



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 05:13 AM GMT

PDB ID : 3D6D
Title : Crystal Structure of the complex between PPARgamma LBD and the
LT175(R-enantiomer)
Authors : Pochetti, G.; Montanari, R.
Deposited on : 2008-05-19
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>

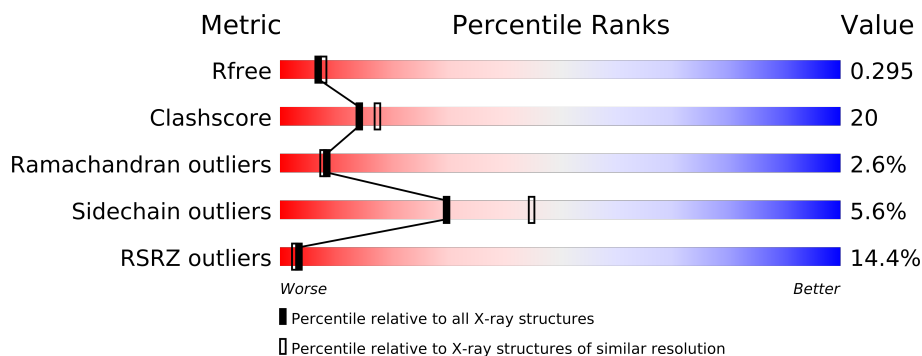
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
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))
R_{free}	66092	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (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.

Mol	Chain	Length	Quality of chain
1	A	286	
1	B	286	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	LRG	A	1	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4504 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.

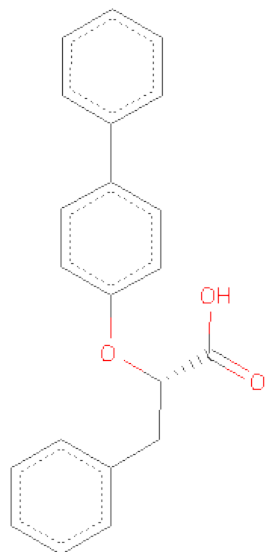
- Molecule 1 is a protein called Peroxisome proliferator-activatedreceptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	119	2	0
			2180	1406	357	407	10			
1	B	270	Total	C	N	O	S	127	0	0
			2166	1397	354	405	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	191	GLY	-	EXPRESSION TAG	UNP P37231
A	192	SER	-	EXPRESSION TAG	UNP P37231
A	193	HIS	-	EXPRESSION TAG	UNP P37231
A	194	MET	-	EXPRESSION TAG	UNP P37231
B	191	GLY	-	EXPRESSION TAG	UNP P37231
B	192	SER	-	EXPRESSION TAG	UNP P37231
B	193	HIS	-	EXPRESSION TAG	UNP P37231
B	194	MET	-	EXPRESSION TAG	UNP P37231

- Molecule 2 is (2S)-2-(BIPHENYL-4-YLOXY)-3-PHENYLPROPANOICACID (three-letter code: LRG) (formula: C₂₁H₁₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			24	21	3		

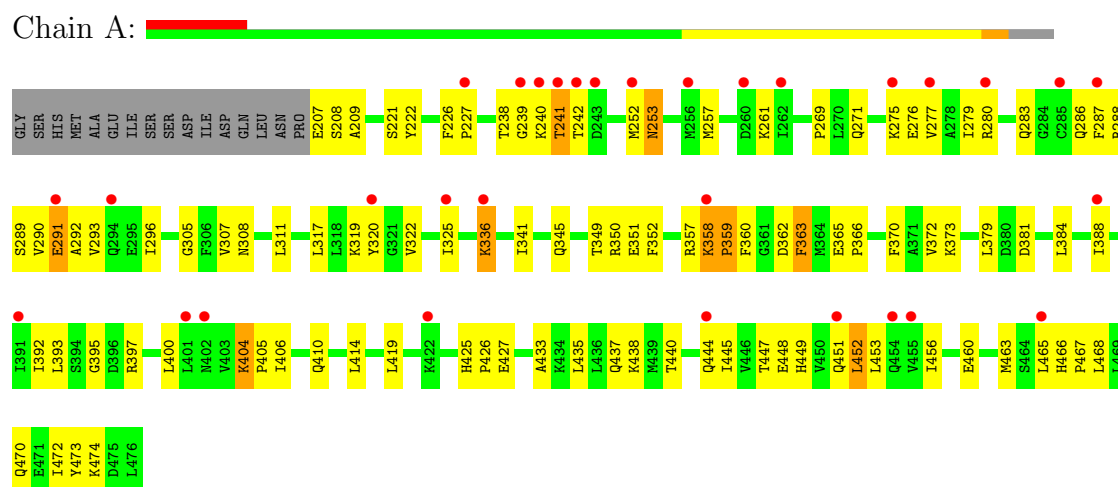
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	68	Total	O	0	0
			68	68		
3	B	66	Total	O	0	0
			66	66		

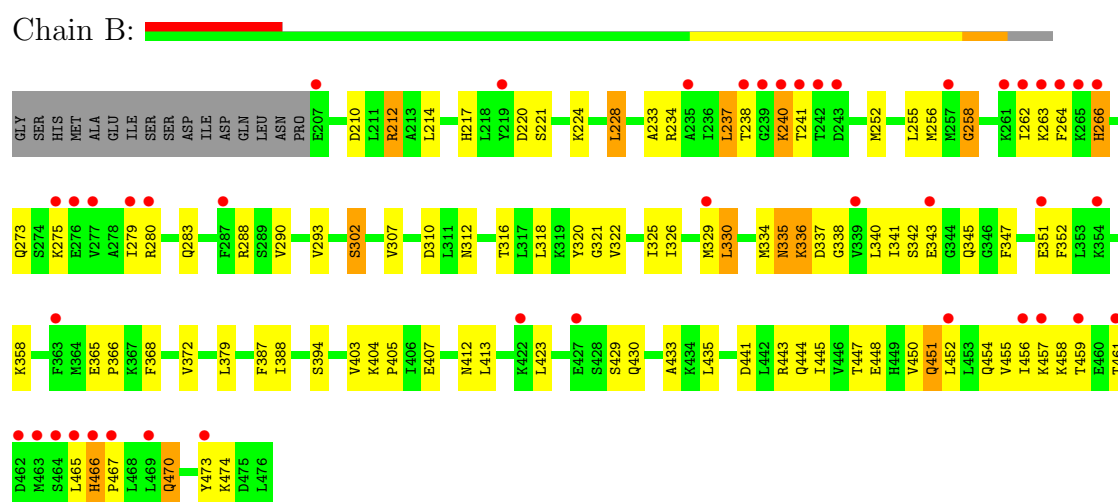
3 Residue-property plots

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: Peroxisome proliferator-activatedreceptor gamma



- Molecule 1: Peroxisome proliferator-activatedreceptor gamma



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	93.54Å 61.05Å 119.24Å 90.00° 103.49° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40 24.79 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.40) 98.1 (24.79-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.41Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.248 , 0.305 0.244 , 0.295	Depositor DCC
R_{free} test set	2476 reflections (9.90%)	DCC
Wilson B-factor (Å ²)	48.8	Xtriage
Anisotropy	0.632	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 65.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 25328 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4504	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.39% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.39	0/2223	0.58	0/2993
1	B	0.35	0/2203	0.56	0/2967
All	All	0.37	0/4426	0.57	0/5960

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	2180	0	2251	87	0
1	B	2166	0	2232	77	0
2	A	24	0	17	5	0
3	A	68	0	0	12	0
3	B	66	0	0	12	0
All	All	4504	0	4500	165	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 20.

All (165) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:341:ILE:HG22	2:A:1:LRG:HAX	1.38	1.01
1:A:288[B]:ARG:NH1	1:A:291[B]:GLU:HG2	1.76	1.01
1:B:325:ILE:HD12	1:B:388:ILE:HG23	1.45	0.99
1:A:226:PHE:HA	3:A:530:HOH:O	1.66	0.94
1:A:288[B]:ARG:HH11	1:A:291[B]:GLU:HG2	1.37	0.89
1:A:286:GLN:NE2	1:A:465:LEU:HD12	1.90	0.86
1:A:325:ILE:HG23	1:A:388:ILE:HD12	1.60	0.83
1:B:336:LYS:HE3	3:B:96:HOH:O	1.77	0.82
1:B:387:PHE:HB2	3:B:37:HOH:O	1.81	0.79
1:A:288[B]:ARG:NH1	1:A:291[B]:GLU:CG	2.46	0.79
1:B:451:GLN:O	1:B:454:GLN:HG2	1.85	0.77
1:B:266:HIS:HB2	3:B:107:HOH:O	1.83	0.76
1:B:465:LEU:C	1:B:467:PRO:HD2	2.07	0.75
1:B:455:VAL:HA	1:B:458:LYS:HE2	1.69	0.74
1:A:473:TYR:O	1:A:474:LYS:HD2	1.89	0.73
1:B:212:ARG:HH11	1:B:212:ARG:HB3	1.52	0.72
1:A:437:GLN:O	1:A:440:THR:HG22	1.88	0.71
1:A:370:PHE:HA	1:A:373:LYS:HE2	1.72	0.71
1:B:279:ILE:O	1:B:283:GLN:HG3	1.92	0.70
1:A:452:LEU:O	1:A:456:ILE:HD13	1.91	0.69
1:A:238:THR:O	1:A:240:LYS:N	2.26	0.69
1:A:349:THR:HG22	1:A:352:PHE:H	1.58	0.68
1:A:286:GLN:HE22	1:A:465:LEU:HD12	1.57	0.68
3:A:534:HOH:O	1:B:430:GLN:HB2	1.94	0.68
1:A:286:GLN:HE22	1:A:465:LEU:HA	1.58	0.67
1:A:287:PHE:O	1:A:290:VAL:HG12	1.94	0.66
1:B:275:LYS:HE3	1:B:275:LYS:HA	1.77	0.65
1:B:447:THR:O	1:B:450:VAL:HG22	1.99	0.63
1:B:335:ASN:ND2	1:B:337:ASP:H	1.98	0.62
1:B:252:MET:O	1:B:256:MET:HG2	2.00	0.61
1:B:307:VAL:HG22	3:B:12:HOH:O	2.02	0.60
1:B:466:HIS:N	1:B:467:PRO:HD2	2.16	0.60
1:B:220:ASP:O	1:B:224:LYS:HG2	2.02	0.59
1:B:457:LYS:NZ	1:B:461:THR:HG22	2.16	0.59
1:A:286:GLN:NE2	1:A:466:HIS:H	2.00	0.59
1:B:336:LYS:HB2	3:B:96:HOH:O	2.03	0.59
1:B:329:MET:HE3	3:B:59:HOH:O	2.02	0.58
1:B:466:HIS:N	1:B:467:PRO:CD	2.66	0.58
1:B:256:MET:HE1	1:B:280:ARG:HH12	1.69	0.58
1:B:335:ASN:ND2	3:B:96:HOH:O	2.37	0.57
1:B:264:PHE:CE2	1:B:266:HIS:HB3	2.39	0.57
1:A:253:ASN:HB2	3:A:504:HOH:O	2.05	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:310:ASP:OD2	1:B:312:ASN:HB2	2.05	0.57
1:B:365:GLU:HB3	3:B:110:HOH:O	2.04	0.57
1:A:276:GLU:CD	1:A:357:ARG:HH21	2.07	0.57
1:A:238:THR:C	1:A:240:LYS:H	2.09	0.56
1:A:207:GLU:HG3	1:A:209:ALA:H	1.70	0.56
1:A:279:ILE:O	1:A:283:GLN:HG2	2.06	0.56
1:B:379:LEU:HD21	1:B:435:LEU:HD13	1.87	0.55
1:A:305:GLY:HA2	1:A:308:ASN:HD22	1.71	0.55
1:B:452:LEU:O	1:B:456:ILE:HG12	2.07	0.55
1:A:276:GLU:HG2	1:A:279:ILE:HG12	1.88	0.55
1:A:336:LYS:HE2	1:A:372:VAL:HG21	1.88	0.54
1:A:444:GLN:O	1:A:448:GLU:HG3	2.08	0.54
1:A:289:SER:O	1:A:293:VAL:HG23	2.07	0.54
1:B:217:HIS:HE1	1:B:302:SER:O	1.90	0.54
1:B:368:PHE:O	1:B:372:VAL:HG23	2.08	0.53
1:A:242:THR:O	1:A:242:THR:HG22	2.08	0.53
1:A:433:ALA:O	1:A:437:GLN:HG3	2.09	0.53
1:B:448:GLU:O	1:B:451:GLN:HB3	2.09	0.52
1:B:351:GLU:HG3	3:B:51:HOH:O	2.10	0.52
1:A:435:LEU:O	1:A:438:LYS:HB2	2.09	0.52
2:A:1:LRG:HAQA	2:A:1:LRG:CAN	2.38	0.52
1:B:455:VAL:HA	1:B:458:LYS:CE	2.39	0.52
1:A:363:PHE:CD1	1:A:452:LEU:HD13	2.45	0.51
1:A:456:ILE:HG21	1:A:463:MET:CE	2.41	0.51
1:A:292:ALA:O	1:A:296:ILE:HG13	2.11	0.51
1:A:466:HIS:ND1	1:A:467:PRO:HD2	2.26	0.51
1:B:365:GLU:N	1:B:366:PRO:HD2	2.26	0.51
1:A:320:TYR:CB	1:A:397:ARG:HD2	2.39	0.51
1:B:441:ASP:O	1:B:445:ILE:HG12	2.10	0.51
1:B:340:LEU:O	1:B:341:ILE:HD12	2.11	0.50
1:A:226:PHE:HD1	3:A:530:HOH:O	1.94	0.50
3:A:512:HOH:O	1:B:429:SER:HB3	2.11	0.50
1:B:430:GLN:HG3	1:B:433:ALA:HB3	1.93	0.50
1:A:325:ILE:HD11	1:A:392:ILE:HG13	1.94	0.50
1:A:325:ILE:HD11	1:A:392:ILE:CG1	2.42	0.50
1:A:336:LYS:HG2	1:A:372:VAL:CG2	2.41	0.50
1:A:277:VAL:HA	1:A:280:ARG:HD3	1.94	0.49
1:A:370:PHE:HB2	1:A:445:ILE:HD11	1.94	0.49
1:A:379:LEU:HD11	1:A:435:LEU:HD13	1.95	0.49
1:B:290:VAL:HG21	1:B:473:TYR:CD1	2.48	0.49
1:A:425:HIS:N	1:A:426:PRO:HD3	2.28	0.49
1:A:393:LEU:O	1:A:410:GLN:HB2	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:257:MET:O	1:A:261:LYS:HG2	2.13	0.48
1:A:404:LYS:N	1:A:405:PRO:HD2	2.29	0.48
1:B:212:ARG:HD2	3:B:66:HOH:O	2.14	0.48
1:B:255:LEU:HD22	1:B:352:PHE:CZ	2.48	0.48
1:A:317:LEU:HD11	1:A:406:ILE:HD13	1.95	0.47
1:A:359:PRO:HG2	1:A:360:PHE:CD1	2.49	0.47
1:B:403:VAL:HG12	1:B:407:GLU:HG3	1.96	0.47
1:B:466:HIS:O	1:B:470:GLN:HB3	2.14	0.47
1:A:242:THR:HA	3:A:541:HOH:O	2.13	0.47
1:A:307:VAL:HG22	3:A:480:HOH:O	2.13	0.47
1:B:214:LEU:HD21	1:B:413:LEU:HD23	1.96	0.47
1:B:275:LYS:CE	1:B:275:LYS:HA	2.45	0.46
1:B:338:GLY:HA3	1:B:347:PHE:CZ	2.51	0.46
1:A:395:GLY:HA2	1:A:400:LEU:CD1	2.45	0.46
1:B:263:LYS:HE3	1:B:263:LYS:HA	1.98	0.46
1:B:293:VAL:HG22	1:B:322:VAL:HG11	1.97	0.46
1:A:320:TYR:HB3	1:A:397:ARG:HD2	1.98	0.46
1:A:336:LYS:CD	1:A:372:VAL:HG21	2.46	0.46
1:A:336:LYS:HD3	1:A:350:ARG:NH1	2.32	0.45
1:B:237:LEU:HB3	1:B:238:THR:H	1.58	0.45
1:B:275:LYS:HG2	1:B:280:ARG:HD2	1.99	0.45
1:A:336:LYS:HD3	1:A:350:ARG:HH12	1.80	0.45
1:A:349:THR:HG22	1:A:351:GLU:N	2.31	0.45
1:A:288[B]:ARG:HD2	1:A:288[B]:ARG:HA	1.65	0.45
1:A:357:ARG:HD2	1:A:359:PRO:HD2	1.98	0.45
1:A:447:THR:HG22	1:A:451:GLN:HE21	1.80	0.45
1:B:451:GLN:HG2	1:B:452:LEU:N	2.30	0.45
1:B:321:GLY:O	1:B:325:ILE:HG12	2.17	0.45
1:B:262:ILE:HD11	1:B:345:GLN:HB2	1.99	0.45
1:B:326:ILE:O	1:B:330:LEU:HB2	2.17	0.44
1:A:474:LYS:HE2	3:A:528:HOH:O	2.17	0.44
1:A:345:GLN:HG3	3:A:535:HOH:O	2.16	0.44
1:A:456:ILE:N	1:A:456:ILE:HD12	2.33	0.44
1:A:252:MET:HG2	3:A:543:HOH:O	2.17	0.44
1:B:470:GLN:HE21	1:B:470:GLN:HB3	1.54	0.43
1:A:208:SER:HB2	1:A:419:LEU:HD21	2.00	0.43
1:A:447:THR:O	1:A:451:GLN:HG2	2.18	0.43
2:A:1:LRG:HAO	2:A:1:LRG:HAK	1.82	0.43
1:B:288:ARG:O	1:B:288:ARG:HD2	2.18	0.43
2:A:1:LRG:HAL	2:A:1:LRG:HAP	1.81	0.43
1:A:384:LEU:O	1:A:388:ILE:HG12	2.18	0.43
1:A:336:LYS:CE	1:A:372:VAL:HG21	2.49	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:468:LEU:HD13	1:A:468:LEU:C	2.39	0.43
1:B:256:MET:C	1:B:258:GLY:H	2.22	0.43
1:B:343:GLU:HG3	1:B:343:GLU:O	2.18	0.43
1:A:414:LEU:HB2	1:B:430:GLN:HG2	2.01	0.42
1:B:337:ASP:N	3:B:96:HOH:O	2.52	0.42
1:A:288[B]:ARG:HH12	1:A:291[B]:GLU:HG2	1.75	0.42
1:B:341:ILE:CG2	1:B:342:SER:N	2.82	0.42
1:B:341:ILE:HG23	1:B:342:SER:N	2.35	0.42
1:A:365:GLU:N	1:A:366:PRO:HD2	2.35	0.42
1:B:435:LEU:HD21	3:B:37:HOH:O	2.20	0.42
1:A:319:LYS:HD3	1:A:320:TYR:CE1	2.55	0.42
1:B:290:VAL:HG21	1:B:473:TYR:CE1	2.55	0.42
1:B:316:THR:O	1:B:320:TYR:HD1	2.02	0.42
1:B:228:LEU:HD22	1:B:233:ALA:HB2	2.03	0.41
1:A:293:VAL:HG22	1:A:322:VAL:HG11	2.02	0.41
1:A:319:LYS:O	1:A:472:ILE:HG23	2.20	0.41
1:B:336:LYS:H	1:B:336:LYS:HD3	1.85	0.41
1:A:336:LYS:HG2	1:A:372:VAL:HG22	2.01	0.41
1:B:457:LYS:HZ3	1:B:461:THR:HG22	1.84	0.41
1:A:288[A]:ARG:HG3	2:A:1:LRG:HAI	2.01	0.41
1:A:465:LEU:HD23	1:A:470:GLN:HG3	2.02	0.41
1:A:456:ILE:HG21	1:A:463:MET:HE1	2.01	0.41
1:A:311:LEU:HD23	1:A:311:LEU:C	2.41	0.41
1:B:336:LYS:HE3	1:B:336:LYS:HB2	1.96	0.41
1:A:349:THR:HG23	3:A:484:HOH:O	2.20	0.41
1:A:358:LYS:HB2	1:A:359:PRO:HD3	2.03	0.41
1:B:457:LYS:HZ2	1:B:461:THR:HG22	1.83	0.41
1:B:240:LYS:O	1:B:241:THR:C	2.60	0.41
1:B:456:ILE:O	1:B:459:THR:HB	2.21	0.40
1:A:227:PRO:HD3	3:A:530:HOH:O	2.22	0.40
1:B:404:LYS:HB3	1:B:405:PRO:CD	2.51	0.40
1:A:222:TYR:CE2	1:A:381:ASP:HB3	2.57	0.40
1:B:212:ARG:HE	1:B:423:LEU:CD1	2.35	0.40
1:A:414:LEU:CB	1:B:430:GLN:HG2	2.51	0.40
1:B:262:ILE:HD11	1:B:345:GLN:CB	2.51	0.40
1:A:208:SER:HB2	1:A:419:LEU:CD2	2.51	0.40
1:B:234:ARG:NH2	1:B:334:MET:O	2.46	0.40
1:A:449:HIS:O	1:A:453:LEU:HD13	2.22	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	270/286 (94%)	251 (93%)	12 (4%)	7 (3%)	8	8
1	B	268/286 (94%)	237 (88%)	24 (9%)	7 (3%)	8	8
All	All	538/572 (94%)	488 (91%)	36 (7%)	14 (3%)	8	8

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	239	GLY
1	B	474	LYS
1	B	237	LEU
1	B	266	HIS
1	B	394	SER
1	A	275	LYS
1	B	240	LYS
1	A	269	PRO
1	A	336	LYS
1	A	241	THR
1	A	358	LYS
1	A	359	PRO
1	B	258	GLY
1	B	466	HIS

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	245/257 (95%)	233 (95%)	12 (5%)	35	53
1	B	243/257 (95%)	227 (93%)	16 (7%)	24	35

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	488/514 (95%)	460 (94%)	28 (6%)	30	44

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	221	SER
1	A	241	THR
1	A	253	ASN
1	A	271	GLN
1	A	291[A]	GLU
1	A	291[B]	GLU
1	A	362	ASP
1	A	363	PHE
1	A	404	LYS
1	A	427	GLU
1	A	452	LEU
1	A	460	GLU
1	B	210	ASP
1	B	212	ARG
1	B	221	SER
1	B	228	LEU
1	B	273	GLN
1	B	302	SER
1	B	318	LEU
1	B	330	LEU
1	B	335	ASN
1	B	336	LYS
1	B	358	LYS
1	B	412	ASN
1	B	443	ARG
1	B	444	GLN
1	B	451	GLN
1	B	470	GLN

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	286	GLN
1	A	308	ASN
1	A	314	GLN
1	A	410	GLN

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Mol	Chain	Res	Type
1	A	451	GLN
1	A	454	GLN
1	A	470	GLN
1	B	217	HIS
1	B	253	ASN
1	B	335	ASN
1	B	412	ASN
1	B	415	GLN
1	B	430	GLN
1	B	437	GLN
1	B	451	GLN
1	B	470	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 ⓘ

1 ligand is 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	LRG	A	1	-	26,26,26	0.88	1 (3%)	34,34,34	1.95	2 (5%)

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	LRG	A	1	-	1/1/2/2	0/16/16/16	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	LRG	CAW-CAV	-3.45	1.39	1.49

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	LRG	CAU-OAR-CAX	-10.44	112.89	118.83
2	A	1	LRG	OAR-CAX-CAQ	2.19	111.74	107.44

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1	LRG	CAX

There are no torsion outliers.

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 ⓘ

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	256/286 (89%)	0.67	31 (12%)	5 4	29, 51, 93, 100	0
1	B	256/286 (89%)	1.01	43 (16%)	2 1	27, 53, 100, 100	2 (0%)
All	All	512/572 (89%)	0.84	74 (14%)	3 2	27, 52, 100, 100	2 (0%)

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	461	THR	11.6
1	B	465	LEU	10.7
1	B	462	ASP	9.5
1	A	240	LYS	9.0
1	B	242	THR	8.4
1	B	463	MET	8.2
1	B	263	LYS	8.1
1	B	257	MET	7.7
1	B	261	LYS	7.4
1	B	241	THR	7.0
1	A	239	GLY	6.7
1	B	264	PHE	6.6
1	B	464	SER	6.1
1	B	238	THR	5.9
1	B	266	HIS	5.9
1	B	240	LYS	5.9
1	B	459	THR	5.9
1	A	260	ASP	4.6
1	B	427	GLU	4.4
1	A	455	VAL	4.2
1	B	207	GLU	4.2
1	B	265	LYS	4.0
1	B	456	ILE	3.9
1	B	243	ASP	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	351	GLU	3.7
1	B	262	ILE	3.7
1	B	343	GLU	3.7
1	A	243	ASP	3.4
1	B	466	HIS	3.4
1	B	276	GLU	3.3
1	A	280	ARG	3.2
1	A	358	LYS	3.2
1	A	256	MET	3.2
1	B	469	LEU	3.1
1	A	287	PHE	3.1
1	A	401	LEU	3.1
1	A	320	TYR	3.1
1	B	452	LEU	3.1
1	A	241	THR	3.0
1	B	239	GLY	3.0
1	B	422	LYS	3.0
1	A	242	THR	2.9
1	B	280	ARG	2.8
1	B	457	LYS	2.8
1	A	275	LYS	2.8
1	B	354	LYS	2.8
1	A	391	ILE	2.8
1	B	275	LYS	2.7
1	A	277	VAL	2.7
1	A	402	ASN	2.7
1	A	285	CYS	2.6
1	B	287	PHE	2.6
1	A	444	GLN	2.6
1	A	454	GLN	2.5
1	B	467	PRO	2.5
1	A	388	ILE	2.5
1	B	279	ILE	2.5
1	B	363	PHE	2.5
1	A	294	GLN	2.4
1	A	227	PRO	2.4
1	A	451	GLN	2.3
1	A	422	LYS	2.3
1	A	262	ILE	2.3
1	A	252	MET	2.2
1	B	473	TYR	2.2
1	B	219	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	277	VAL	2.1
1	B	339	VAL	2.1
1	A	465	LEU	2.1
1	B	235	ALA	2.1
1	A	336	LYS	2.1
1	A	291[A]	GLU	2.0
1	B	329	MET	2.0
1	A	325	ILE	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	LRG	A	1	24/24	0.40	3.32	78,85,87,87	0

6.5 Other polymers ⓘ

There are no such residues in this entry.