



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:10 AM GMT

PDB ID : 2G0H  
Title : Structure-based drug design of a novel family of PPAR partial agonists: virtual screening, x-ray crystallography and in vitro/in vivo biological activities  
Authors : Lu, I.L.; Peng, Y.H.; Huang, C.F.; Lin, Y.T.; Hsu, J.T.A.; Wu, S.Y.  
Deposited on : 2006-02-13  
Resolution : 2.30 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 (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

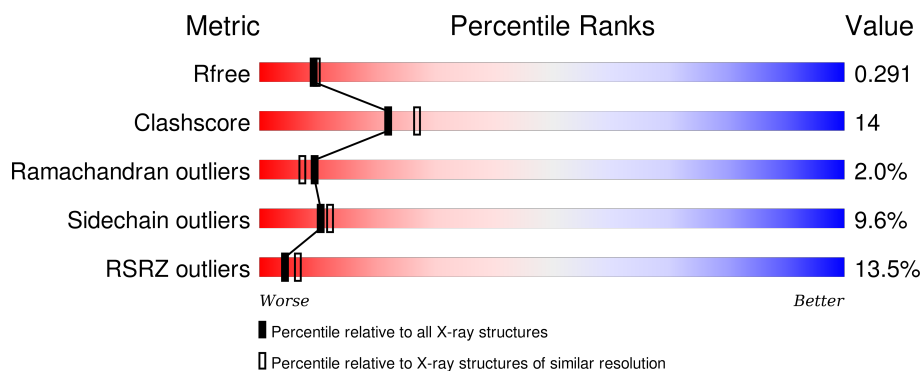
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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.

Mol	Chain	Length	Quality of chain
1	A	271	<div> <div>15%</div> <div>63%</div> <div>30%</div> <div>7%</div> </div>
1	B	271	<div> <div>12%</div> <div>74%</div> <div>21%</div> <div>5%</div> </div>

## 2 Entry composition [i](#)

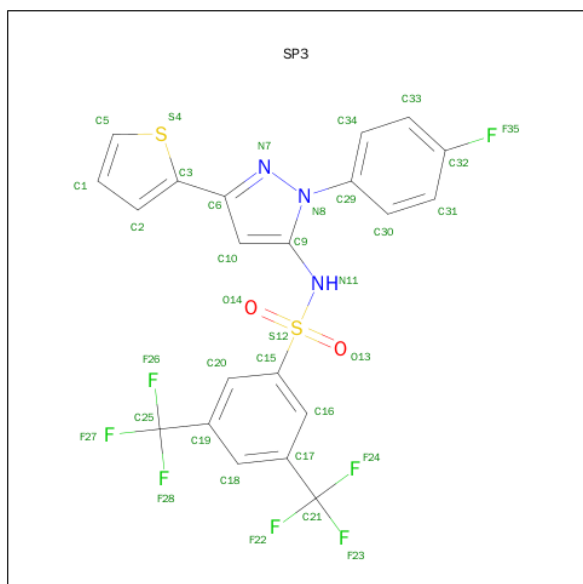
There are 3 unique types of molecules in this entry. The entry contains 4511 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Peroxisome proliferator-activated receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2177	1406	355	406	10			
1	B	271	Total	C	N	O	S	0	0	0
			2177	1406	355	406	10			

- Molecule 2 is N-[1-(4-FLUOROPHENYL)-3-(2-THIENYL)-1H-PYRAZOL-5-YL]-3,5-BIS(TRIFLUOROMETHYL)BENZENESULFONAMIDE (three-letter code: SP3) (formula:  $C_{21}H_{12}F_7N_3O_2S_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0
			35	21	7	3	2	2	
2	B	1	Total	C	F	N	O	S	0
			35	21	7	3	2	2	

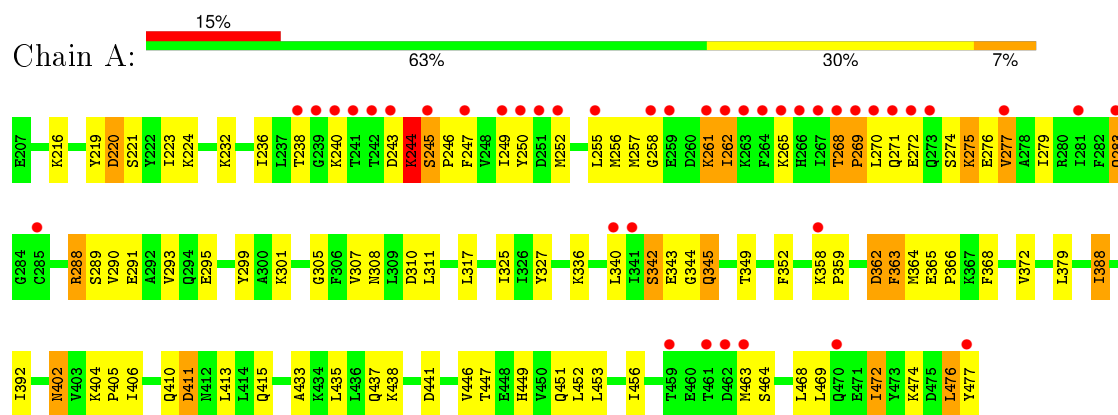
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	23	Total 23	O 23	0	0
3	B	64	Total 64	O 64	0	0

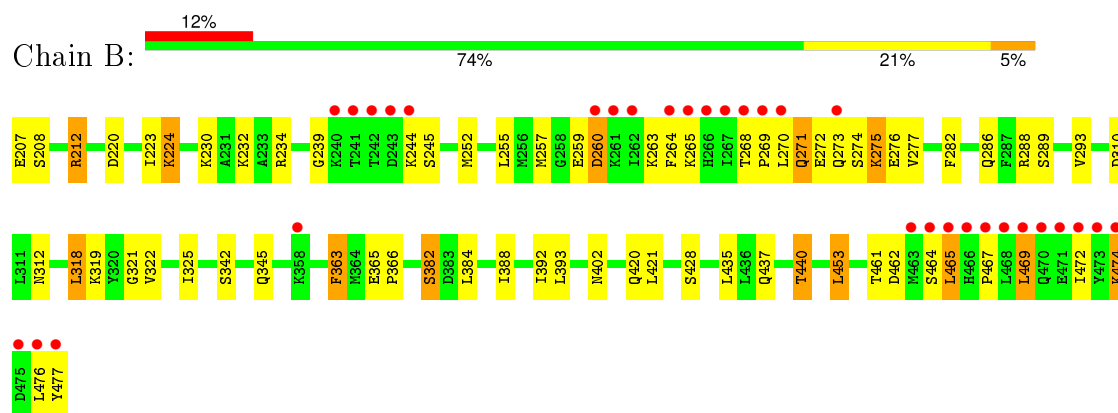
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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-activated receptor gamma



- Molecule 1: Peroxisome proliferator-activated receptor gamma



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.30Å 88.90Å 58.45Å 90.00° 90.80° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.90 – 2.19	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.30) 99.1 (19.90-2.19)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.00 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.238 , 0.293 0.237 , 0.291	Depositor DCC
$R_{free}$ test set	1320 reflections (5.44%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.7	Xtriage
Anisotropy	0.148	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 45.9	EDS
Estimated twinning fraction	0.000 for l,k,-h 0.029 for h,-k,-l 0.015 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 29417 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4511	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.12% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: SP3

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	0.40	0/2215	0.83	1/2985 (0.0%)
1	B	0.48	0/2215	0.90	2/2985 (0.1%)
All	All	0.44	0/4430	0.87	3/5970 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	243	ASP	N-CA-CB	7.37	123.87	110.60
1	B	212	ARG	NE-CZ-NH2	-5.69	117.46	120.30
1	B	207	GLU	OE1-CD-OE2	-5.18	117.08	123.30

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	2177	0	2241	67	0
1	B	2177	0	2241	56	0
2	A	35	0	12	4	0
2	B	35	0	12	4	0
3	A	23	0	0	2	0
3	B	64	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4511	0	4506	124	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 14.

All (124) 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:268:THR:HG22	1:B:268:THR:O	1.68	0.94
1:B:259:GLU:O	1:B:263:LYS:HB3	1.78	0.83
1:B:271:GLN:HA	1:B:271:GLN:HE21	1.44	0.80
1:B:275:LYS:HE3	1:B:276:GLU:H	1.44	0.80
1:B:271:GLN:HA	1:B:271:GLN:NE2	1.97	0.79
1:B:286:GLN:HE22	1:B:465:LEU:HB3	1.47	0.78
1:B:263:LYS:HE2	1:B:269:PRO:HG3	1.66	0.77
1:B:293:VAL:HG22	1:B:322:VAL:HG11	1.67	0.76
1:B:363:PHE:HZ	1:B:453:LEU:HD13	1.51	0.75
1:B:268:THR:O	1:B:268:THR:CG2	2.35	0.74
2:B:201:SP3:H11	2:B:201:SP3:H34	1.52	0.74
2:A:101:SP3:H11	2:A:101:SP3:H34	1.54	0.72
1:B:469:LEU:HA	1:B:472:ILE:HB	1.71	0.72
1:B:275:LYS:HD3	1:B:462:ASP:OD1	1.91	0.71
1:A:262:ILE:HG23	1:A:265:LYS:O	1.94	0.67
2:B:201:SP3:H11	2:B:201:SP3:C34	2.06	0.67
1:A:311:LEU:HD23	1:B:271:GLN:HG2	1.77	0.66
1:A:438:LYS:HE3	3:A:79:HOH:O	1.94	0.66
1:A:247:PHE:CZ	1:A:257:MET:HG2	2.31	0.65
1:A:301:LYS:NZ	1:B:467:PRO:HD2	2.11	0.65
1:A:364:MET:HE3	2:A:101:SP3:H10	1.81	0.62
1:A:305:GLY:HA2	1:A:308:ASN:HD22	1.63	0.62
1:A:247:PHE:HZ	1:A:257:MET:HG2	1.66	0.61
1:A:325:ILE:HG23	1:A:388:ILE:HG12	1.83	0.60
1:A:410:GLN:HA	1:A:413:LEU:HD12	1.82	0.60
2:A:101:SP3:H11	2:A:101:SP3:C34	2.14	0.59
1:A:255:LEU:CD2	1:A:277:VAL:HG13	2.33	0.59
1:A:275:LYS:HG3	1:A:283:GLN:NE2	2.18	0.58
1:A:238:THR:CG2	1:A:240:LYS:HG2	2.33	0.58
1:A:365:GLU:HB3	1:A:366:PRO:CD	2.33	0.58
1:B:220:ASP:O	1:B:224:LYS:HG3	2.05	0.56
1:A:255:LEU:HD21	1:A:277:VAL:HG13	1.86	0.56
2:B:201:SP3:N11	2:B:201:SP3:H34	2.16	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:ASP:OD2	1:B:269:PRO:HD2	2.06	0.56
1:B:472:ILE:O	1:B:472:ILE:HG22	2.06	0.56
1:A:220:ASP:O	1:A:224:LYS:HG3	2.06	0.56
1:A:252:MET:O	1:A:256:MET:HG3	2.06	0.55
1:A:325:ILE:HD11	1:A:392:ILE:HG13	1.89	0.54
1:B:293:VAL:HG22	1:B:322:VAL:CG1	2.38	0.53
1:B:363:PHE:CZ	1:B:453:LEU:HD13	2.39	0.53
1:A:238:THR:HG21	1:A:240:LYS:HE3	1.89	0.53
1:A:472:ILE:O	1:A:476:LEU:HD13	2.08	0.53
1:A:379:LEU:HD11	1:A:435:LEU:HD13	1.90	0.53
1:B:255:LEU:CD2	1:B:277:VAL:HG13	2.38	0.53
1:B:274:SER:O	1:B:275:LYS:HB3	2.09	0.52
1:A:327:TYR:HE2	1:A:446:VAL:HG22	1.74	0.52
1:B:271:GLN:CD	1:B:272:GLU:H	2.13	0.52
1:B:325:ILE:HD12	1:B:388:ILE:HG23	1.92	0.52
1:B:474:LYS:HB3	1:B:474:LYS:NZ	2.25	0.51
1:A:435:LEU:O	1:A:438:LYS:HB2	2.11	0.51
1:A:301:LYS:HZ2	1:B:467:PRO:HD2	1.75	0.51
1:A:342:SER:O	1:A:343:GLU:HB2	2.11	0.51
1:A:404:LYS:N	1:A:405:PRO:HD2	2.26	0.51
1:A:238:THR:HG22	1:A:240:LYS:HG2	1.92	0.51
1:A:447:THR:O	1:A:451:GLN:HB2	2.11	0.50
1:A:290:VAL:HG13	1:A:468:LEU:HD23	1.92	0.50
1:A:301:LYS:HZ1	1:B:467:PRO:HD2	1.77	0.49
1:B:252:MET:SD	1:B:277:VAL:HG11	2.52	0.49
1:B:382:SER:OG	1:B:420:GLN:NE2	2.46	0.49
1:B:437:GLN:O	1:B:440:THR:HG23	2.13	0.48
1:B:363:PHE:HZ	1:B:453:LEU:CD1	2.22	0.48
1:B:472:ILE:CG2	1:B:472:ILE:O	2.62	0.48
1:B:319:LYS:HE3	1:B:472:ILE:HG23	1.95	0.48
1:B:321:GLY:O	1:B:325:ILE:HG12	2.14	0.48
1:B:365:GLU:N	1:B:366:PRO:CD	2.77	0.47
1:A:288:ARG:HA	1:A:288:ARG:HD2	1.64	0.47
1:A:275:LYS:HB3	1:A:276:GLU:H	1.54	0.47
1:B:318:LEU:O	1:B:322:VAL:HG23	2.14	0.47
1:B:384:LEU:O	1:B:388:ILE:HG12	2.16	0.46
1:A:411:ASP:O	1:A:415:GLN:HG3	2.16	0.46
1:A:275:LYS:HG3	1:A:283:GLN:HE22	1.80	0.46
1:A:258:GLY:O	1:A:261:LYS:HB3	2.16	0.45
1:A:272:GLU:O	1:A:274:SER:N	2.48	0.45
1:B:282:PHE:CE1	2:B:201:SP3:H30	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:101:SP3:H34	2:A:101:SP3:N11	2.26	0.45
1:A:238:THR:HG21	1:A:240:LYS:HG2	1.98	0.45
1:A:291:GLU:O	1:A:295:GLU:HG3	2.16	0.45
1:A:232:LYS:O	1:A:236:ILE:HG12	2.16	0.45
1:B:402:ASN:N	1:B:402:ASN:ND2	2.63	0.45
1:B:208:SER:O	1:B:212:ARG:HG2	2.16	0.45
1:B:342:SER:O	1:B:345:GLN:HB2	2.17	0.45
1:A:247:PHE:HZ	1:A:257:MET:CG	2.27	0.45
1:A:433:ALA:O	1:A:437:GLN:HG3	2.17	0.45
1:A:247:PHE:CZ	1:A:257:MET:CG	2.98	0.44
1:B:274:SER:O	1:B:275:LYS:CB	2.65	0.44
1:A:438:LYS:HD3	1:A:438:LYS:HA	1.78	0.44
1:A:289:SER:O	1:A:293:VAL:HG23	2.16	0.44
1:B:325:ILE:HD12	1:B:388:ILE:HD12	2.00	0.44
1:A:245:SER:HA	1:A:246:PRO:HD3	1.81	0.44
1:B:255:LEU:HD23	1:B:277:VAL:CG1	2.48	0.44
1:B:476:LEU:HB3	1:B:477:TYR:H	1.67	0.43
1:A:250:TYR:HA	1:A:352:PHE:HB2	2.00	0.43
1:B:273:GLN:OE1	1:B:273:GLN:HA	2.19	0.43
1:B:264:PHE:O	1:B:265:LYS:C	2.56	0.43
1:B:310:ASP:OD1	1:B:312:ASN:N	2.51	0.43
1:A:340:LEU:HB3	1:A:344:GLY:HA2	2.01	0.43
1:A:244:LYS:HB2	1:A:245:SER:H	1.53	0.42
1:B:230:LYS:O	1:B:234:ARG:HG2	2.19	0.42
1:A:265:LYS:NZ	1:A:345:GLN:HG3	2.35	0.42
1:B:255:LEU:HD23	1:B:277:VAL:HG13	2.02	0.42
1:A:279:ILE:HG22	1:A:283:GLN:HE22	1.83	0.42
1:A:469:LEU:O	1:A:472:ILE:HG22	2.18	0.42
1:A:307:VAL:HG22	3:A:80:HOH:O	2.19	0.42
1:A:261:LYS:HZ1	1:A:262:ILE:HD13	1.84	0.42
1:A:402:ASN:HD22	1:A:402:ASN:HA	1.65	0.42
1:A:268:THR:HB	1:A:269:PRO:HD2	2.01	0.42
1:A:362:ASP:OD1	1:A:362:ASP:N	2.53	0.42
1:A:268:THR:O	1:A:270:LEU:N	2.52	0.41
1:B:392:ILE:HG22	1:B:393:LEU:HD22	2.02	0.41
1:A:365:GLU:HB3	1:A:366:PRO:HD3	2.02	0.41
1:A:327:TYR:CE2	1:A:446:VAL:HG22	2.55	0.41
1:A:249:ILE:HG23	1:A:255:LEU:HA	2.01	0.41
1:A:317:LEU:HD11	1:A:406:ILE:HG12	2.01	0.41
1:B:421:LEU:HD11	1:B:435:LEU:HD23	2.03	0.41
1:B:271:GLN:NE2	1:B:271:GLN:CA	2.76	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:293:VAL:CG2	1:B:322:VAL:HG11	2.44	0.41
1:A:299:TYR:CE2	1:A:388:ILE:HG22	2.56	0.41
1:B:260:ASP:HA	1:B:264:PHE:HB3	2.03	0.41
1:B:232:LYS:NZ	3:B:76:HOH:O	2.53	0.41
1:A:368:PHE:O	1:A:372:VAL:HG23	2.21	0.41
1:A:219:TYR:O	1:A:223:ILE:HG12	2.21	0.41
1:B:289:SER:O	1:B:293:VAL:HG23	2.22	0.40
1:A:363:PHE:HE1	1:A:449:HIS:CE1	2.38	0.40
1:A:325:ILE:HD11	1:A:392:ILE:CG1	2.52	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	269/271 (99%)	228 (85%)	31 (12%)	10 (4%)	4	2
1	B	269/271 (99%)	242 (90%)	26 (10%)	1 (0%)	39	48
All	All	538/542 (99%)	470 (87%)	57 (11%)	11 (2%)	9	7

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	244	LYS
1	A	262	ILE
1	A	277	VAL
1	A	453	LEU
1	B	239	GLY
1	A	349	THR
1	A	359	PRO
1	A	464	SER
1	A	245	SER

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Mol	Chain	Res	Type
1	A	269	PRO
1	A	358	LYS

### 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	244/244 (100%)	218 (89%)	26 (11%)	8	9
1	B	244/244 (100%)	223 (91%)	21 (9%)	13	15
All	All	488/488 (100%)	441 (90%)	47 (10%)	10	12

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	216	LYS
1	A	220	ASP
1	A	221	SER
1	A	244	LYS
1	A	261	LYS
1	A	268	THR
1	A	271	GLN
1	A	275	LYS
1	A	283	GLN
1	A	288	ARG
1	A	336	LYS
1	A	342	SER
1	A	345	GLN
1	A	362	ASP
1	A	363	PHE
1	A	388	ILE
1	A	402	ASN
1	A	411	ASP
1	A	441	ASP
1	A	452	LEU
1	A	456	ILE

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Mol	Chain	Res	Type
1	A	463	MET
1	A	472	ILE
1	A	474	LYS
1	A	476	LEU
1	A	477	TYR
1	B	223	ILE
1	B	224	LYS
1	B	244	LYS
1	B	245	SER
1	B	257	MET
1	B	260	ASP
1	B	270	LEU
1	B	271	GLN
1	B	275	LYS
1	B	288	ARG
1	B	318	LEU
1	B	363	PHE
1	B	382	SER
1	B	428	SER
1	B	440	THR
1	B	453	LEU
1	B	461	THR
1	B	464	SER
1	B	465	LEU
1	B	469	LEU
1	B	474	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	283	GLN
1	A	294	GLN
1	A	308	ASN
1	A	312	ASN
1	A	402	ASN
1	A	412	ASN
1	A	437	GLN
1	A	451	GLN
1	B	271	GLN
1	B	283	GLN
1	B	286	GLN
1	B	402	ASN

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Mol	Chain	Res	Type
1	B	451	GLN

### 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 carbohydrates 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	SP3	A	101	-	36,38,38	2.58	8 (22%)	48,59,59	2.16	13 (27%)
2	SP3	B	201	-	36,38,38	2.55	8 (22%)	48,59,59	2.22	12 (25%)

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	SP3	A	101	-	-	0/26/31/31	0/4/4/4
2	SP3	B	201	-	-	0/26/31/31	0/4/4/4

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	101	SP3	N7-N8	-8.75	1.23	1.39
2	B	201	SP3	N7-N8	-8.63	1.23	1.39
2	A	101	SP3	C29-N8	-4.85	1.34	1.44
2	B	201	SP3	C29-N8	-4.72	1.34	1.44
2	A	101	SP3	C6-C3	-4.08	1.39	1.49
2	A	101	SP3	C10-C6	-3.72	1.33	1.40
2	B	201	SP3	C6-C3	-3.68	1.40	1.49
2	B	201	SP3	C10-C6	-3.47	1.34	1.40
2	A	101	SP3	C15-S12	-2.93	1.72	1.76
2	B	201	SP3	C21-C17	2.18	1.54	1.49
2	A	101	SP3	C25-C19	2.22	1.54	1.49
2	B	201	SP3	C25-C19	2.39	1.55	1.49
2	A	101	SP3	O13-S12	6.02	1.50	1.43
2	B	201	SP3	O13-S12	6.03	1.50	1.43
2	A	101	SP3	O14-S12	6.71	1.50	1.43
2	B	201	SP3	O14-S12	6.84	1.51	1.43

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	SP3	O14-S12-O13	-7.37	109.76	119.54
2	A	101	SP3	C1-C5-S4	-7.32	105.12	113.23
2	A	101	SP3	O14-S12-O13	-5.30	112.50	119.54
2	B	201	SP3	C1-C5-S4	-4.54	108.21	113.23
2	A	101	SP3	F24-C21-C17	-3.41	105.66	112.95
2	A	101	SP3	C16-C17-C21	-3.35	115.14	119.59
2	B	201	SP3	C29-N8-C9	-3.32	125.21	129.02
2	A	101	SP3	C16-C15-S12	-3.13	115.58	119.09
2	A	101	SP3	F28-C25-C19	-2.90	106.74	112.95
2	B	201	SP3	C10-C6-C3	-2.60	125.32	129.24
2	A	101	SP3	C29-N8-C9	-2.37	126.30	129.02
2	A	101	SP3	F23-C21-F22	2.12	113.36	105.71
2	B	201	SP3	C15-S12-N11	2.14	109.62	106.87
2	A	101	SP3	F26-C25-F28	2.15	113.48	105.71
2	B	201	SP3	C16-C15-S12	2.17	121.51	119.09
2	B	201	SP3	F22-C21-F24	2.42	114.44	105.71
2	B	201	SP3	F26-C25-F28	2.58	115.04	105.71
2	B	201	SP3	O13-S12-C15	2.89	111.61	107.96
2	A	101	SP3	C20-C15-S12	3.00	122.44	119.09
2	B	201	SP3	C17-C16-C15	3.01	120.23	118.84
2	A	101	SP3	C18-C17-C21	3.39	124.10	119.59
2	A	101	SP3	C3-C6-N7	3.41	125.34	120.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	101	SP3	O13-S12-C15	4.63	113.81	107.96
2	B	201	SP3	C3-C6-N7	4.99	127.47	120.75
2	B	201	SP3	C19-C20-C15	6.76	121.96	118.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	101	SP3	4	0
2	B	201	SP3	4	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	271/271 (100%)	0.75	41 (15%) 3 5	23, 45, 119, 133	0
1	B	271/271 (100%)	0.54	32 (11%) 6 10	17, 34, 104, 122	0
All	All	542/542 (100%)	0.64	73 (13%) 4 6	17, 39, 109, 133	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	267	ILE	15.1
1	A	267	ILE	11.8
1	B	468	LEU	11.8
1	B	476	LEU	11.6
1	B	474	LYS	10.6
1	B	477	TYR	10.2
1	A	270	LEU	9.6
1	A	263	LYS	9.1
1	B	473	TYR	8.9
1	B	241	THR	8.1
1	B	470	GLN	7.5
1	B	266	HIS	7.1
1	A	269	PRO	7.0
1	B	469	LEU	6.8
1	A	477	TYR	6.6
1	B	260	ASP	6.4
1	A	264	PHE	6.3
1	A	242	THR	6.0
1	B	465	LEU	5.9
1	A	266	HIS	5.9
1	A	271	GLN	5.7
1	B	467	PRO	5.6
1	B	471	GLU	5.3
1	A	272	GLU	5.2

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Mol	Chain	Res	Type	RSRZ
1	A	268	THR	5.2
1	B	265	LYS	5.1
1	B	243	ASP	5.1
1	A	463	MET	5.0
1	A	283	GLN	4.8
1	B	273	GLN	4.7
1	B	268	THR	4.7
1	B	472	ILE	4.6
1	A	252	MET	4.6
1	A	262	ILE	4.4
1	A	459	THR	4.4
1	B	269	PRO	4.2
1	A	358	LYS	4.1
1	B	242	THR	4.1
1	B	240	LYS	4.0
1	B	264	PHE	4.0
1	A	250	TYR	4.0
1	B	463	MET	3.9
1	A	255	LEU	3.8
1	A	277	VAL	3.8
1	A	265	LYS	3.7
1	A	461	THR	3.7
1	A	462	ASP	3.7
1	B	261	LYS	3.5
1	A	273	GLN	3.4
1	B	262	ILE	3.4
1	A	241	THR	3.3
1	A	341	ILE	3.1
1	A	245	SER	3.0
1	B	466	HIS	2.9
1	B	358	LYS	2.9
1	B	475	ASP	2.8
1	A	281	ILE	2.7
1	B	464	SER	2.6
1	A	239	GLY	2.6
1	A	259	GLU	2.6
1	A	249	ILE	2.5
1	A	261	LYS	2.5
1	B	244	LYS	2.5
1	A	340	LEU	2.4
1	A	240	LYS	2.4
1	A	258	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	238	THR	2.3
1	A	470	GLN	2.2
1	A	243	ASP	2.2
1	A	285	CYS	2.2
1	A	251	ASP	2.1
1	B	270	LEU	2.1
1	A	247	PHE	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 carbohydrates 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. 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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
2	SP3	A	101	35/35	0.91	0.19	0.15	55,65,72,74	0
2	SP3	B	201	35/35	0.87	0.18	0.05	49,60,67,71	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.