



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 01:09 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 X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

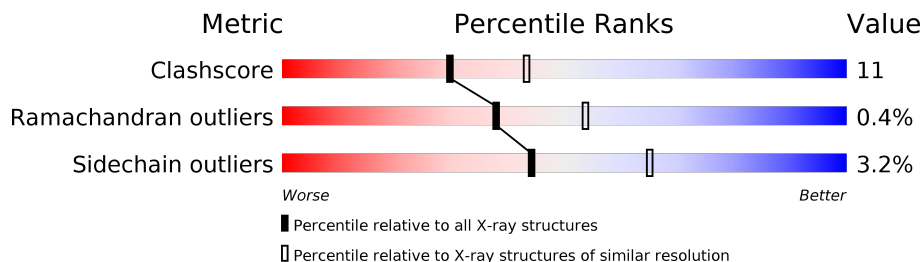
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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.

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 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.

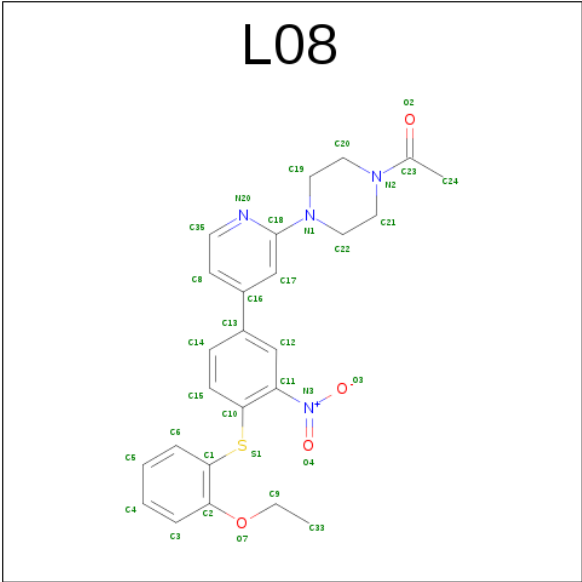
- Molecule 1 is a protein called Integrin alpha-L.

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: C<sub>25</sub>H<sub>26</sub>N<sub>4</sub>O<sub>4</sub>S).



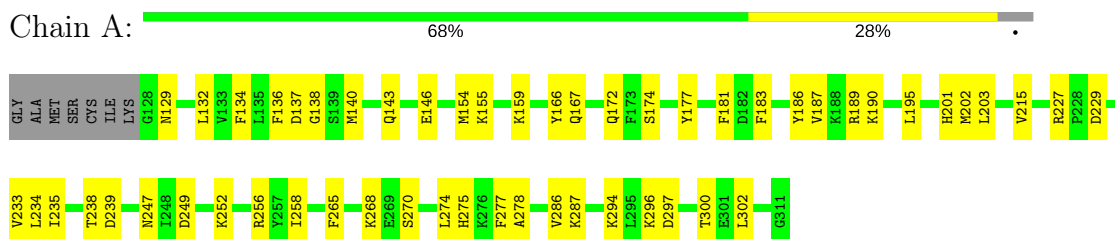
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	B	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	C	1	Total	C	N	O	S	0	0
			34	25	4	4	1		
2	D	1	Total	C	N	O	S	0	0
			34	25	4	4	1		

### 3 Residue-property plots

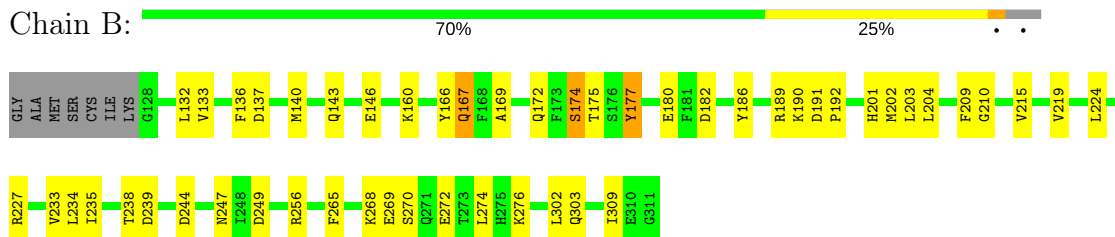
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

Note EDS was not executed.

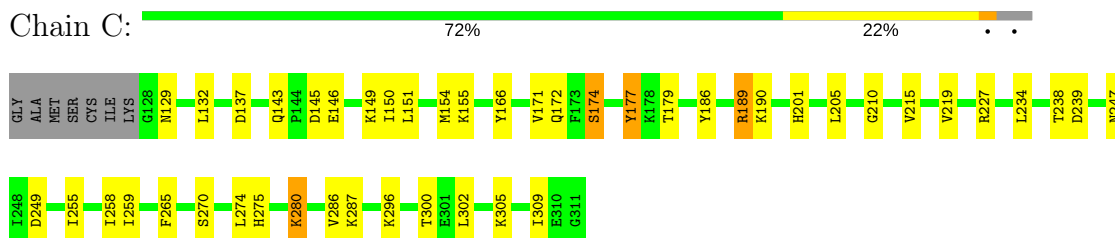
#### • Molecule 1: Integrin alpha-L



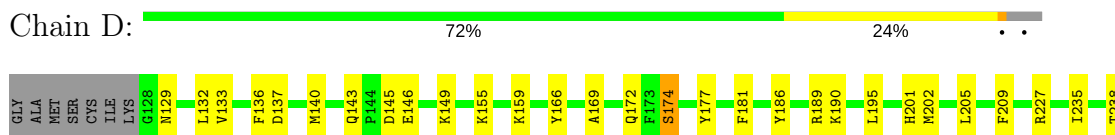
#### • Molecule 1: Integrin alpha-L



#### • Molecule 1: Integrin alpha-L



#### • Molecule 1: Integrin alpha-L



D244	D249	K252	R256	T257	L258	F265	K268	E269	S270	Q271	E272	T273	L274	H275	K276	F277	L288	L302	I309	E310	G311
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## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.96 Å   64.41 Å   66.21 Å 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 <sup>2</sup> )	14.0	wwPDB-VP

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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 Too-close contacts [i](#)

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	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
All	All	6052	0	6052	130	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 11.

All (130) 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: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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:136:PHE:CZ	1:B:172:GLN:HB2	2.48	0.49
1:D:132:LEU:HD11	1:D:235:ILE:CD1	2.42	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:A:258:ILE:HG22	1:A:278:ALA:HB2	1.96	0.47
1:C:177:TYR:OH	1:C:210:GLY:HA3	2.14	0.47
1:D:172:GLN:HE22	1:D:201:HIS:HA	1.78	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
1:B:215:VAL:HA	1:B:219:VAL:HG23	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:134:PHE:CD2	1:A:154:MET:HG2	2.54	0.43
1:C:215:VAL:HA	1:C:219:VAL:CG2	2.48	0.43
1:C:186:TYR:O	1:C:190:LYS:N	2.51	0.43
1:A:268:LYS:HE2	1:A:268:LYS:HB3	1.89	0.42
1:B:186:TYR:O	1:B:190:LYS:N	2.50	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	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:LYS:HA	1:A:297:ASP:OD1	2.21	0.41
1:B:132:LEU:HD23	1:B:132:LEU:C	2.42	0.41
1:B:140:MET:HB3	1:B:203:LEU:O	2.21	0.41
1:A:215:VAL:HG11	1:A:234:LEU:HD13	2.03	0.40
1:D:174:SER:HA	1:D:205:LEU:O	2.22	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:151:LEU:HD23	1:C:154:MET:HE2	2.03	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

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%)	32	46
1	C	182/191 (95%)	171 (94%)	10 (6%)	1 (0%)	32	46
1	D	182/191 (95%)	171 (94%)	10 (6%)	1 (0%)	32	46
All	All	728/764 (95%)	683 (94%)	42 (6%)	3 (0%)	38	54

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%)	45	66
1	B	163/168 (97%)	157 (96%)	6 (4%)	39	59
1	C	163/168 (97%)	157 (96%)	6 (4%)	39	59
1	D	163/168 (97%)	159 (98%)	4 (2%)	53	73
All	All	652/672 (97%)	631 (97%)	21 (3%)	44	65

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 [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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	-	35,37,37	3.77	25 (71%)	44,51,51	2.50	14 (31%)
2	L08	B	1328	-	35,37,37	3.51	23 (65%)	44,51,51	2.60	16 (36%)
2	L08	C	2328	-	35,37,37	3.71	24 (68%)	44,51,51	2.46	15 (34%)
2	L08	D	3328	-	35,37,37	3.60	23 (65%)	44,51,51	2.60	15 (34%)

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	-	-	0/21/33/33	0/4/4/4
2	L08	B	1328	-	-	0/21/33/33	0/4/4/4
2	L08	C	2328	-	-	0/21/33/33	0/4/4/4
2	L08	D	3328	-	-	2/21/33/33	0/4/4/4

All (95) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	328	L08	C1-S1	-2.19	1.75	1.78
2	B	1328	L08	C22-C21	2.02	1.59	1.51
2	C	2328	L08	C22-C21	2.03	1.59	1.51
2	D	3328	L08	C23-N2	2.03	1.41	1.35
2	D	3328	L08	C22-C21	2.03	1.59	1.51
2	B	1328	L08	C18-N1	2.07	1.41	1.37
2	A	328	L08	C22-C21	2.11	1.59	1.51
2	B	1328	L08	C10-S1	2.12	1.81	1.78
2	C	2328	L08	C23-N2	2.14	1.42	1.35
2	D	3328	L08	C14-C15	2.28	1.42	1.38
2	A	328	L08	C23-N2	2.45	1.43	1.35
2	B	1328	L08	C20-N2	2.54	1.51	1.47
2	A	328	L08	C14-C15	2.59	1.43	1.38
2	D	3328	L08	C20-N2	2.66	1.51	1.47
2	A	328	L08	C20-N2	2.70	1.51	1.47
2	C	2328	L08	C20-N2	2.76	1.51	1.47
2	C	2328	L08	C8-C35	2.81	1.44	1.38
2	B	1328	L08	C14-C15	2.84	1.43	1.38
2	C	2328	L08	C14-C15	2.89	1.43	1.38
2	A	328	L08	C8-C35	2.94	1.44	1.38
2	C	2328	L08	C21-N2	2.97	1.52	1.47
2	D	3328	L08	C18-N1	3.01	1.43	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1328	L08	C17-C16	3.03	1.45	1.39
2	C	2328	L08	C10-S1	3.04	1.82	1.78
2	A	328	L08	C10-S1	3.04	1.82	1.78
2	B	1328	L08	C5-C6	3.07	1.44	1.38
2	D	3328	L08	C8-C35	3.07	1.45	1.38
2	B	1328	L08	C8-C35	3.09	1.45	1.38
2	C	2328	L08	C18-N1	3.11	1.44	1.37
2	D	3328	L08	C21-N2	3.16	1.52	1.47
2	D	3328	L08	C8-C16	3.22	1.46	1.39
2	D	3328	L08	C5-C6	3.25	1.45	1.38
2	A	328	L08	C17-C16	3.31	1.45	1.39
2	B	1328	L08	C21-N2	3.31	1.52	1.47
2	A	328	L08	C5-C6	3.37	1.45	1.38
2	C	2328	L08	C5-C6	3.38	1.45	1.38
2	C	2328	L08	C17-C16	3.48	1.45	1.39
2	A	328	L08	C18-N1	3.50	1.44	1.37
2	C	2328	L08	C8-C16	3.51	1.46	1.39
2	B	1328	L08	C11-C10	3.52	1.47	1.40
2	D	3328	L08	C17-C16	3.53	1.46	1.39
2	A	328	L08	C12-C13	3.55	1.46	1.39
2	A	328	L08	C21-N2	3.55	1.53	1.47
2	B	1328	L08	C8-C16	3.56	1.46	1.39
2	C	2328	L08	C12-C13	3.56	1.46	1.39
2	A	328	L08	C8-C16	3.57	1.46	1.39
2	B	1328	L08	C12-C13	3.63	1.46	1.39
2	B	1328	L08	C14-C13	3.74	1.47	1.39
2	D	3328	L08	C12-C13	3.87	1.46	1.39
2	A	328	L08	C14-C13	3.90	1.47	1.39
2	D	3328	L08	C11-C10	3.92	1.48	1.40
2	D	3328	L08	C14-C13	3.97	1.47	1.39
2	C	2328	L08	C14-C13	4.06	1.48	1.39
2	A	328	L08	C11-C10	4.23	1.48	1.40
2	D	3328	L08	O7-C2	4.43	1.46	1.37
2	A	328	L08	C6-C1	4.49	1.48	1.39
2	B	1328	L08	O7-C2	4.57	1.46	1.37
2	C	2328	L08	C11-C10	4.59	1.49	1.40
2	B	1328	L08	C6-C1	4.59	1.48	1.39
2	B	1328	L08	C15-C10	4.59	1.48	1.39
2	D	3328	L08	C5-C4	4.61	1.49	1.38
2	D	3328	L08	C3-C2	4.62	1.49	1.39
2	D	3328	L08	C6-C1	4.71	1.48	1.39
2	B	1328	L08	C5-C4	4.74	1.49	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1328	L08	C3-C2	4.75	1.49	1.39
2	D	3328	L08	C35-N20	4.82	1.45	1.34
2	C	2328	L08	C15-C10	4.83	1.48	1.39
2	A	328	L08	C5-C4	4.85	1.49	1.38
2	B	1328	L08	C35-N20	4.87	1.45	1.34
2	C	2328	L08	C5-C4	4.88	1.49	1.38
2	D	3328	L08	C15-C10	4.88	1.48	1.39
2	C	2328	L08	C6-C1	4.90	1.48	1.39
2	C	2328	L08	O7-C2	4.92	1.47	1.37
2	C	2328	L08	C35-N20	4.93	1.45	1.34
2	A	328	L08	C3-C2	4.98	1.50	1.39
2	A	328	L08	C35-N20	5.01	1.45	1.34
2	A	328	L08	O7-C2	5.03	1.47	1.37
2	C	2328	L08	C3-C2	5.10	1.50	1.39
2	A	328	L08	C15-C10	5.14	1.49	1.39
2	C	2328	L08	C4-C3	5.63	1.49	1.38
2	B	1328	L08	C4-C3	5.66	1.49	1.38
2	C	2328	L08	C2-C1	5.72	1.49	1.40
2	D	3328	L08	C4-C3	5.74	1.49	1.38
2	A	328	L08	C4-C3	5.81	1.50	1.38
2	B	1328	L08	C2-C1	6.02	1.49	1.40
2	D	3328	L08	C2-C1	6.07	1.49	1.40
2	B	1328	L08	C12-C11	6.16	1.51	1.39
2	A	328	L08	C2-C1	6.20	1.49	1.40
2	A	328	L08	C12-C11	6.36	1.51	1.39
2	C	2328	L08	C12-C11	6.65	1.52	1.39
2	D	3328	L08	C12-C11	6.73	1.52	1.39
2	B	1328	L08	C17-C18	7.40	1.51	1.39
2	D	3328	L08	C17-C18	7.86	1.52	1.39
2	C	2328	L08	C17-C18	8.30	1.52	1.39
2	A	328	L08	C17-C18	8.66	1.53	1.39

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3328	L08	C21-N2-C23	-5.82	107.34	122.96
2	D	3328	L08	C12-C11-C10	-5.66	117.28	122.93
2	C	2328	L08	C21-N2-C23	-5.52	108.13	122.96
2	A	328	L08	C21-N2-C23	-5.50	108.18	122.96
2	A	328	L08	C12-C11-C10	-5.42	117.52	122.93
2	B	1328	L08	C21-N2-C23	-5.35	108.58	122.96
2	B	1328	L08	C20-N2-C23	-5.08	109.31	122.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2328	L08	C12-C11-C10	-5.08	117.86	122.93
2	B	1328	L08	C12-C11-C10	-4.94	118.01	122.93
2	C	2328	L08	C20-N2-C23	-4.80	110.06	122.96
2	A	328	L08	C20-N2-C23	-4.79	110.10	122.96
2	D	3328	L08	C20-N2-C23	-4.52	110.81	122.96
2	B	1328	L08	C17-C18-N1	-4.01	117.42	122.15
2	D	3328	L08	C17-C18-N1	-3.92	117.52	122.15
2	A	328	L08	C17-C18-N1	-3.57	117.94	122.15
2	C	2328	L08	C17-C18-N1	-3.50	118.02	122.15
2	B	1328	L08	C19-N1-C18	-3.26	112.65	120.33
2	A	328	L08	C19-N1-C18	-2.95	113.38	120.33
2	C	2328	L08	C19-N1-C18	-2.82	113.69	120.33
2	A	328	L08	C21-C22-N1	-2.78	105.19	110.68
2	D	3328	L08	C19-N1-C18	-2.67	114.03	120.33
2	C	2328	L08	C21-C22-N1	-2.51	105.73	110.68
2	B	1328	L08	C21-C22-N1	-2.49	105.78	110.68
2	D	3328	L08	C21-C22-N1	-2.47	105.82	110.68
2	C	2328	L08	C17-C18-N20	-2.41	118.21	122.45
2	D	3328	L08	C17-C18-N20	-2.39	118.24	122.45
2	A	328	L08	C17-C18-N20	-2.32	118.35	122.45
2	B	1328	L08	C8-C35-N20	-2.27	121.30	123.92
2	B	1328	L08	C17-C16-C13	-2.23	117.11	120.88
2	D	3328	L08	C17-C16-C13	-2.18	117.18	120.88
2	C	2328	L08	C17-C16-C13	-2.15	117.24	120.88
2	B	1328	L08	C17-C18-N20	-2.12	118.71	122.45
2	A	328	L08	C17-C16-C13	-2.03	117.45	120.88
2	B	1328	L08	C13-C12-C11	2.01	122.42	119.56
2	D	3328	L08	C9-O7-C2	2.02	122.75	118.04
2	C	2328	L08	C21-N2-C20	2.07	116.51	112.61
2	B	1328	L08	C21-N2-C20	2.08	116.54	112.61
2	D	3328	L08	C13-C12-C11	2.12	122.57	119.56
2	C	2328	L08	C13-C12-C11	2.15	122.61	119.56
2	A	328	L08	C13-C12-C11	2.32	122.86	119.56
2	C	2328	L08	C6-C1-C2	2.77	122.04	118.61
2	D	3328	L08	C6-C1-C2	2.84	122.13	118.61
2	B	1328	L08	C6-C1-C2	3.06	122.40	118.61
2	A	328	L08	C6-C1-C2	3.24	122.63	118.61
2	C	2328	L08	C12-C11-N3	3.40	119.77	115.82
2	A	328	L08	C12-C11-N3	4.09	120.57	115.82
2	A	328	L08	N20-C18-N1	4.56	123.70	116.89
2	C	2328	L08	N20-C18-N1	4.60	123.77	116.89
2	D	3328	L08	C12-C11-N3	4.62	121.18	115.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1328	L08	N20-C18-N1	4.67	123.87	116.89
2	B	1328	L08	C12-C11-N3	4.67	121.24	115.82
2	A	328	L08	C35-N20-C18	4.80	123.41	116.88
2	C	2328	L08	C35-N20-C18	4.86	123.49	116.88
2	D	3328	L08	N20-C18-N1	4.92	124.24	116.89
2	B	1328	L08	C35-N20-C18	4.96	123.63	116.88
2	D	3328	L08	C35-N20-C18	5.09	123.80	116.88
2	A	328	L08	C22-N1-C19	6.75	125.88	111.57
2	D	3328	L08	C22-N1-C19	7.00	126.40	111.57
2	C	2328	L08	C22-N1-C19	7.10	126.61	111.57
2	B	1328	L08	C22-N1-C19	7.51	127.50	111.57

There are no chirality outliers.

All (2) torsion outliers are listed below:

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

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	328	L08	2	0
2	B	1328	L08	3	0
2	C	2328	L08	5	0
2	D	3328	L08	2	0

## 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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.