



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:46 PM GMT

PDB ID : 4AG4
Title : Crystal structure of a DDR1-Fab complex
Authors : Carafoli, F.; Mayer, M.C.; Shiraishi, K.; Pecheva, M.A.; Chan, L.Y.; Nan, R.;
Leitinger, B.; Hohenester, E.
Deposited on : 2012-01-24
Resolution : 2.80 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

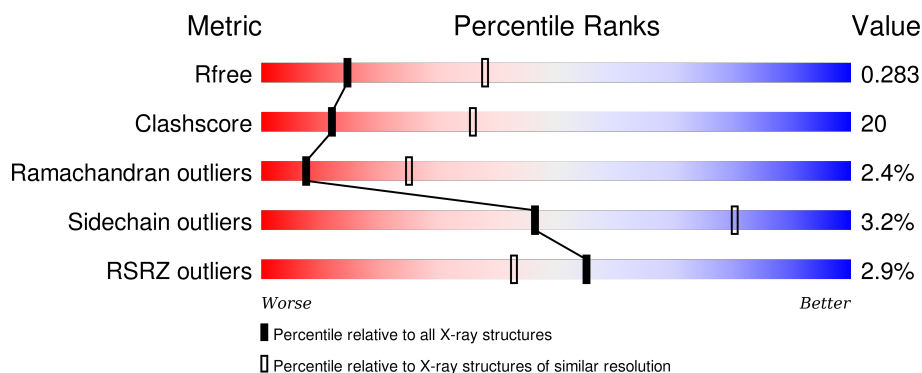
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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. 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	351	<div> <div>4%</div> <div>64%</div> <div>32%</div> <div>..</div> </div>
2	H	215	<div> <div>%</div> <div>63%</div> <div>31%</div> <div>..</div> </div>
3	L	213	<div> <div>2%</div> <div>53%</div> <div>44%</div> <div>..</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5885 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 EPITHELIAL DISCOIDIN DOMAIN-CONTAINING RECEPTOR 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	1
			2644	1661	481	485	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	ALA	-	EXPRESSION TAG	UNP Q08345
A	27	PRO	-	EXPRESSION TAG	UNP Q08345
A	28	LEU	-	EXPRESSION TAG	UNP Q08345
A	368	ALA	-	EXPRESSION TAG	UNP Q08345
A	369	ALA	-	EXPRESSION TAG	UNP Q08345
A	370	ALA	-	EXPRESSION TAG	UNP Q08345
A	371	HIS	-	EXPRESSION TAG	UNP Q08345
A	372	HIS	-	EXPRESSION TAG	UNP Q08345
A	373	HIS	-	EXPRESSION TAG	UNP Q08345
A	374	HIS	-	EXPRESSION TAG	UNP Q08345
A	375	HIS	-	EXPRESSION TAG	UNP Q08345
A	376	HIS	-	EXPRESSION TAG	UNP Q08345

- Molecule 2 is a protein called MONOCLONAL ANTIBODY 3E3 HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	209	Total	C	N	O	S	0	0	0
			1567	995	254	313	5			

- Molecule 3 is a protein called MONOCLONAL ANTIBODY 3E3 LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	1
			1602	1000	263	330	9			

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	0	0
			28	16	2	10		
4	A	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Ca	0	0
			2	2		

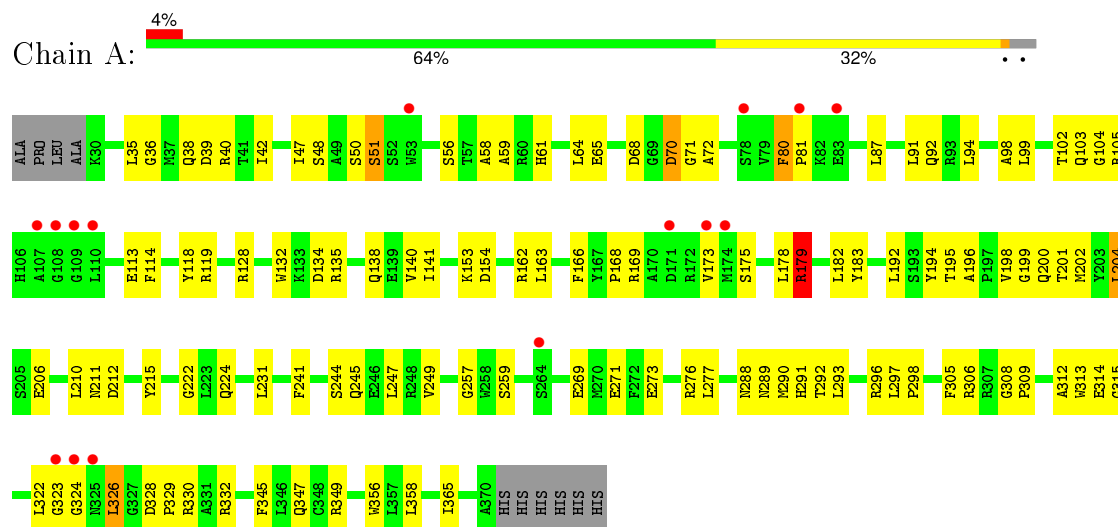
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	6	Total	O	0	0
			6	6		
6	H	5	Total	O	0	0
			5	5		
6	L	3	Total	O	0	0
			3	3		

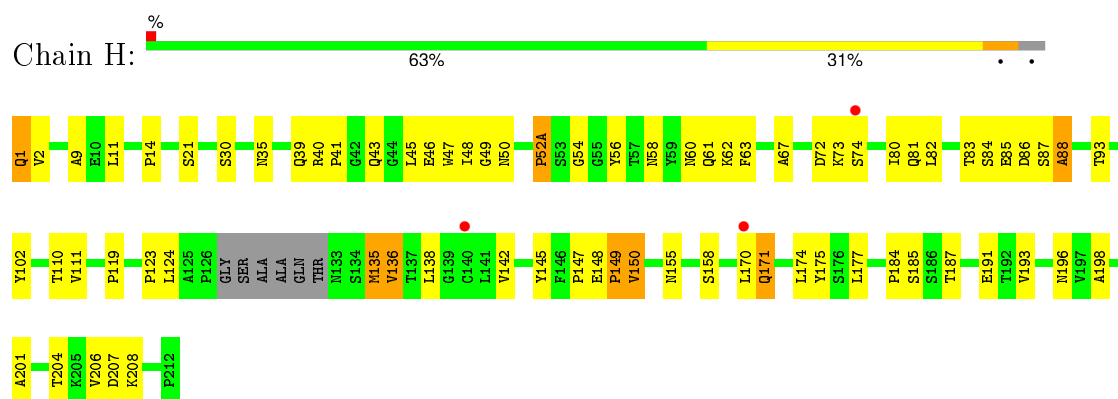
3 Residue-property plots

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.

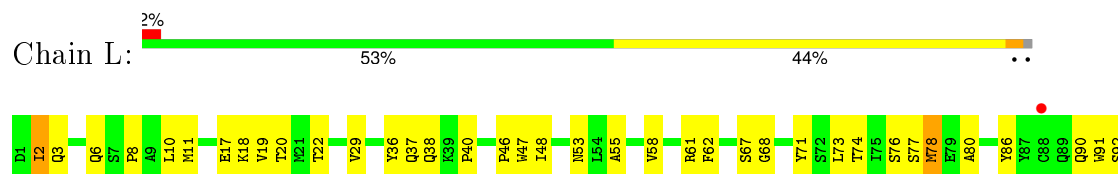
• Molecule 1: EPITHELIAL DISCOIDIN DOMAIN-CONTAINING RECEPTOR 1

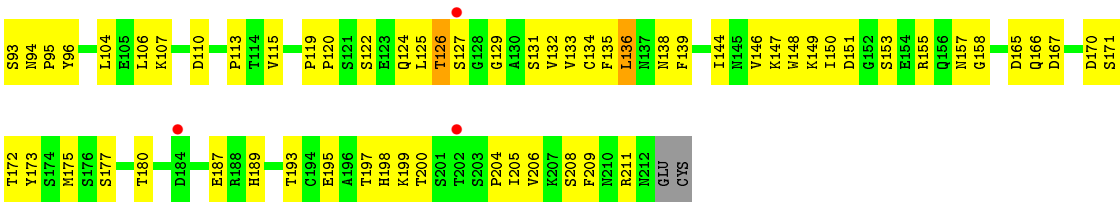


• Molecule 2: MONOCLONAL ANTIBODY 3E3 HEAVY CHAIN



• Molecule 3: MONOCLONAL ANTIBODY 3E3 LIGHT CHAIN





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	102.51Å 251.48Å 75.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 19.92 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.5 (20.00-2.80) 98.5 (19.92-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.20 (at 2.79Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.216 , 0.285 0.217 , 0.283	Depositor DCC
R_{free} test set	2388 reflections (11.03%)	DCC
Wilson B-factor (Å ²)	57.5	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 24034 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5885	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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: CA, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.41	0/2714	0.69	1/3679 (0.0%)
2	H	0.40	0/1611	0.68	0/2211
3	L	0.38	0/1643	0.62	0/2239
All	All	0.40	0/5968	0.67	1/8129 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	ARG	N-CA-C	-5.62	95.82	111.00

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	2644	0	2486	95	0
2	H	1567	0	1497	60	0
3	L	1602	0	1506	74	0
4	A	56	0	50	0	0
5	A	2	0	0	0	0
6	A	6	0	0	0	0
6	H	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	L	3	0	0	0	0
All	All	5885	0	5539	228	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 20.

All (228) 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:80:PHE:HB3	1:A:81:PRO:HD2	1.29	1.13
1:A:138:GLN:HE21	1:A:140:VAL:H	1.12	0.95
1:A:72:ALA:HB2	1:A:179:ARG:HD3	1.54	0.89
3:L:195:GLU:HG2	3:L:206:VAL:HG22	1.57	0.84
2:H:1:GLN:NE2	2:H:1:GLN:N	2.28	0.82
1:A:138:GLN:NE2	1:A:140:VAL:H	1.78	0.82
1:A:80:PHE:HB3	1:A:81:PRO:CD	2.08	0.81
1:A:199:GLY:HA3	1:A:211:ASN:HA	1.63	0.81
3:L:8:PRO:HD2	3:L:11:MET:HE2	1.64	0.80
3:L:8:PRO:HD2	3:L:11:MET:CE	2.12	0.79
3:L:155:ARG:HE	3:L:157:ASN:HB2	1.47	0.79
1:A:210:LEU:HD13	1:A:358:LEU:HD11	1.65	0.79
1:A:138:GLN:HE21	1:A:140:VAL:N	1.80	0.78
1:A:135:ARG:HG2	1:A:135:ARG:HH11	1.54	0.73
1:A:298:PRO:O	1:A:322:LEU:HD12	1.88	0.73
3:L:2:ILE:HD12	3:L:93:SER:HB3	1.72	0.72
1:A:141:ILE:CG2	1:A:153:LYS:HG3	2.20	0.72
1:A:290:MET:HG2	1:A:292:THR:HG22	1.71	0.71
1:A:80:PHE:CB	1:A:81:PRO:HD2	2.17	0.71
1:A:141:ILE:HG23	1:A:153:LYS:HG3	1.73	0.71
2:H:1:GLN:HE21	2:H:1:GLN:H1	1.38	0.70
2:H:1:GLN:NE2	2:H:1:GLN:H3	1.90	0.70
2:H:83:THR:O	2:H:85:GLU:N	2.23	0.69
3:L:78:MET:HE1	3:L:104:LEU:HD21	1.73	0.69
3:L:193:THR:HG23	3:L:208:SER:HB2	1.73	0.69
1:A:36:GLY:HA2	1:A:39:ASP:HB3	1.75	0.68
2:H:123:PRO:O	2:H:124:LEU:HD23	1.93	0.68
1:A:58:ALA:H	1:A:61:HIS:HD2	1.42	0.68
2:H:67:ALA:HA	2:H:81:GLN:O	1.93	0.67
1:A:50:SER:O	1:A:51:SER:HB3	1.94	0.67
1:A:70:ASP:O	1:A:105:ARG:HD2	1.95	0.67
2:H:40:ARG:HH21	2:H:85:GLU:HB3	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:90:GLN:HE21	3:L:92:SER:H	1.43	0.66
2:H:35:ASN:OD1	2:H:50:ASN:HB3	1.96	0.65
2:H:150:VAL:CG1	2:H:177:LEU:HD13	2.28	0.64
2:H:119:PRO:HB3	2:H:145:TYR:HB3	1.78	0.64
1:A:39:ASP:O	1:A:40:ARG:HB2	1.97	0.64
1:A:202:MET:HB3	1:A:356:TRP:CZ2	2.32	0.64
1:A:103:GLN:HB3	1:A:179:ARG:HB3	1.81	0.63
3:L:90:GLN:NE2	3:L:93:SER:H	1.97	0.63
1:A:113:GLU:HB3	1:A:175:SER:O	1.99	0.63
1:A:72:ALA:CB	1:A:179:ARG:HD3	2.29	0.62
1:A:326:LEU:HD21	1:A:332:ARG:HD2	1.81	0.62
1:A:292:THR:HG23	1:A:293:LEU:HG	1.82	0.61
3:L:20:THR:OG1	3:L:74:THR:HG23	2.00	0.61
2:H:47:TRP:CZ2	2:H:49:GLY:HA2	2.35	0.61
2:H:187:THR:O	2:H:191:GLU:HB2	2.00	0.61
3:L:78:MET:CE	3:L:104:LEU:HD21	2.31	0.60
2:H:83:THR:H	2:H:86:ASP:HB2	1.66	0.60
3:L:125:LEU:C	3:L:127:SER:H	2.05	0.60
2:H:1:GLN:NE2	2:H:1:GLN:H1	1.92	0.60
2:H:87:SER:HB3	2:H:111:VAL:HG23	1.84	0.59
2:H:50:ASN:OD1	2:H:58:ASN:HB2	2.03	0.59
2:H:147:PRO:HD2	2:H:201:ALA:CB	2.32	0.58
3:L:36:TYR:CD1	3:L:46:PRO:HA	2.38	0.58
3:L:187:GLU:HG2	3:L:211:ARG:NH2	2.19	0.58
2:H:11:LEU:HD12	2:H:110:THR:O	2.03	0.58
1:A:99:LEU:HB2	1:A:183:TYR:HB2	1.84	0.58
1:A:99:LEU:HD23	1:A:153:LYS:O	2.04	0.58
1:A:212:ASP:HB3	1:A:224:GLN:HE22	1.69	0.58
3:L:8:PRO:HD2	3:L:11:MET:HE1	1.87	0.57
1:A:271:GLU:HG3	1:A:347:GLN:HG3	1.86	0.57
1:A:56:SER:HB2	1:A:70:ASP:OD1	2.04	0.57
2:H:46:GLU:OE2	2:H:63:PHE:HE1	1.89	0.56
2:H:54:GLY:HA3	2:H:56:TYR:CE2	2.41	0.56
1:A:276:ARG:HG2	1:A:277:LEU:H	1.69	0.56
2:H:2:VAL:HB	2:H:102:TYR:CE1	2.41	0.55
1:A:94:LEU:HD11	1:A:192:LEU:CD2	2.37	0.55
1:A:92:GLN:HA	1:A:162:ARG:NH1	2.21	0.55
2:H:1:GLN:N	2:H:1:GLN:CD	2.59	0.55
1:A:365:ILE:HD12	1:A:365:ILE:N	2.22	0.55
1:A:198:VAL:HG23	1:A:222:GLY:C	2.28	0.55
3:L:136:LEU:HD11	3:L:146:VAL:HG22	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:72:ASP:C	2:H:74:SER:H	2.10	0.54
1:A:119:ARG:NH1	1:A:169:ARG:HD2	2.23	0.54
1:A:105:ARG:NE	1:A:113:GLU:OE1	2.36	0.54
3:L:120:PRO:HD3	3:L:132:VAL:HG22	1.89	0.54
1:A:204:LEU:O	1:A:206:GLU:N	2.40	0.54
3:L:29:VAL:HG11	3:L:90:GLN:HG3	1.89	0.54
3:L:90:GLN:NE2	3:L:93:SER:N	2.55	0.54
2:H:142:VAL:HB	2:H:177:LEU:HB3	1.89	0.54
1:A:215:TYR:CE2	1:A:224:GLN:HB3	2.43	0.53
1:A:273:GLU:HG3	1:A:345:PHE:CE2	2.42	0.53
1:A:210:LEU:HD22	1:A:358:LEU:CD1	2.39	0.53
3:L:167:ASP:HB3	3:L:170:ASP:OD1	2.08	0.53
2:H:1:GLN:H3	2:H:1:GLN:CD	2.12	0.53
1:A:71:GLY:HA3	1:A:105:ARG:HG3	1.90	0.53
2:H:150:VAL:HG12	2:H:177:LEU:HD13	1.90	0.53
2:H:14:PRO:HG3	2:H:111:VAL:HG12	1.89	0.53
1:A:166:PHE:O	1:A:168:PRO:HD3	2.09	0.53
3:L:166:GLN:HG3	3:L:173:TYR:CZ	2.44	0.53
1:A:61:HIS:O	1:A:72:ALA:HA	2.08	0.52
1:A:135:ARG:NH1	1:A:135:ARG:HG2	2.17	0.52
3:L:106:LEU:HD23	3:L:171:SER:OG	2.09	0.52
3:L:90:GLN:HE21	3:L:92:SER:N	2.07	0.52
3:L:19:VAL:HG12	3:L:20:THR:N	2.24	0.52
1:A:241:PHE:HB3	1:A:288:ASN:ND2	2.24	0.52
3:L:205:ILE:N	3:L:205:ILE:HD12	2.24	0.52
1:A:200:GLN:NE2	1:A:259:SER:H	2.08	0.52
3:L:148:TRP:HE1	3:L:177:SER:HB3	1.75	0.52
1:A:271:GLU:CG	1:A:347:GLN:HG3	2.41	0.51
1:A:257:GLY:HA3	1:A:358:LEU:HD13	1.91	0.51
3:L:10:LEU:HD12	3:L:11:MET:N	2.25	0.51
1:A:276:ARG:HG2	1:A:277:LEU:N	2.25	0.51
2:H:30:SER:O	2:H:52(A):PRO:HG2	2.11	0.51
1:A:194:TYR:CZ	1:A:231:LEU:HB2	2.46	0.50
2:H:40:ARG:HG2	2:H:88:ALA:HB2	1.93	0.50
1:A:308:GLY:HA2	1:A:345:PHE:CE1	2.47	0.50
3:L:205:ILE:H	3:L:205:ILE:HD12	1.76	0.50
2:H:204:THR:CG2	2:H:206:VAL:HG23	2.41	0.50
2:H:47:TRP:CH2	2:H:49:GLY:HA2	2.47	0.49
1:A:102:THR:HG21	1:A:118:TYR:CE2	2.46	0.49
3:L:62:PHE:HD1	3:L:73:LEU:HD21	1.76	0.49
1:A:36:GLY:HA3	1:A:42:ILE:HG13	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:125:LEU:O	3:L:127:SER:N	2.43	0.49
3:L:91:TRP:HA	3:L:96:TYR:CD1	2.47	0.49
2:H:170:LEU:HG	2:H:175:TYR:CE1	2.48	0.49
3:L:94:ASN:OD1	3:L:95:PRO:CA	2.60	0.49
2:H:155:ASN:O	2:H:158:SER:HB3	2.12	0.49
2:H:40:ARG:HG3	2:H:40:ARG:HH11	1.78	0.48
1:A:289:ASN:HB2	1:A:297:LEU:HD21	1.94	0.48
1:A:308:GLY:HA2	1:A:345:PHE:HE1	1.78	0.48
1:A:119:ARG:HH12	1:A:169:ARG:HD2	1.77	0.48
2:H:60:ASN:ND2	2:H:62:LYS:HB3	2.29	0.48
3:L:78:MET:HE1	3:L:104:LEU:CD2	2.41	0.48
3:L:144:ILE:HG23	3:L:175:MET:HE3	1.94	0.48
3:L:147:LYS:N	3:L:195:GLU:O	2.40	0.48
2:H:41:PRO:O	2:H:43:GLN:HG2	2.14	0.47
2:H:40:ARG:HG3	2:H:40:ARG:NH1	2.29	0.47
3:L:189:HIS:O	3:L:211:ARG:NH1	2.47	0.47
1:A:72:ALA:HB1	1:A:178:LEU:O	2.15	0.47
1:A:305:PHE:N	1:A:305:PHE:CD1	2.82	0.47
1:A:50:SER:O	1:A:51:SER:CB	2.62	0.47
1:A:210:LEU:CD1	1:A:358:LEU:HD11	2.40	0.47
3:L:150:ILE:O	3:L:151:ASP:HB2	2.14	0.47
3:L:6:GLN:HA	3:L:22:THR:O	2.14	0.47
2:H:135:MET:HE1	2:H:184:PRO:HG3	1.96	0.47
3:L:136:LEU:HD11	3:L:146:VAL:CG2	2.45	0.47
1:A:128:ARG:HH11	1:A:128:ARG:HG3	1.79	0.47
2:H:39:GLN:HB2	2:H:45:LEU:HD23	1.96	0.47
3:L:147:LYS:HB2	3:L:195:GLU:HB2	1.97	0.46
3:L:11:MET:HG3	3:L:104:LEU:HD12	1.97	0.46
2:H:35:ASN:HB2	2:H:93:THR:HG23	1.96	0.46
2:H:204:THR:HG22	2:H:206:VAL:HG23	1.96	0.46
1:A:313:TRP:O	1:A:315:GLY:N	2.49	0.46
1:A:289:ASN:O	1:A:329:PRO:HA	2.15	0.46
1:A:65:GLU:OE1	1:A:65:GLU:HA	2.15	0.46
3:L:158:GLY:O	3:L:180:THR:HG22	2.16	0.46
2:H:81:GLN:O	2:H:82:LEU:HD23	2.16	0.45
1:A:200:GLN:HE22	1:A:259:SER:H	1.64	0.45
3:L:138:ASN:HA	3:L:172:THR:OG1	2.16	0.45
2:H:123:PRO:HD3	2:H:208:LYS:HE2	1.98	0.45
3:L:67:SER:HA	3:L:71:TYR:CE1	2.51	0.45
3:L:80:ALA:HA	3:L:106:LEU:HD22	1.98	0.45
2:H:198:ALA:HA	2:H:204:THR:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:VAL:O	1:A:173:VAL:HG12	2.17	0.45
3:L:198:HIS:CE1	3:L:200:THR:HG23	2.51	0.45
3:L:125:LEU:HA	3:L:129:GLY:O	2.16	0.45
1:A:91:LEU:O	1:A:92:GLN:HB2	2.16	0.45
1:A:141:ILE:HG21	1:A:153:LYS:HG3	1.96	0.44
3:L:62:PHE:CD1	3:L:73:LEU:HD21	2.51	0.44
3:L:133:VAL:HG12	3:L:134:CYS:N	2.32	0.44
3:L:195:GLU:HG2	3:L:206:VAL:CG2	2.39	0.44
3:L:115:VAL:HA	3:L:135:PHE:O	2.17	0.44
1:A:322:LEU:O	1:A:324:GLY:N	2.50	0.44
1:A:244:SER:OG	1:A:245:GLN:N	2.50	0.44
1:A:39:ASP:O	1:A:40:ARG:CB	2.62	0.44
1:A:48:SER:O	1:A:87:LEU:HD12	2.18	0.44
3:L:106:LEU:HD23	3:L:171:SER:CB	2.47	0.44
2:H:123:PRO:C	2:H:124:LEU:HD23	2.38	0.44
2:H:80:ILE:O	2:H:80:ILE:HG23	2.18	0.44
3:L:119:PRO:HB3	3:L:209:PHE:CE2	2.52	0.44
3:L:94:ASN:OD1	3:L:95:PRO:HA	2.17	0.44
2:H:148:GLU:HG2	2:H:175:TYR:CE2	2.53	0.44
1:A:306:ARG:HG3	1:A:312:ALA:O	2.17	0.44
1:A:210:LEU:HD22	1:A:358:LEU:HD12	1.99	0.43
1:A:328:ASP:OD2	1:A:330:ARG:HG2	2.18	0.43
1:A:296:ARG:HG2	1:A:296:ARG:HH11	1.82	0.43
3:L:17:GLU:HG2	3:L:18:LYS:N	2.34	0.43
3:L:124:GLN:OE1	3:L:131:SER:N	2.48	0.43
3:L:149:LYS:HA	3:L:153:SER:O	2.18	0.43
1:A:58:ALA:H	1:A:61:HIS:CD2	2.29	0.43
2:H:46:GLU:OE2	2:H:63:PHE:CE1	2.71	0.43
2:H:150:VAL:HG11	2:H:177:LEU:HD13	1.99	0.43
3:L:107:LYS:HB2	3:L:107:LYS:HE3	1.76	0.43
3:L:55:ALA:O	3:L:58:VAL:HB	2.19	0.43
2:H:60:ASN:O	2:H:61:GLN:C	2.56	0.42
1:A:103:GLN:HG2	1:A:179:ARG:NH2	2.34	0.42
1:A:99:LEU:HD23	1:A:153:LYS:C	2.40	0.42
1:A:291:HIS:HB3	1:A:296:ARG:CG	2.49	0.42
2:H:83:THR:HG22	2:H:83:THR:O	2.18	0.42
2:H:2:VAL:HG11	2:H:102:TYR:CD2	2.54	0.42
3:L:73:LEU:HD23	3:L:73:LEU:C	2.39	0.42
3:L:48:ILE:HG23	3:L:53:ASN:O	2.20	0.42
2:H:60:ASN:OD1	3:L:95:PRO:HB3	2.19	0.42
1:A:195:THR:O	1:A:196:ALA:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:72:ASP:O	2:H:74:SER:N	2.50	0.42
1:A:296:ARG:HG2	1:A:296:ARG:NH1	2.35	0.42
2:H:170:LEU:O	2:H:171:GLN:HB2	2.20	0.42
1:A:80:PHE:O	1:A:81:PRO:C	2.58	0.42
1:A:132:TRP:CH2	1:A:134:ASP:HB3	2.54	0.42
1:A:47:ILE:HG22	1:A:59:ALA:HB1	2.01	0.42
2:H:48:ILE:HD12	2:H:63:PHE:CE2	2.55	0.42
3:L:106:LEU:HG	3:L:107:LYS:N	2.35	0.42
3:L:125:LEU:C	3:L:127:SER:N	2.72	0.41
1:A:104:GLY:HA3	1:A:114:PHE:HB3	2.02	0.41
1:A:98:ALA:O	1:A:99:LEU:HG	2.20	0.41
1:A:330:ARG:NE	1:A:332:ARG:NH1	2.68	0.41
3:L:10:LEU:HD12	3:L:11:MET:H	1.85	0.41
3:L:20:THR:HG1	3:L:74:THR:HG23	1.85	0.41
3:L:113:PRO:HB3	3:L:139:PHE:HB3	2.03	0.41
1:A:332:ARG:HG2	1:A:332:ARG:HH11	1.86	0.41
3:L:90:GLN:HE22	3:L:93:SER:H	1.65	0.41
2:H:170:LEU:HD23	2:H:174:LEU:O	2.21	0.41
1:A:247:LEU:O	1:A:249:VAL:HG13	2.20	0.41
3:L:86:TYR:N	3:L:86:TYR:CD1	2.89	0.41
2:H:138:LEU:HD12	2:H:193:VAL:HG11	2.02	0.41
2:H:9:ALA:HB2	2:H:149:PRO:HD2	2.01	0.41
1:A:35:LEU:HG	1:A:182:LEU:O	2.21	0.41
2:H:136:VAL:CG2	2:H:185:SER:HA	2.50	0.41
3:L:122:SER:O	3:L:126:THR:HG23	2.21	0.41
3:L:37:GLN:HG2	3:L:38:GLN:N	2.36	0.40
3:L:197:THR:CG2	3:L:204:PRO:HB3	2.50	0.40
1:A:38:GLN:HG3	1:A:64:LEU:O	2.21	0.40
3:L:61:ARG:HD2	3:L:77:SER:O	2.20	0.40
1:A:269:GLU:OE2	1:A:349:ARG:NH1	2.53	0.40
1:A:135:ARG:NH1	1:A:135:ARG:CG	2.84	0.40
3:L:67:SER:HA	3:L:71:TYR:CZ	2.56	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	339/351 (97%)	303 (89%)	29 (9%)	7 (2%)	9	29
2	H	205/215 (95%)	185 (90%)	15 (7%)	5 (2%)	7	25
3	L	209/213 (98%)	185 (88%)	18 (9%)	6 (3%)	6	19
All	All	753/779 (97%)	673 (89%)	62 (8%)	18 (2%)	7	25

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	314	GLU
1	A	323	GLY
2	H	84	SER
1	A	51	SER
2	H	171	GLN
3	L	68	GLY
1	A	80	PHE
1	A	309	PRO
3	L	126	THR
1	A	68	ASP
2	H	73	LYS
2	H	88	ALA
3	L	76	SER
1	A	326	LEU
3	L	199	LYS
3	L	40	PRO
2	H	52(A)	PRO
3	L	2	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	268/284 (94%)	262 (98%)	6 (2%)	60	89
2	H	176/184 (96%)	168 (96%)	8 (4%)	34	68
3	L	181/188 (96%)	175 (97%)	6 (3%)	45	79
All	All	625/656 (95%)	605 (97%)	20 (3%)	46	80

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ASP
1	A	154	ASP
1	A	163	LEU
1	A	179	ARG
1	A	201	THR
1	A	204	LEU
2	H	1	GLN
2	H	21	SER
2	H	135	MET
2	H	136	VAL
2	H	149	PRO
2	H	150	VAL
2	H	196	ASN
2	H	207	ASP
3	L	3	GLN
3	L	47	TRP
3	L	78	MET
3	L	110	ASP
3	L	136	LEU
3	L	165	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	HIS
1	A	138	GLN
1	A	144	ASN
1	A	200	GLN
1	A	224	GLN
1	A	281	GLN
2	H	1	GLN
2	H	3	GLN
2	H	5	GLN

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Mol	Chain	Res	Type
2	H	81	GLN
3	L	90	GLN
3	L	138	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

4 carbohydrates 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	$\# Z > 2$
4	NAG	A	4001	1,4	14,14,15	0.54	0	15,19,21	1.09	1 (6%)
4	NAG	A	4002	4	14,14,15	0.52	0	15,19,21	0.69	0
4	NAG	A	4003	1,4	14,14,15	0.52	0	15,19,21	0.64	0
4	NAG	A	4004	4	14,14,15	0.79	1 (7%)	15,19,21	0.72	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	4001	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	4002	4	-	2/6/23/26	0/1/1/1
4	NAG	A	4003	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	4004	4	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	4004	NAG	C1-C2	2.22	1.55	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	4001	NAG	C2-N2-C7	-3.62	118.39	123.04

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	4002	NAG	O7-C7-N2-C2
4	A	4002	NAG	C8-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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(Å ²)	Q<0.9
1	A	341/351 (97%)	-0.06	15 (4%) 38 26	30, 47, 75, 95	0
2	H	209/215 (97%)	-0.12	3 (1%) 78 69	30, 50, 70, 82	0
3	L	211/213 (99%)	0.06	4 (1%) 70 59	33, 61, 88, 92	0
All	All	761/779 (97%)	-0.05	22 (2%) 55 43	30, 51, 81, 95	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	171	ASP	6.4
1	A	110	LEU	4.0
1	A	109	GLY	3.7
1	A	174	MET	3.3
1	A	325	ASN	3.2
1	A	81	PRO	3.2
1	A	323	GLY	3.2
1	A	324	GLY	2.9
3	L	127	SER	2.8
1	A	108	GLY	2.7
3	L	202	THR	2.7
1	A	83	GLU	2.4
2	H	74	SER	2.3
2	H	140	CYS	2.3
1	A	264	SER	2.3
2	H	170	LEU	2.3
1	A	173	VAL	2.1
1	A	78	SER	2.1
1	A	53	TRP	2.1
3	L	184	ASP	2.0
1	A	107	ALA	2.0
3	L	88	CYS	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	NAG	A	4003	14/15	0.85	0.22	1.46	60,63,65,67	0
4	NAG	A	4001	14/15	0.94	0.15	-	55,60,64,66	0
4	NAG	A	4004	14/15	0.88	0.43	-	70,72,75,75	0
4	NAG	A	4002	14/15	0.92	0.26	-	64,68,70,70	0

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	CA	A	5002	1/1	0.96	0.17	-0.70	69,69,69,69	0
5	CA	A	5001	1/1	0.97	0.09	-1.92	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.