



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:09 AM UTC

PDB ID : 4LI1 / pdb_00004li1
Title : Crystal Structures of Lgr4 and its complex with R-spondin1
Authors : Xu, Y.; Rajashankar, K.; Robev, D.
Deposited on : 2013-07-02
Resolution : 2.66 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

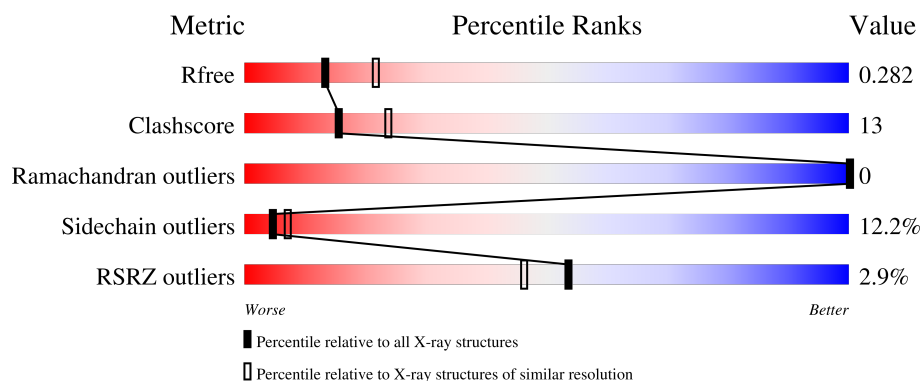
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.66 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	432	5% 55% 37% 6% ..
1	B	432	69% 25% ..

2 Entry composition [i](#)

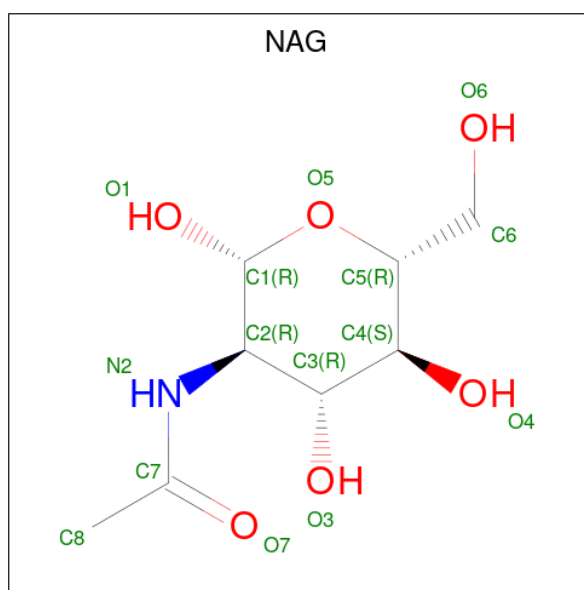
There are 3 unique types of molecules in this entry. The entry contains 6803 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Leucine-rich repeat-containing G-protein coupled receptor 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	425	Total	C	N	O	S	0	0	0
			3303	2102	565	625	11			
1	A	424	Total	C	N	O	S	0	0	0
			3298	2099	564	624	11			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

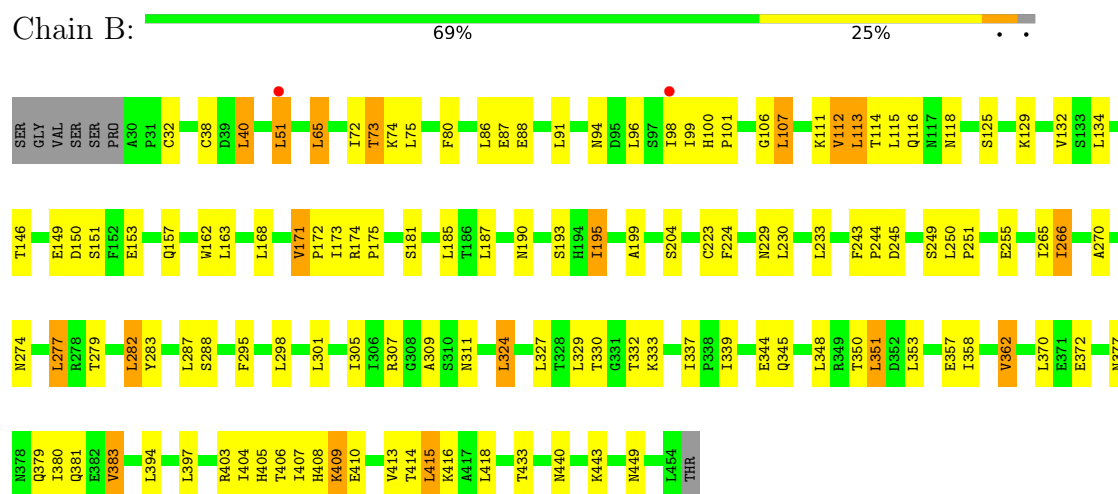
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	114	Total 114	O 114	0	0
3	A	60	Total 60	O 60	0	0

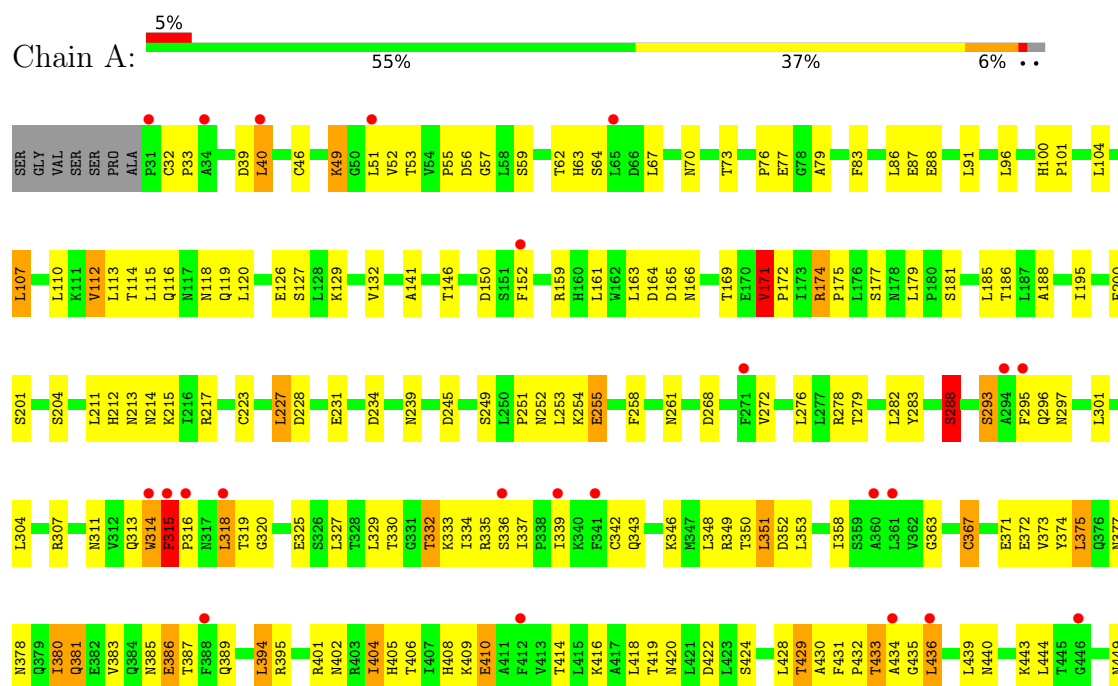
3 Residue-property plots

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

- Molecule 1: Leucine-rich repeat-containing G-protein coupled receptor 4



- Molecule 1: Leucine-rich repeat-containing G-protein coupled receptor 4



VAL
KAS
ETD
TRD
LAS
THR

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	60.29Å 158.58Å 227.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	79.29 – 2.66 79.29 – 2.66	Depositor EDS
% Data completeness (in resolution range)	99.7 (79.29-2.66) 93.9 (79.29-2.66)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.216 , 0.272 0.229 , 0.282	Depositor DCC
R_{free} test set	1610 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	59.6	Xtriage
Anisotropy	0.413	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6803	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.11% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.58	0/3365	1.08	15/4578 (0.3%)
1	B	0.63	1/3370 (0.0%)	1.07	6/4586 (0.1%)
All	All	0.61	1/6735 (0.0%)	1.08	21/9164 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	266	ILE	CA-CB	5.67	1.59	1.53

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	314	TRP	N-CA-C	9.01	121.18	108.74
1	B	195	ILE	CA-C-N	8.21	128.35	120.31
1	B	195	ILE	C-N-CA	8.21	128.35	120.31
1	A	171	VAL	CA-C-N	7.00	128.59	119.84
1	A	171	VAL	C-N-CA	7.00	128.59	119.84
1	A	433	THR	N-CA-C	-6.87	98.63	109.07
1	A	436	LEU	N-CA-C	-6.72	104.70	113.17
1	A	83	PHE	CA-C-N	6.55	126.14	119.19
1	A	83	PHE	C-N-CA	6.55	126.14	119.19
1	A	268	ASP	N-CA-C	-6.43	100.76	110.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	381	GLN	N-CA-C	5.79	120.49	112.45
1	A	336	SER	N-CA-C	5.78	116.03	107.88
1	A	288	SER	N-CA-C	-5.68	106.19	113.23
1	B	409	LYS	N-CA-C	-5.67	103.05	110.53
1	A	315	PHE	CA-C-N	5.67	126.14	120.14
1	A	315	PHE	C-N-CA	5.67	126.14	120.14
1	B	305	ILE	N-CA-C	5.52	115.81	107.75
1	A	239	ASN	N-CA-C	5.45	118.93	112.72
1	A	319	THR	N-CA-C	-5.24	101.61	109.95
1	B	380	ILE	N-CA-C	5.18	117.10	109.63
1	B	344	GLU	N-CA-C	5.15	119.62	113.23

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	106	GLY	Peptide

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	3298	0	3317	110	0
1	B	3303	0	3319	69	0
2	B	28	0	26	1	0
3	A	60	0	0	5	0
3	B	114	0	0	4	0
All	All	6803	0	6662	178	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 13.

All (178) 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:B:408:HIS:ND1	1:B:409:LYS:O	2.05	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:ILE:HG22	1:B:223:CYS:HB2	1.64	0.78
1:A:255:GLU:HG3	1:A:279:THR:HB	1.65	0.78
1:A:119:GLN:O	3:A:524:HOH:O	2.06	0.73
1:B:255:GLU:HG3	1:B:279:THR:HB	1.70	0.72
1:B:348:LEU:HD21	1:B:351:LEU:HG	1.74	0.68
1:B:174:ARG:HG2	1:B:175:PRO:HD3	1.75	0.68
1:A:350:THR:HG22	1:A:372:GLU:HB2	1.76	0.66
1:B:204:SER:O	1:B:229:ASN:ND2	2.29	0.66
1:B:91:LEU:O	1:B:94:ASN:ND2	2.29	0.65
1:A:88:GLU:HG2	1:A:112:VAL:HG13	1.79	0.65
1:B:171:VAL:HG22	1:B:173:ILE:HD13	1.77	0.64
1:A:419:THR:OG1	1:A:420:ASN:OD1	2.12	0.64
1:A:283:TYR:HB2	1:A:307:ARG:HB2	1.80	0.64
1:B:383:VAL:HG13	1:B:407:ILE:HG12	1.79	0.63
1:A:315:PHE:CG	1:A:316:PRO:HD2	2.33	0.63
1:A:434:ALA:HB1	1:A:435:GLY:HA2	1.79	0.62
1:A:258:PHE:O	1:A:261:ASN:ND2	2.31	0.62
1:B:350:THR:HG22	1:B:372:GLU:HB2	1.81	0.62
1:A:293:SER:HB3	1:A:296:GLN:HE21	1.66	0.61
1:B:274:ASN:HB2	1:B:277:LEU:HD22	1.82	0.61
1:B:283:TYR:HB2	1:B:307:ARG:HB2	1.83	0.60
1:A:348:LEU:HD21	1:A:351:LEU:HG	1.83	0.60
1:B:73:THR:O	1:B:96:LEU:HA	2.02	0.60
1:A:76:PRO:HG2	1:A:79:ALA:HB2	1.82	0.60
1:A:159:ARG:NH2	1:A:181:SER:O	2.32	0.60
1:B:440:ASN:O	1:A:440:ASN:ND2	2.35	0.59
1:A:371:GLU:HA	1:A:394:LEU:HA	1.84	0.59
1:B:287:LEU:HB2	1:B:309:ALA:HB1	1.83	0.59
1:A:381:GLN:NE2	3:A:526:HOH:O	2.28	0.59
1:B:414:THR:O	1:B:416:LYS:NZ	2.36	0.58
1:A:231:GLU:HB3	1:A:254:LYS:HG2	1.83	0.58
1:A:386:GLU:HA	1:A:389:GLN:HB2	1.84	0.58
1:A:316:PRO:HB2	1:A:318:LEU:HD23	1.85	0.58
1:A:215:LYS:HE3	1:A:217:ARG:HH12	1.66	0.58
1:A:174:ARG:O	1:A:177:SER:OG	2.22	0.58
1:A:342:CYS:HB3	1:A:367:CYS:N	2.19	0.57
1:B:415:LEU:HD23	1:B:418:LEU:HD22	1.86	0.56
1:A:186:THR:HG22	1:A:188:ALA:H	1.69	0.56
1:A:432:PRO:O	3:A:540:HOH:O	2.17	0.56
1:A:46:CYS:SG	1:A:55:PRO:HG2	2.45	0.56
1:A:228:ASP:O	1:A:252:ASN:ND2	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:385:ASN:OD1	1:A:410:GLU:HG3	2.06	0.56
1:B:125:SER:HB2	1:B:151:SER:HA	1.88	0.55
1:A:212:HIS:HB3	1:A:234:ASP:OD1	2.06	0.55
1:A:181:SER:HB2	3:A:544:HOH:O	2.06	0.54
1:B:174:ARG:H	1:B:174:ARG:HD3	1.72	0.54
1:A:314:TRP:HA	1:A:334:ILE:HG13	1.88	0.54
1:A:385:ASN:HD21	1:A:408:HIS:CE1	2.25	0.54
1:B:80:PHE:HB3	1:B:107:LEU:HD21	1.89	0.54
1:B:333:LYS:HA	1:B:357:GLU:HG2	1.90	0.53
1:A:164:ASP:O	1:A:166:ASN:ND2	2.41	0.53
1:A:450:PHE:O	1:A:453:THR:OG1	2.25	0.53
1:B:96:LEU:HB2	1:B:118:ASN:OD1	2.08	0.53
1:A:174:ARG:HB2	1:A:175:PRO:HD3	1.90	0.53
1:A:254:LYS:HE3	1:A:278:ARG:HG3	1.91	0.53
1:B:88:GLU:HG2	1:B:112:VAL:HG13	1.89	0.52
1:A:49:LYS:H	1:A:70:ASN:HD21	1.55	0.52
1:B:65:LEU:HB2	1:B:86:LEU:HD11	1.91	0.52
1:A:380:ILE:HG13	1:A:404:ILE:HG22	1.92	0.52
1:B:381:GLN:O	1:B:404:ILE:HA	2.10	0.52
1:A:282:LEU:HD21	1:A:295:PHE:HZ	1.75	0.52
1:A:293:SER:HB3	1:A:296:GLN:NE2	2.24	0.51
1:B:224:PHE:HB3	1:B:250:LEU:HD21	1.92	0.51
1:A:307:ARG:HG2	1:A:330:THR:OG1	2.10	0.51
1:A:381:GLN:O	1:A:404:ILE:HA	2.11	0.51
1:A:297:ASN:H	1:A:320:GLY:H	1.58	0.51
1:A:114:THR:HB	1:A:116:GLN:OE1	2.11	0.50
1:A:377:ASN:OD1	1:A:401:ARG:NE	2.30	0.50
1:A:405:HIS:CE1	1:A:406:THR:HG22	2.46	0.50
1:B:403:ARG:NH1	3:B:648:HOH:O	2.44	0.50
1:A:51:LEU:HD21	1:A:55:PRO:HG3	1.94	0.50
1:A:385:ASN:O	3:A:557:HOH:O	2.20	0.50
1:B:87:GLU:OE2	1:B:111:LYS:NZ	2.44	0.50
1:A:195:ILE:HG22	1:A:223:CYS:HB2	1.94	0.50
1:B:40:LEU:H	1:B:40:LEU:HD12	1.76	0.49
1:B:187:LEU:O	1:B:190:ASN:ND2	2.37	0.49
1:B:114:THR:HB	1:B:116:GLN:OE1	2.12	0.49
1:B:287:LEU:HB2	1:B:309:ALA:CB	2.43	0.49
1:A:405:HIS:ND1	1:A:406:THR:HG22	2.28	0.49
1:A:126:GLU:O	1:A:129:LYS:HB2	2.12	0.49
1:A:49:LYS:H	1:A:70:ASN:ND2	2.11	0.48
1:A:297:ASN:H	1:A:320:GLY:N	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:ILE:HG23	1:A:402:ASN:OD1	2.13	0.48
1:A:126:GLU:HA	1:A:129:LYS:HD3	1.95	0.48
1:B:243:PHE:CZ	1:B:270:ALA:HB1	2.48	0.48
1:A:46:CYS:HB2	1:A:67:LEU:HD23	1.95	0.47
1:A:429:THR:O	1:A:449:ASN:ND2	2.47	0.47
1:A:372:GLU:HB3	1:A:374:TYR:CE2	2.50	0.47
1:A:428:LEU:HD13	1:A:430:ALA:O	2.15	0.47
1:A:371:GLU:HB3	1:A:395:ARG:HE	1.80	0.46
1:A:373:VAL:HG12	1:A:375:LEU:HD22	1.98	0.46
1:A:386:GLU:H	1:A:386:GLU:CD	2.23	0.46
1:A:87:GLU:C	1:A:110:LEU:HD12	2.41	0.46
1:A:293:SER:HA	1:A:296:GLN:HG3	1.97	0.46
1:A:408:HIS:HD2	1:A:409:LYS:H	1.62	0.46
1:B:51:LEU:O	1:B:72:ILE:HG13	2.16	0.46
1:A:51:LEU:CD2	1:A:55:PRO:HG3	2.46	0.46
1:B:32:CYS:SG	1:B:38:CYS:N	2.89	0.46
1:B:91:LEU:HD23	1:B:91:LEU:HA	1.77	0.46
1:A:307:ARG:NH2	1:A:352:ASP:OD2	2.48	0.46
1:A:152:PHE:HB3	1:A:179:LEU:HD21	1.97	0.46
1:A:363:GLY:HA2	1:A:387:THR:HA	1.98	0.46
1:B:113:LEU:HB2	1:B:134:LEU:HD11	1.97	0.45
1:B:171:VAL:HA	1:B:172:PRO:HD3	1.82	0.45
1:B:149:GLU:CD	1:B:174:ARG:HG3	2.40	0.45
1:B:249:SER:O	1:B:251:PRO:HD3	2.16	0.45
1:A:39:ASP:HB2	1:A:40:LEU:HD12	1.97	0.45
1:A:49:LYS:HA	1:A:49:LYS:HD3	1.70	0.45
1:A:171:VAL:HA	1:A:172:PRO:HD3	1.57	0.45
1:A:211:LEU:O	1:A:214:ASN:ND2	2.49	0.45
1:A:343:GLN:O	1:A:346:LYS:HE2	2.17	0.45
1:B:174:ARG:H	1:B:174:ARG:CD	2.29	0.45
1:A:342:CYS:HB3	1:A:367:CYS:H	1.81	0.45
1:A:163:LEU:HD23	1:A:163:LEU:HA	1.83	0.45
1:A:325:GLU:HG2	1:A:349:ARG:HD3	1.99	0.45
1:B:337:ILE:HD11	1:B:358:ILE:HD13	1.98	0.44
1:A:63:HIS:O	1:A:86:LEU:HD12	2.17	0.44
1:A:358:ILE:HB	1:A:378:ASN:ND2	2.31	0.44
1:A:107:LEU:HB3	1:A:110:LEU:HB2	2.00	0.44
1:A:96:LEU:HB2	1:A:118:ASN:OD1	2.18	0.44
1:A:335:ARG:HA	1:A:358:ILE:HA	1.98	0.44
1:A:451:LYS:HA	1:A:454:LEU:HD22	2.00	0.44
1:A:431:PHE:CE1	1:A:444:LEU:HB3	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:GLU:HG2	1:A:100:HIS:CE1	2.53	0.44
1:B:149:GLU:HG2	1:B:174:ARG:CZ	2.48	0.43
1:B:243:PHE:HA	1:B:244:PRO:HD3	1.75	0.43
1:B:288:SER:HA	1:B:311:ASN:HB2	1.99	0.43
1:A:200:PHE:HB3	1:A:227:LEU:HD21	1.99	0.43
1:A:422:ASP:OD1	1:A:424:SER:OG	2.26	0.43
1:B:162:TRP:CD1	1:B:162:TRP:N	2.86	0.43
1:B:324:LEU:HB2	1:B:345:GLN:OE1	2.19	0.43
1:B:80:PHE:CB	1:B:107:LEU:HD21	2.49	0.43
1:B:91:LEU:HB2	1:B:115:LEU:HD23	1.99	0.43
1:B:339:ILE:HA	1:B:362:VAL:HG11	2.00	0.43
1:A:449:ASN:C	1:A:449:ASN:HD22	2.27	0.43
1:B:307:ARG:HG2	1:B:330:THR:OG1	2.18	0.42
1:A:91:LEU:HB2	1:A:115:LEU:HD23	2.01	0.42
1:B:100:HIS:HA	1:B:101:PRO:HD2	1.88	0.42
1:A:315:PHE:CD1	1:A:316:PRO:HD2	2.54	0.42
1:A:428:LEU:HD11	1:A:431:PHE:CD1	2.54	0.42
1:A:59:SER:O	1:A:62:THR:HG23	2.19	0.42
1:A:249:SER:O	1:A:251:PRO:HD3	2.20	0.42
1:A:353:LEU:HA	1:A:353:LEU:HD23	1.77	0.42
1:B:72:ILE:O	1:B:94:ASN:HB3	2.19	0.42
1:B:74:LYS:HE3	1:B:98:ILE:H	1.85	0.42
1:B:150:ASP:O	1:B:153:GLU:HB2	2.19	0.42
1:A:51:LEU:HG	1:A:53:THR:O	2.19	0.42
1:A:301:LEU:HD21	1:A:304:LEU:HD22	2.00	0.42
1:B:73:THR:O	1:B:96:LEU:HD23	2.20	0.42
1:A:315:PHE:CZ	1:A:337:ILE:HG22	2.55	0.42
1:B:173:ILE:HD12	1:B:199:ALA:HA	2.02	0.42
1:B:409:LYS:O	1:B:410:GLU:HB2	2.20	0.42
1:A:351:LEU:HD22	1:A:353:LEU:HG	2.02	0.42
1:B:351:LEU:HD22	1:B:353:LEU:HG	2.01	0.41
1:A:101:PRO:O	1:A:127:SER:HA	2.20	0.41
1:B:282:LEU:N	1:B:282:LEU:HD12	2.35	0.41
1:B:370:LEU:HD23	1:B:370:LEU:HA	1.94	0.41
1:A:313:GLN:HA	1:A:333:LYS:HB3	2.01	0.41
1:A:254:LYS:NZ	1:A:276:LEU:O	2.44	0.41
1:A:418:LEU:O	1:A:439:LEU:HD22	2.20	0.41
1:A:288:SER:HA	1:A:311:ASN:HB2	2.01	0.41
1:A:325:GLU:O	1:A:348:LEU:HD12	2.21	0.41
1:A:332:THR:OG1	1:A:334:ILE:HD12	2.19	0.41
1:B:149:GLU:OE1	1:B:174:ARG:NE	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:LEU:HD21	1:B:295:PHE:HZ	1.85	0.41
1:B:405:HIS:HD2	3:B:709:HOH:O	2.04	0.41
1:B:408:HIS:CE1	2:B:502:NAG:H82	2.55	0.41
1:A:141:ALA:H	1:A:165:ASP:HB3	1.86	0.41
1:A:150:ASP:OD1	1:A:150:ASP:N	2.52	0.41
1:B:397:LEU:HD12	1:B:397:LEU:HA	1.86	0.40
1:B:408:HIS:HD2	3:B:710:HOH:O	2.04	0.40
1:A:380:ILE:HG13	1:A:380:ILE:O	2.20	0.40
1:B:163:LEU:HD23	1:B:163:LEU:HA	1.90	0.40
1:B:174:ARG:HB2	3:B:665:HOH:O	2.21	0.40
1:A:33:PRO:HD3	1:A:57:GLY:O	2.21	0.40
1:A:245:ASP:N	1:A:245:ASP:OD1	2.53	0.40
1:A:161:LEU:HB3	1:A:185:LEU:HD23	2.04	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	422/432 (98%)	409 (97%)	13 (3%)	0	100	100
1	B	423/432 (98%)	410 (97%)	13 (3%)	0	100	100
All	All	845/864 (98%)	819 (97%)	26 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	378/384 (98%)	329 (87%)	49 (13%)	4	6
1	B	378/384 (98%)	335 (89%)	43 (11%)	5	8
All	All	756/768 (98%)	664 (88%)	92 (12%)	5	7

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	40	LEU
1	B	51	LEU
1	B	65	LEU
1	B	73	THR
1	B	75	LEU
1	B	99	ILE
1	B	107	LEU
1	B	112	VAL
1	B	113	LEU
1	B	129	LYS
1	B	132	VAL
1	B	146	THR
1	B	157	GLN
1	B	168	LEU
1	B	171	VAL
1	B	181	SER
1	B	185	LEU
1	B	193	SER
1	B	230	LEU
1	B	233	LEU
1	B	245	ASP
1	B	265	ILE
1	B	266	ILE
1	B	277	LEU
1	B	282	LEU
1	B	298	LEU
1	B	301	LEU
1	B	324	LEU
1	B	327	LEU
1	B	329	LEU
1	B	332	THR
1	B	351	LEU
1	B	362	VAL

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Mol	Chain	Res	Type
1	B	377	ASN
1	B	379	GLN
1	B	383	VAL
1	B	394	LEU
1	B	406	THR
1	B	413	VAL
1	B	415	LEU
1	B	433	THR
1	B	443	LYS
1	B	449	ASN
1	A	32	CYS
1	A	40	LEU
1	A	49	LYS
1	A	52	VAL
1	A	56	ASP
1	A	64	SER
1	A	73	THR
1	A	104	LEU
1	A	107	LEU
1	A	112	VAL
1	A	113	LEU
1	A	120	LEU
1	A	132	VAL
1	A	146	THR
1	A	169	THR
1	A	171	VAL
1	A	174	ARG
1	A	201	SER
1	A	204	SER
1	A	213	ASN
1	A	227	LEU
1	A	253	LEU
1	A	255	GLU
1	A	272	VAL
1	A	288	SER
1	A	293	SER
1	A	315	PHE
1	A	318	LEU
1	A	327	LEU
1	A	329	LEU
1	A	332	THR
1	A	339	ILE

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Mol	Chain	Res	Type
1	A	351	LEU
1	A	367	CYS
1	A	375	LEU
1	A	380	ILE
1	A	383	VAL
1	A	386	GLU
1	A	394	LEU
1	A	404	ILE
1	A	410	GLU
1	A	414	THR
1	A	416	LYS
1	A	429	THR
1	A	433	THR
1	A	436	LEU
1	A	443	LYS
1	A	452	GLU
1	A	454	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	119	GLN
1	B	143	HIS
1	B	213	ASN
1	B	274	ASN
1	B	384	GLN
1	B	449	ASN
1	A	143	HIS
1	A	222	HIS
1	A	239	ASN
1	A	296	GLN
1	A	313	GLN
1	A	408	HIS
1	A	440	ASN
1	A	449	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	NAG	B	501	1	14,14,15	0.24	0	17,19,21	0.64	0
2	NAG	B	502	1	14,14,15	0.58	0	17,19,21	0.57	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
2	NAG	B	501	1	-	2/6/23/26	0/1/1/1
2	NAG	B	502	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	NAG	C8-C7-N2-C2
2	B	501	NAG	O7-C7-N2-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	NAG	1	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

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	424/432 (98%)	0.46	23 (5%) 31 25	36, 70, 104, 112	0
1	B	425/432 (98%)	-0.09	2 (0%) 87 85	32, 52, 85, 106	0
All	All	849/864 (98%)	0.19	25 (2%) 53 46	32, 60, 100, 112	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	412	PHE	3.4
1	A	341	PHE	3.3
1	A	316	PRO	3.2
1	A	315	PHE	2.9
1	A	295	PHE	2.8
1	A	314	TRP	2.8
1	A	34	ALA	2.8
1	A	388	PHE	2.7
1	A	360	ALA	2.6
1	A	152	PHE	2.5
1	A	318	LEU	2.5
1	A	31	PRO	2.5
1	A	436	LEU	2.4
1	A	446	GLY	2.4
1	A	65	LEU	2.3
1	A	294	ALA	2.3
1	B	98	ILE	2.3
1	A	336	SER	2.3
1	A	361	LEU	2.3
1	A	434	ALA	2.2
1	A	40	LEU	2.2
1	B	51	LEU	2.1
1	A	271	PHE	2.1
1	A	51	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	339	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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. 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	B-factors(Å ²)	Q<0.9
2	NAG	B	501	14/15	0.70	0.11	60,73,77,81	0
2	NAG	B	502	14/15	0.88	0.08	62,68,72,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.