B:133:VAL:HG22	1:B:169:ALA:HB3	1.84	0.60
1:B:247:ASN:ND2	1:B:249:ASP:H	2.00	0.59
1:B:132:LEU:HD11	1:B:235:ILE:CD1	2.33	0.58
1:C:215:VAL:HG11	1:C:234:LEU:HD13	1.85	0.58
1:C:296:LYS:O	1:C:300:THR:HG23	2.03	0.58
1:C:258:ILE:CG2	1:C:286:VAL:HG22	2.34	0.57
1:C:265:PHE:CE2	1:C:274:LEU:HD11	2.39	0.57
1:B:177:TYR:OH	1:B:210:GLY:HA3	2.05	0.57
1:B:167:GLN:HB2	1:B:227:ARG:NH2	2.19	0.57
1:D:258:ILE:HD13	1:D:274:LEU:HB3	1.87	0.56
1:C:247:ASN:HD22	1:C:249:ASP:H	1.52	0.56
1:A:189:ARG:C	1:A:190:LYS:HD3	2.25	0.56
1:A:167:GLN:HG3	1:A:227:ARG:CZ	2.36	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:280:LYS:N	1:C:280:LYS:HD2	2.21	0.56
1:D:249:ASP:O	1:D:252:LYS:HG3	2.06	0.55
1:D:129:ASN:HB3	1:D:227:ARG:CZ	2.36	0.55
1:D:189:ARG:C	1:D:190:LYS:HD3	2.27	0.55
1:C:302:LEU:HD23	1:C:302:LEU:C	2.27	0.54
1:B:234:LEU:HD23	1:B:256:ARG:HG2	1.89	0.54
1:A:233:VAL:HG21	2:A:328:L08:H191	1.90	0.54
1:A:296:LYS:O	1:A:300:THR:HG23	2.09	0.53
1:A:265:PHE:CZ	1:A:274:LEU:HD11	2.43	0.53
1:A:166:TYR:CE2	2:A:328:L08:H35	2.44	0.52
1:B:265:PHE:CZ	1:B:274:LEU:HD11	2.45	0.52
1:A:138:GLY:HA3	1:A:172:GLN:HE21	1.75	0.52
1:B:180:GLU:OE2	1:B:180:GLU:HA	2.10	0.52
1:D:145:ASP:O	1:D:149:LYS:HG3	2.09	0.51
1:B:233:VAL:HG21	2:B:1328:L08:H191	1.92	0.51
1:C:143:GLN:HB2	1:C:146:GLU:HG3	1.91	0.51
1:D:181:PHE:CD1	1:D:195:LEU:HB3	2.46	0.51
1:D:174:SER:HB3	1:D:202:MET:HE1	1.93	0.51
1:B:172:GLN:HE22	1:B:201:HIS:HA	1.77	0.50
1:D:133:VAL:HG22	1:D:169:ALA:HB3	1.95	0.49
1:D:132:LEU:HD11	1:D:235:ILE:CD1	2.42	0.49
1:B:136:PHE:CZ	1:B:172:GLN:HB2	2.48	0.49
1:D:143:GLN:HG3	1:D:146:GLU:OE1	2.12	0.49
1:C:166:TYR:HE2	2:C:2328:L08:H35	1.73	0.49
1:A:265:PHE:HA	1:A:270:SER:OG	2.13	0.48
1:B:175:THR:OG1	1:B:204:LEU:HB3	2.14	0.48
2:D:3328:L08:O3	2:D:3328:L08:S1	2.71	0.48
1:C:172:GLN:HE22	1:C:201:HIS:HA	1.79	0.48
1:C:265:PHE:HA	1:C:270:SER:OG	2.14	0.48
1:D:271:GLN:HG2	1:D:288:ILE:HD12	1.94	0.48
1:C:150:ILE:O	1:C:154:MET:HG3	2.14	0.48
1:D:172:GLN:HE22	1:D:201:HIS:HA	1.78	0.47
1:C:177:TYR:OH	1:C:210:GLY:HA3	2.14	0.47
1:A:258:ILE:HG22	1:A:278:ALA:HB2	1.96	0.47
1:C:151:LEU:HD23	1:C:154:MET:CE	2.45	0.47
1:A:172:GLN:HE22	1:A:201:HIS:HA	1.78	0.47
1:B:265:PHE:HA	1:B:270:SER:OG	2.15	0.47
1:B:166:TYR:CE2	2:B:1328:L08:H35	2.50	0.47
1:D:256:ARG:HD2	1:D:277:PHE:O	2.15	0.47
1:D:166:TYR:CE2	2:D:3328:L08:H35	2.50	0.47
1:C:129:ASN:HB3	1:C:227:ARG:CZ	2.45	0.47
1:B:268:LYS:O	1:B:272:GLU:HG2	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:215:VAL:HA	1:B:219:VAL:HG23	1.97	0.46
2:B:1328:L08:S1	2:B:1328:L08:O3	2.73	0.46
1:A:174:SER:HB2	1:A:202:MET:HE1	1.98	0.46
1:D:209:PHE:HB2	1:D:244:ASP:OD2	2.16	0.46
1:B:137:ASP:CB	1:B:238:THR:HA	2.43	0.46
1:C:215:VAL:HA	1:C:219:VAL:HG23	1.97	0.46
1:B:174:SER:HB3	1:B:202:MET:CE	2.46	0.45
1:A:247:ASN:ND2	1:A:249:ASP:HB2	2.31	0.45
1:C:145:ASP:O	1:C:149:LYS:HG3	2.16	0.45
1:C:151:LEU:O	1:C:155:LYS:HG3	2.17	0.45
1:A:143:GLN:HB2	1:A:146:GLU:HG3	1.99	0.45
1:A:258:ILE:CG2	1:A:286:VAL:HG22	2.47	0.45
1:B:209:PHE:HB2	1:B:244:ASP:OD2	2.17	0.44
1:B:309:ILE:O	1:B:309:ILE:HD12	2.17	0.44
1:C:137:ASP:CB	1:C:238:THR:HA	2.47	0.44
1:A:132:LEU:HD23	1:A:132:LEU:C	2.38	0.44
1:A:247:ASN:ND2	1:A:249:ASP:OD2	2.50	0.44
1:A:181:PHE:CD1	1:A:195:LEU:HB3	2.52	0.44
1:A:137:ASP:CB	1:A:238:THR:HA	2.47	0.44
1:C:265:PHE:CZ	1:C:274:LEU:HD11	2.52	0.44
1:D:155:LYS:O	1:D:159:LYS:HG3	2.16	0.44
1:C:255:ILE:HD12	2:C:2328:L08:H202	1.99	0.43
1:C:174:SER:HA	1:C:205:LEU:O	2.18	0.43
1:A:136:PHE:CZ	1:A:172:GLN:HB2	2.53	0.43
1:A:140:MET:HA	1:A:203:LEU:HD23	2.00	0.43
1:C:171:VAL:HG22	1:C:179:THR:HA	2.00	0.43
1:C:215:VAL:HA	1:C:219:VAL:CG2	2.48	0.43
1:A:134:PHE:CD2	1:A:154:MET:HG2	2.54	0.43
1:C:186:TYR:O	1:C:190:LYS:N	2.51	0.43
1:B:186:TYR:O	1:B:190:LYS:N	2.50	0.42
1:A:268:LYS:HE2	1:A:268:LYS:HB3	1.89	0.42
1:C:189:ARG:O	1:C:190:LYS:HB2	2.19	0.42
1:B:215:VAL:HA	1:B:219:VAL:CG2	2.48	0.42
1:A:129:ASN:ND2	1:A:229:ASP:OD2	2.51	0.42
1:B:143:GLN:HG3	1:B:146:GLU:OE1	2.20	0.42
1:D:136:PHE:CZ	1:D:172:GLN:HB2	2.55	0.42
1:D:265:PHE:HA	1:D:270:SER:OG	2.21	0.41
1:B:189:ARG:O	1:B:190:LYS:HB2	2.20	0.41
1:C:132:LEU:C	1:C:132:LEU:HD23	2.40	0.41
1:A:186:TYR:O	1:A:190:LYS:N	2.52	0.41
1:B:182:ASP:OD1	1:B:224:LEU:HD22	2.20	0.41
1:C:259:ILE:HD11	2:C:2328:L08:S1	2.60	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:132:LEU:HD23	1:B:132:LEU:C	2.42	0.41
1:A:294:LYS:HA	1:A:297:ASP:OD1	2.21	0.41
1:B:140:MET:HB3	1:B:203:LEU:O	2.21	0.41
1:D:174:SER:HA	1:D:205:LEU:O	2.22	0.40
1:A:215:VAL:HG11	1:A:234:LEU:HD13	2.03	0.40
1:A:256:ARG:HD2	1:A:277:PHE:O	2.21	0.40
1:B:191:ASP:HA	1:B:192:PRO:HD3	1.84	0.40
1:A:183:PHE:O	1:A:187:VAL:HG23	2.22	0.40
1:C:166:TYR:CZ	2:C:2328:L08:H35	2.56	0.40
1:D:186:TYR:O	1:D:190:LYS:N	2.52	0.40
1:C:151:LEU:HD23	1:C:154:MET:HE2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	182/191 (95%)	171 (94%)	11 (6%)	0	100	100
1	B	182/191 (95%)	170 (93%)	11 (6%)	1 (0%)	38	53
1	C	182/191 (95%)	171 (94%)	10 (6%)	1 (0%)	38	53
1	D	182/191 (95%)	171 (94%)	10 (6%)	1 (0%)	38	53
All	All	728/764 (95%)	683 (94%)	42 (6%)	3 (0%)	43	61

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	174	SER
1	B	174	SER
1	C	174	SER

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	163/168 (97%)	158 (97%)	5 (3%)	52	74
1	B	163/168 (97%)	157 (96%)	6 (4%)	45	66
1	C	163/168 (97%)	157 (96%)	6 (4%)	45	66
1	D	163/168 (97%)	159 (98%)	4 (2%)	60	80
All	All	652/672 (97%)	631 (97%)	21 (3%)	51	72

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	177	TYR
1	A	239	ASP
1	A	275	HIS
1	A	287	LYS
1	A	302	LEU
1	B	167	GLN
1	B	177	TYR
1	B	239	ASP
1	B	269	GLU
1	B	276	LYS
1	B	302	LEU
1	C	177	TYR
1	C	189	ARG
1	C	239	ASP
1	C	275	HIS
1	C	280	LYS
1	C	287	LYS
1	D	140	MET
1	D	177	TYR
1	D	275	HIS
1	D	302	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	163	ASN
1	A	172	GLN
1	A	213	ASN
1	A	247	ASN
1	B	172	GLN
1	B	247	ASN
1	B	303	GLN
1	C	163	ASN
1	C	172	GLN
1	C	213	ASN
1	C	247	ASN
1	C	303	GLN
1	D	163	ASN
1	D	172	GLN
1	D	247	ASN
1	D	303	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	L08	A	328	-	37,37,37	3.45	24 (64%)	51,51,51	2.45	16 (31%)
2	L08	B	1328	-	37,37,37	3.23	23 (62%)	51,51,51	2.54	19 (37%)
2	L08	C	2328	-	37,37,37	3.40	23 (62%)	51,51,51	2.40	18 (35%)
2	L08	D	3328	-	37,37,37	3.29	22 (59%)	51,51,51	2.54	18 (35%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	L08	A	328	-	-	0/22/33/33	0/4/4/4
2	L08	B	1328	-	-	0/22/33/33	0/4/4/4
2	L08	C	2328	-	-	0/22/33/33	0/4/4/4
2	L08	D	3328	-	-	2/22/33/33	0/4/4/4

All (92) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	328	L08	C17-C18	8.71	1.53	1.39
2	C	2328	L08	C17-C18	8.34	1.52	1.39
2	D	3328	L08	C17-C18	7.90	1.52	1.39
2	B	1328	L08	C17-C18	7.45	1.51	1.39
2	D	3328	L08	C12-C11	5.96	1.52	1.39
2	C	2328	L08	C12-C11	5.89	1.52	1.39
2	A	328	L08	C12-C11	5.63	1.51	1.39
2	A	328	L08	C2-C1	5.57	1.49	1.40
2	B	1328	L08	C12-C11	5.46	1.51	1.39
2	D	3328	L08	C2-C1	5.45	1.49	1.40
2	B	1328	L08	C2-C1	5.41	1.49	1.40
2	C	2328	L08	C2-C1	5.14	1.49	1.40
2	A	328	L08	O7-C2	5.07	1.47	1.37
2	C	2328	L08	C3-C2	5.00	1.50	1.39
2	C	2328	L08	O7-C2	4.96	1.47	1.37
2	A	328	L08	C3-C2	4.88	1.50	1.39
2	A	328	L08	C15-C10	4.88	1.49	1.40
2	A	328	L08	C4-C3	4.74	1.50	1.39
2	A	328	L08	C35-N20	4.70	1.45	1.34
2	D	3328	L08	C4-C3	4.69	1.49	1.39
2	B	1328	L08	C3-C2	4.66	1.49	1.39
2	C	2328	L08	C6-C1	4.65	1.48	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3328	L08	C15-C10	4.63	1.48	1.40
2	B	1328	L08	C4-C3	4.62	1.49	1.39
2	C	2328	L08	C35-N20	4.62	1.45	1.34
2	B	1328	L08	O7-C2	4.60	1.46	1.37
2	C	2328	L08	C4-C3	4.60	1.49	1.39
2	C	2328	L08	C15-C10	4.58	1.48	1.40
2	B	1328	L08	C35-N20	4.56	1.45	1.34
2	D	3328	L08	C3-C2	4.54	1.49	1.39
2	D	3328	L08	C35-N20	4.52	1.45	1.34
2	D	3328	L08	O7-C2	4.47	1.46	1.37
2	D	3328	L08	C6-C1	4.46	1.48	1.40
2	B	1328	L08	C15-C10	4.36	1.48	1.40
2	B	1328	L08	C6-C1	4.35	1.48	1.40
2	A	328	L08	C6-C1	4.26	1.48	1.40
2	C	2328	L08	C5-C4	4.12	1.49	1.37
2	A	328	L08	C5-C4	4.10	1.49	1.37
2	C	2328	L08	C14-C13	4.04	1.48	1.39
2	B	1328	L08	C5-C4	4.01	1.49	1.37
2	D	3328	L08	C14-C13	3.94	1.47	1.39
2	D	3328	L08	C5-C4	3.90	1.49	1.37
2	A	328	L08	C14-C13	3.87	1.47	1.39
2	C	2328	L08	C11-C10	3.85	1.49	1.40
2	B	1328	L08	C14-C13	3.71	1.47	1.39
2	D	3328	L08	C12-C13	3.59	1.46	1.39
2	A	328	L08	C11-C10	3.55	1.48	1.40
2	A	328	L08	C8-C16	3.55	1.46	1.39
2	B	1328	L08	C8-C16	3.54	1.46	1.39
2	C	2328	L08	C8-C16	3.49	1.46	1.39
2	A	328	L08	C18-N1	3.48	1.44	1.37
2	A	328	L08	C21-N2	3.47	1.53	1.47
2	B	1328	L08	C12-C13	3.37	1.46	1.39
2	C	2328	L08	C12-C13	3.31	1.46	1.39
2	D	3328	L08	C11-C10	3.31	1.48	1.40
2	A	328	L08	C12-C13	3.30	1.46	1.39
2	D	3328	L08	C17-C16	3.28	1.46	1.39
2	C	2328	L08	C17-C16	3.24	1.45	1.39
2	B	1328	L08	C21-N2	3.23	1.52	1.47
2	D	3328	L08	C8-C16	3.20	1.46	1.39
2	D	3328	L08	C21-N2	3.08	1.52	1.47
2	C	2328	L08	C18-N1	3.08	1.44	1.37
2	A	328	L08	C17-C16	3.07	1.45	1.39
2	B	1328	L08	C8-C35	3.01	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3328	L08	C8-C35	2.99	1.45	1.38
2	B	1328	L08	C11-C10	2.98	1.47	1.40
2	D	3328	L08	C18-N1	2.98	1.43	1.37
2	A	328	L08	C10-S1	2.95	1.82	1.78
2	C	2328	L08	C10-S1	2.95	1.82	1.78
2	C	2328	L08	C21-N2	2.90	1.52	1.47
2	A	328	L08	C8-C35	2.86	1.44	1.38
2	B	1328	L08	C17-C16	2.81	1.45	1.39
2	C	2328	L08	C5-C6	2.74	1.45	1.39
2	C	2328	L08	C8-C35	2.74	1.44	1.38
2	A	328	L08	C5-C6	2.73	1.45	1.39
2	C	2328	L08	C14-C15	2.69	1.43	1.38
2	C	2328	L08	C20-N2	2.69	1.51	1.47
2	B	1328	L08	C14-C15	2.65	1.43	1.38
2	A	328	L08	C20-N2	2.64	1.51	1.47
2	D	3328	L08	C5-C6	2.63	1.45	1.39
2	D	3328	L08	C20-N2	2.60	1.51	1.47
2	B	1328	L08	C11-N3	-2.55	1.42	1.46
2	B	1328	L08	C5-C6	2.48	1.44	1.39
2	B	1328	L08	C20-N2	2.48	1.51	1.47
2	A	328	L08	C14-C15	2.42	1.43	1.38
2	C	2328	L08	C11-N3	-2.28	1.42	1.46
2	A	328	L08	C23-N2	2.21	1.43	1.35
2	A	328	L08	C1-S1	-2.20	1.75	1.78
2	D	3328	L08	C14-C15	2.13	1.42	1.38
2	B	1328	L08	C10-S1	2.04	1.81	1.78
2	B	1328	L08	C18-N1	2.03	1.41	1.37
2	D	3328	L08	C18-N20	-2.03	1.29	1.34

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1328	L08	C22-N1-C19	7.70	127.50	111.67
2	C	2328	L08	C22-N1-C19	7.26	126.61	111.67
2	D	3328	L08	C22-N1-C19	7.16	126.40	111.67
2	A	328	L08	C22-N1-C19	6.91	125.88	111.67
2	D	3328	L08	C21-N2-C23	-5.68	107.34	122.85
2	B	1328	L08	C12-C11-N3	5.40	121.24	115.82
2	C	2328	L08	C21-N2-C23	-5.39	108.13	122.85
2	A	328	L08	C21-N2-C23	-5.38	108.18	122.85
2	D	3328	L08	C12-C11-N3	5.34	121.18	115.82
2	B	1328	L08	C21-N2-C23	-5.23	108.58	122.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3328	L08	C12-C11-C10	-5.23	117.28	122.91
2	D	3328	L08	C35-N20-C18	5.14	123.80	116.95
2	D	3328	L08	N20-C18-N1	5.02	124.24	117.09
2	B	1328	L08	C35-N20-C18	5.01	123.63	116.95
2	A	328	L08	C12-C11-C10	-5.01	117.52	122.91
2	B	1328	L08	C20-N2-C23	-4.96	109.31	122.85
2	C	2328	L08	C35-N20-C18	4.90	123.49	116.95
2	A	328	L08	C35-N20-C18	4.84	123.41	116.95
2	B	1328	L08	N20-C18-N1	4.76	123.87	117.09
2	A	328	L08	C12-C11-N3	4.73	120.57	115.82
2	C	2328	L08	N20-C18-N1	4.69	123.77	117.09
2	C	2328	L08	C12-C11-C10	-4.69	117.86	122.91
2	C	2328	L08	C20-N2-C23	-4.68	110.06	122.85
2	A	328	L08	C20-N2-C23	-4.67	110.10	122.85
2	A	328	L08	N20-C18-N1	4.65	123.70	117.09
2	B	1328	L08	C12-C11-C10	-4.55	118.01	122.91
2	D	3328	L08	C20-N2-C23	-4.41	110.81	122.85
2	C	2328	L08	C12-C11-N3	3.93	119.77	115.82
2	D	3328	L08	O3-N3-O4	-3.93	113.40	121.35
2	A	328	L08	O3-N3-O4	-3.93	113.40	121.35
2	B	1328	L08	O3-N3-O4	-3.74	113.79	121.35
2	C	2328	L08	O3-N3-O4	-3.70	113.86	121.35
2	B	1328	L08	C17-C18-N1	-3.62	117.42	122.08
2	D	3328	L08	C17-C18-N1	-3.54	117.52	122.08
2	A	328	L08	C17-C18-N1	-3.21	117.94	122.08
2	B	1328	L08	C19-N1-C18	-3.20	112.65	120.19
2	C	2328	L08	C17-C18-N1	-3.15	118.02	122.08
2	A	328	L08	C6-C1-C2	3.13	122.63	118.64
2	B	1328	L08	C6-C1-C2	2.96	122.40	118.64
2	A	328	L08	C19-N1-C18	-2.89	113.38	120.19
2	A	328	L08	C21-C22-N1	-2.88	105.19	110.62
2	C	2328	L08	C19-N1-C18	-2.76	113.69	120.19
2	D	3328	L08	C6-C1-C2	2.74	122.13	118.64
2	D	3328	L08	C16-C17-C18	2.73	120.16	117.84
2	C	2328	L08	C16-C17-C18	2.69	120.13	117.84
2	C	2328	L08	C6-C1-C2	2.68	122.04	118.64
2	B	1328	L08	C16-C17-C18	2.67	120.12	117.84
2	D	3328	L08	C19-N1-C18	-2.61	114.03	120.19
2	C	2328	L08	C21-C22-N1	-2.59	105.73	110.62
2	B	1328	L08	C21-C22-N1	-2.57	105.78	110.62
2	D	3328	L08	C21-C22-N1	-2.55	105.82	110.62
2	A	328	L08	C16-C17-C18	2.53	119.99	117.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2328	L08	C17-C18-N20	-2.50	118.21	122.69
2	D	3328	L08	C17-C18-N20	-2.48	118.24	122.69
2	B	1328	L08	C11-C10-S1	-2.46	117.76	122.36
2	A	328	L08	C17-C18-N20	-2.42	118.35	122.69
2	A	328	L08	O3-N3-C11	2.33	124.42	117.25
2	B	1328	L08	C8-C35-N20	-2.32	121.30	123.88
2	C	2328	L08	O3-N3-C11	2.27	124.25	117.25
2	D	3328	L08	C9-O7-C2	2.25	122.75	118.02
2	B	1328	L08	C17-C18-N20	-2.22	118.71	122.69
2	D	3328	L08	O3-N3-C11	2.17	123.93	117.25
2	B	1328	L08	C17-C16-C13	-2.15	117.11	120.89
2	C	2328	L08	C9-O7-C2	2.12	122.48	118.02
2	B	1328	L08	C21-N2-C20	2.11	116.54	112.52
2	D	3328	L08	C17-C16-C13	-2.11	117.18	120.89
2	C	2328	L08	C21-N2-C20	2.10	116.51	112.52
2	C	2328	L08	C17-C16-C13	-2.08	117.24	120.89
2	A	328	L08	C13-C12-C11	2.06	122.86	119.91
2	D	3328	L08	C11-C10-S1	-2.05	118.53	122.36
2	B	1328	L08	O3-N3-C11	2.01	123.43	117.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	3328	L08	O2-C23-N2-C21
2	D	3328	L08	C24-C23-N2-C21

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.