



# Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 06:14 am BST

PDB ID : 6DGP  
Title : Crystal Structure of Human PPARgamma Ligand Binding Domain in Complex with TRAP220 Coactivator Peptide  
Authors : Shang, J.; Kojetin, D.J.  
Deposited on : 2018-05-17  
Resolution : 3.10 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

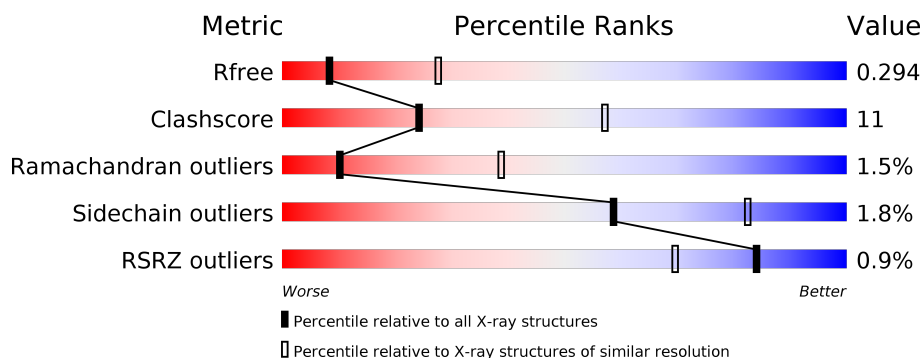
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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))
$R_{free}$	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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 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	297	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>%</span> <div style="width: 100%; height: 10px; background-color: green;"></div> </div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>61%</span> <span>25%</span> <span>• 13%</span> </div> </div>
1	B	297	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>%</span> <div style="width: 100%; height: 10px; background-color: green;"></div> </div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>67%</span> <span>20%</span> <span>13%</span> </div> </div>
2	C	19	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>42%</span> <span>21%</span> <span>37%</span> </div> </div>
2	D	19	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>5%</span> <div style="width: 100%; height: 10px; background-color: green;"></div> </div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>21%</span> <span>42%</span> <span>37%</span> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4346 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	259	Total	C	N	O	S	0	0	0
			2074	1340	337	387	10			
1	B	259	Total	C	N	O	S	0	0	0
			2074	1340	337	387	10			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	MET	-	initiating methionine	UNP P37231
A	182	ALA	-	expression tag	UNP P37231
A	183	HIS	-	expression tag	UNP P37231
A	184	HIS	-	expression tag	UNP P37231
A	185	HIS	-	expression tag	UNP P37231
A	186	HIS	-	expression tag	UNP P37231
A	187	HIS	-	expression tag	UNP P37231
A	188	HIS	-	expression tag	UNP P37231
A	189	VAL	-	expression tag	UNP P37231
A	190	ASP	-	expression tag	UNP P37231
A	191	ASP	-	expression tag	UNP P37231
A	192	ASP	-	expression tag	UNP P37231
A	193	ASP	-	expression tag	UNP P37231
A	194	LYS	-	expression tag	UNP P37231
A	195	MET	-	expression tag	UNP P37231
A	196	GLU	-	expression tag	UNP P37231
A	197	ASN	-	expression tag	UNP P37231
A	198	LEU	-	expression tag	UNP P37231
A	199	TYR	-	expression tag	UNP P37231
A	200	PHE	-	expression tag	UNP P37231
A	201	GLN	-	expression tag	UNP P37231
A	202	GLY	-	expression tag	UNP P37231
B	181	MET	-	initiating methionine	UNP P37231
B	182	ALA	-	expression tag	UNP P37231
B	183	HIS	-	expression tag	UNP P37231

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Chain	Residue	Modelled	Actual	Comment	Reference
B	184	HIS	-	expression tag	UNP P37231
B	185	HIS	-	expression tag	UNP P37231
B	186	HIS	-	expression tag	UNP P37231
B	187	HIS	-	expression tag	UNP P37231
B	188	HIS	-	expression tag	UNP P37231
B	189	VAL	-	expression tag	UNP P37231
B	190	ASP	-	expression tag	UNP P37231
B	191	ASP	-	expression tag	UNP P37231
B	192	ASP	-	expression tag	UNP P37231
B	193	ASP	-	expression tag	UNP P37231
B	194	LYS	-	expression tag	UNP P37231
B	195	MET	-	expression tag	UNP P37231
B	196	GLU	-	expression tag	UNP P37231
B	197	ASN	-	expression tag	UNP P37231
B	198	LEU	-	expression tag	UNP P37231
B	199	TYR	-	expression tag	UNP P37231
B	200	PHE	-	expression tag	UNP P37231
B	201	GLN	-	expression tag	UNP P37231
B	202	GLY	-	expression tag	UNP P37231

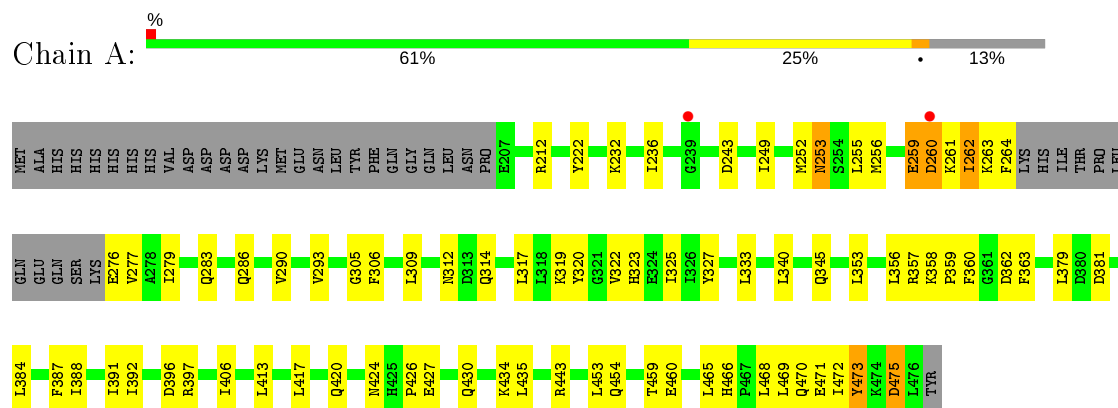
- Molecule 2 is a protein called TRAP220 Coactivator Peptide (Mediator of RNA polymerase II transcription subunit 1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	S	0	0	0
			99	63	18	16	2			
2	D	12	Total	C	N	O	S	0	0	0
			99	63	18	16	2			

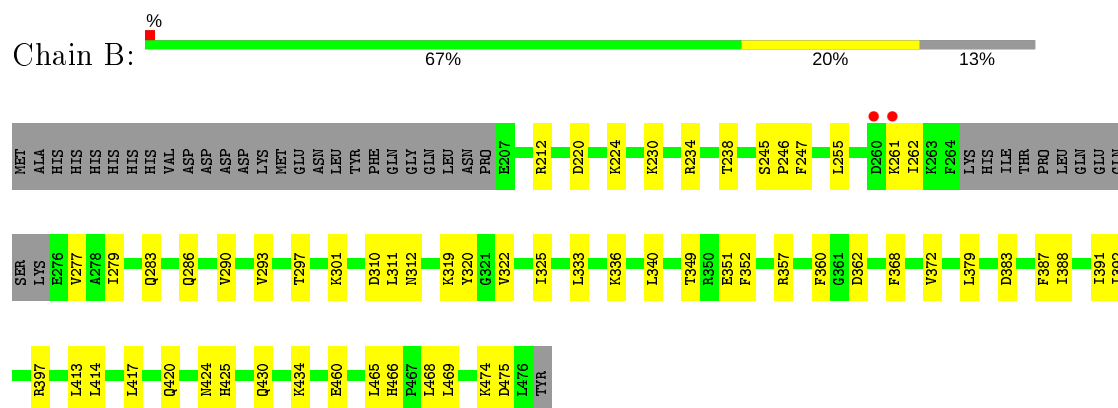
### 3 Residue-property plots

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

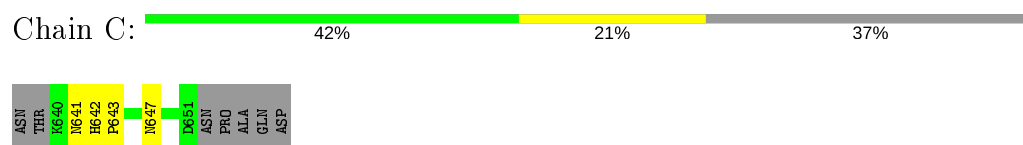
- Molecule 1: Peroxisome proliferator-activated receptor gamma



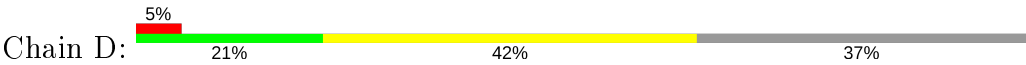
- Molecule 1: Peroxisome proliferator-activated receptor gamma



- Molecule 2: TRAP220 Coactivator Peptide (Mediator of RNA polymerase II transcription subunit 1)



- Molecule 2: TRAP220 Coactivator Peptide (Mediator of RNA polymerase II transcription subunit 1)



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.09 Å 54.73 Å 123.47 Å 90.00° 90.05° 90.00°	Depositor
Resolution (Å)	44.05 – 3.10 44.04 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.05-3.10) 99.6 (44.04-3.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.98 (at 3.12 Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, $R_{free}$	0.222 , 0.294 0.224 , 0.294	Depositor DCC
$R_{free}$ test set	1134 reflections (10.41%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	77.6	Xtriage
Anisotropy	0.317	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 36.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.460 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4346	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.85% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

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

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	0.49	0/2108	0.68	0/2839
1	B	0.47	0/2108	0.67	0/2839
2	C	0.33	0/100	0.62	0/132
2	D	0.91	0/100	1.05	0/132
All	All	0.49	0/4416	0.69	0/5942

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2074	0	2135	51	0
1	B	2074	0	2135	40	0
2	C	99	0	106	3	0
2	D	99	0	106	8	0
All	All	4346	0	4482	97	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (97) 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)
2:D:646:MET:HA	2:D:649:LEU:HD12	1.53	0.88
1:A:459:THR:HG23	1:A:460:GLU:HG2	1.58	0.84
1:A:325:ILE:HG23	1:A:388:ILE:HD12	1.68	0.73
1:B:234:ARG:O	1:B:238:THR:HG22	1.89	0.72
1:A:293:VAL:HG22	1:A:322:VAL:HG21	1.71	0.72
1:A:357:ARG:NH1	1:A:460:GLU:OE2	2.25	0.70
1:A:262:ILE:O	1:A:262:ILE:HG13	1.92	0.69
1:B:430:GLN:O	1:B:434:LYS:HG3	1.92	0.69
1:B:293:VAL:HG22	1:B:322:VAL:HG21	1.75	0.67
1:A:286:GLN:HG3	1:A:469:LEU:HD12	1.78	0.66
1:A:290:VAL:HG23	1:A:468:LEU:HD22	1.76	0.66
1:B:286:GLN:HG3	1:B:469:LEU:HD12	1.76	0.65
1:A:259:GLU:O	1:A:260:ASP:C	2.35	0.65
1:B:368:PHE:O	1:B:372:VAL:HG23	2.02	0.60
1:A:430:GLN:O	1:A:434:LYS:HG3	2.00	0.60
1:B:325:ILE:HD12	1:B:388:ILE:HB	1.85	0.59
1:A:325:ILE:HD12	1:A:388:ILE:HD12	1.86	0.58
1:A:333:LEU:HB3	1:A:340:LEU:HB2	1.84	0.58
1:B:212:ARG:HH12	1:B:420:GLN:HA	1.68	0.58
1:A:279:ILE:O	1:A:283:GLN:HG2	2.04	0.57
1:B:247:PHE:HB3	1:B:262:ILE:HD11	1.85	0.57
1:B:333:LEU:HB3	1:B:340:LEU:HB2	1.86	0.57
1:B:212:ARG:NH1	1:B:420:GLN:HA	2.19	0.56
1:B:357:ARG:HH21	1:B:360:PHE:HE2	1.53	0.56
1:A:325:ILE:HD12	1:A:388:ILE:HG23	1.87	0.55
1:A:259:GLU:N	1:A:259:GLU:OE1	2.29	0.55
1:A:286:GLN:HG3	1:A:469:LEU:CD1	2.36	0.55
1:B:279:ILE:O	1:B:283:GLN:HG2	2.07	0.55
1:B:413:LEU:O	1:B:417:LEU:N	2.34	0.54
1:B:255:LEU:HD22	1:B:277:VAL:HG23	1.88	0.54
1:A:454:GLN:NE2	1:A:475:ASP:OD1	2.39	0.54
1:B:286:GLN:HG3	1:B:469:LEU:CD1	2.39	0.53
1:B:311:LEU:HD21	2:D:646:MET:HG2	1.92	0.52
1:A:413:LEU:O	1:A:417:LEU:N	2.36	0.52
1:A:312:ASN:OD1	2:C:641:ASN:ND2	2.43	0.52
1:B:230:LYS:HE2	1:B:379:LEU:O	2.10	0.51
2:D:645:LEU:O	2:D:648:LEU:N	2.43	0.51
1:A:255:LEU:HD22	1:A:277:VAL:HG23	1.91	0.51
1:A:357:ARG:HH11	1:A:360:PHE:HE1	1.57	0.51
1:A:471:GLU:CD	2:C:642:HIS:HB3	2.31	0.51
1:B:297:THR:HG21	2:D:648:LEU:HB3	1.92	0.51
1:A:465:LEU:HD22	1:A:470:GLN:HB3	1.91	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:LEU:O	1:A:388:ILE:HG12	2.11	0.50
1:B:349:THR:HG22	1:B:351:GLU:H	1.76	0.50
1:B:290:VAL:HG23	1:B:468:LEU:HD23	1.94	0.49
1:B:383:ASP:OD2	1:B:425:HIS:NE2	2.32	0.49
1:A:319:LYS:HE3	1:A:472:ILE:O	2.12	0.49
2:D:650:LYS:O	2:D:651:ASP:C	2.50	0.49
1:A:259:GLU:HG3	1:A:264:PHE:HB2	1.94	0.49
2:D:645:LEU:O	2:D:648:LEU:HB2	2.13	0.48
1:A:252:MET:SD	1:A:277:VAL:HB	2.54	0.47
1:B:387:PHE:CZ	1:B:391:ILE:HD11	2.49	0.47
1:B:320:TYR:HB3	1:B:397:ARG:HD2	1.97	0.47
1:B:357:ARG:NH2	1:B:360:PHE:HE2	2.14	0.46
1:B:220:ASP:O	1:B:224:LYS:HG2	2.16	0.46
1:B:293:VAL:HG22	1:B:322:VAL:CG2	2.45	0.46
1:A:249:ILE:HG21	1:A:255:LEU:HD12	1.97	0.46
1:A:305:GLY:O	1:A:309:LEU:HD13	2.17	0.45
1:A:357:ARG:NH1	1:A:360:PHE:HE1	2.15	0.45
1:A:286:GLN:OE1	1:A:466:HIS:N	2.42	0.45
1:B:286:GLN:OE1	1:B:465:LEU:HA	2.17	0.45
1:B:312:ASN:ND2	2:D:641:ASN:OD1	2.50	0.45
1:A:353:LEU:HA	1:A:356:LEU:HD13	1.99	0.45
1:B:325:ILE:HG13	1:B:392:ILE:HD11	1.99	0.45
1:A:253:ASN:O	1:A:256:MET:HG2	2.16	0.45
1:A:306:PHE:O	1:A:314:GLN:NE2	2.49	0.44
1:A:469:LEU:O	1:A:472:ILE:HG12	2.17	0.44
1:A:212:ARG:HH12	1:A:420:GLN:HA	1.82	0.44
1:B:349:THR:HG22	1:B:351:GLU:N	2.31	0.44
1:A:468:LEU:HD23	1:A:469:LEU:HD23	2.00	0.44
2:C:643:PRO:O	2:C:647:ASN:ND2	2.48	0.44
1:B:349:THR:HB	1:B:352:PHE:HB3	2.00	0.43
1:A:232:LYS:O	1:A:236:ILE:HG13	2.18	0.43
1:B:357:ARG:NH2	1:B:460:GLU:OE2	2.50	0.43
1:A:323:HIS:HB3	1:A:327:TYR:CE2	2.54	0.43
1:B:336:LYS:HE2	1:B:372:VAL:HG11	2.00	0.43
1:A:325:ILE:HG13	1:A:392:ILE:HD11	1.99	0.43
1:A:426:PRO:HD2	1:A:427:GLU:OE1	2.18	0.42
1:A:276:GLU:OE2	1:A:357:ARG:HD3	2.19	0.42
1:B:320:TYR:CB	1:B:397:ARG:HD2	2.49	0.42
1:A:222:TYR:CE2	1:A:381:ASP:HB3	2.54	0.42
1:B:286:GLN:OE1	1:B:466:HIS:N	2.43	0.42
1:A:453:LEU:O	1:A:453:LEU:HG	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:LYS:HZ1	1:B:474:LYS:HB3	1.85	0.41
1:A:317:LEU:HD21	1:A:406:ILE:HD13	2.02	0.41
1:A:358:LYS:HB2	1:A:359:PRO:HD3	2.02	0.41
1:A:387:PHE:CZ	1:A:391:ILE:HD11	2.56	0.41
1:B:245:SER:HA	1:B:246:PRO:HD3	1.88	0.41
1:B:414:LEU:HD23	1:B:414:LEU:HA	1.84	0.41
1:A:396:ASP:HB2	1:A:443:ARG:HH22	1.87	0.40
1:A:379:LEU:HD11	1:A:435:LEU:HD11	2.02	0.40
1:B:301:LYS:HB3	1:B:301:LYS:HE2	1.84	0.40
1:A:320:TYR:CB	1:A:397:ARG:HD2	2.51	0.40
1:B:362:ASP:OD1	1:B:362:ASP:N	2.51	0.40
1:A:262:ILE:HG21	1:A:345:GLN:O	2.21	0.40
1:A:470:GLN:HA	1:A:473:TYR:HB2	2.03	0.40
2:D:646:MET:CA	2:D:649:LEU:HD12	2.37	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	255/297 (86%)	232 (91%)	18 (7%)	5 (2%)	7	31
1	B	255/297 (86%)	235 (92%)	18 (7%)	2 (1%)	19	54
2	C	10/19 (53%)	8 (80%)	2 (20%)	0	100	100
2	D	10/19 (53%)	7 (70%)	2 (20%)	1 (10%)	0	3
All	All	530/632 (84%)	482 (91%)	40 (8%)	8 (2%)	10	39

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	475	ASP

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Mol	Chain	Res	Type
1	B	475	ASP
1	A	261	LYS
1	A	260	ASP
1	A	262	ILE
1	A	259	GLU
2	D	642	HIS
1	B	310	ASP

### 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	232/268 (87%)	225 (97%)	7 (3%)	41	71
1	B	232/268 (87%)	230 (99%)	2 (1%)	78	91
2	C	12/18 (67%)	12 (100%)	0	100	100
2	D	12/18 (67%)	12 (100%)	0	100	100
All	All	488/572 (85%)	479 (98%)	9 (2%)	59	82

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	243	ASP
1	A	253	ASN
1	A	263	LYS
1	A	362	ASP
1	A	363	PHE
1	A	424	ASN
1	A	473	TYR
1	B	261	LYS
1	B	424	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	312	ASN
1	B	444	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	259/297 (87%)	-0.46	2 (0%) 86 72	30, 66, 113, 138	0
1	B	259/297 (87%)	-0.45	2 (0%) 86 72	37, 67, 111, 134	0
2	C	12/19 (63%)	-0.34	0 100 100	92, 100, 113, 124	0
2	D	12/19 (63%)	-0.32	1 (8%) 11 4	76, 84, 96, 122	0
All	All	542/632 (85%)	-0.45	5 (0%) 84 69	30, 68, 113, 138	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	260	ASP	3.2
1	B	261	LYS	2.9
1	A	239	GLY	2.9
1	B	260	ASP	2.8
2	D	651	ASP	2.3

### 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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.