



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 08:50 am GMT

PDB ID : 2PIZ
Title : CRYSTAL STRUCTURE OF ACTIVATED PORCINE PANCREATIC CARBOXYPEPTIDASE B 2-(3-Guanidino-phenyl)-3-[hydroxy-(3-phenyl-propyl)-phosphinoyl]-propionic acid COMPLEX
Authors : Adler, M.; Whitlow, M.
Deposited on : 2007-04-15
Resolution : 1.60 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

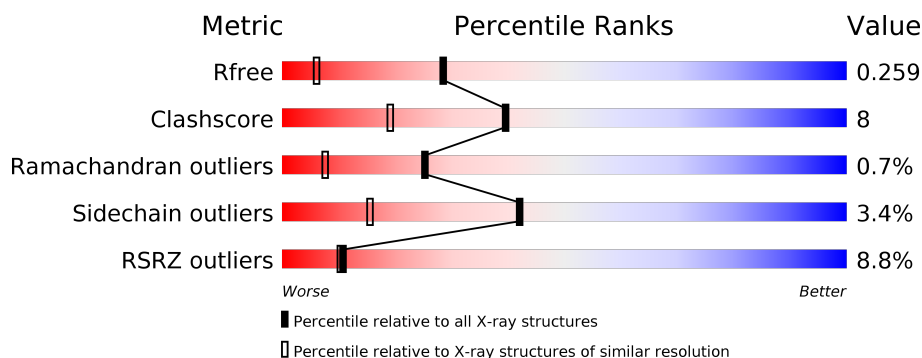
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	100719	2696 (1.60-1.60)
Clashscore	112137	2967 (1.60-1.60)
Ramachandran outliers	110173	2887 (1.60-1.60)
Sidechain outliers	110143	2886 (1.60-1.60)
RSRZ outliers	101464	2714 (1.60-1.60)

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

Mol	Chain	Length	Quality of chain
1	A	306	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 90%, yellow 8%, orange 1%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 90% 8% • </div> </div>
1	B	306	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 88%, yellow 10%, orange 1%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 88% 10% •• </div> </div>
1	C	306	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 24%, green 71%, yellow 26%, orange 1%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 24% 71% 26% • </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	606	B	501	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7973 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.

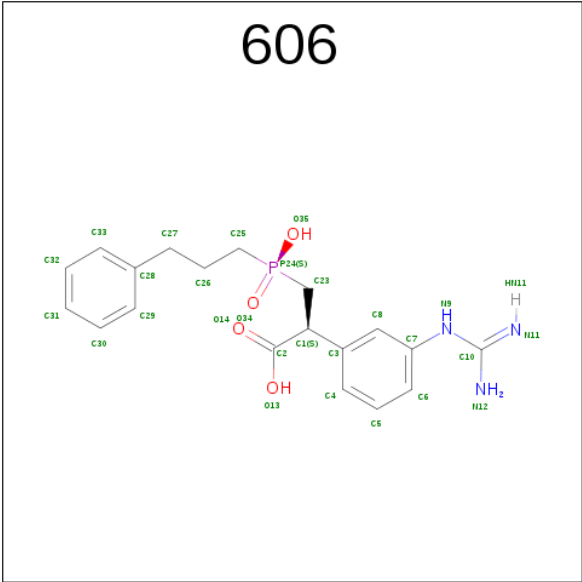
- Molecule 1 is a protein called Carboxypeptidase B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2443	1564	398	469	12			
1	B	304	Total	C	N	O	S	0	0	0
			2436	1560	397	467	12			
1	C	305	Total	C	N	O	S	140	0	0
			2443	1564	398	469	12			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is (2S)-2-(3-{[AMINO(IMINO)METHYL]AMINO}PHENYL)-3-[(S)-HYDROXY(3-PHENYLPROPYL)PHOSPHORYL]PROPANOIC ACID (three-letter code: 606) (formula: C₁₉H₂₄N₃O₄P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	19	3	4	1		
3	B	1	Total	C	N	O	P	0	0
			27	19	3	4	1		
3	C	1	Total	C	N	O	P	0	0
			27	19	3	4	1		

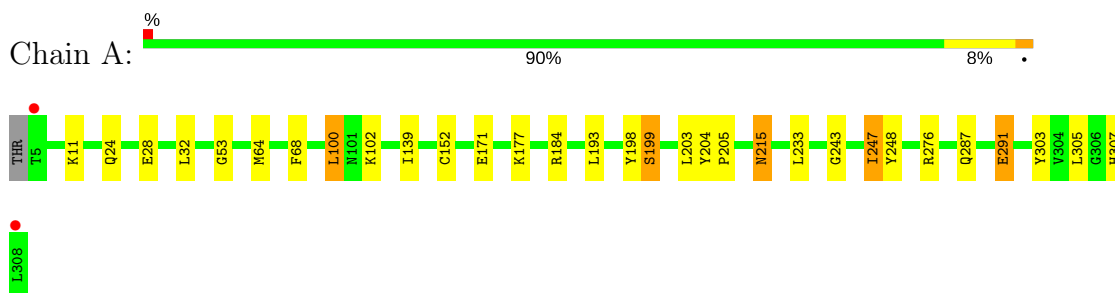
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	249	Total	O	0	0
			249	249		
4	B	223	Total	O	0	0
			223	223		
4	C	95	Total	O	0	0
			95	95		

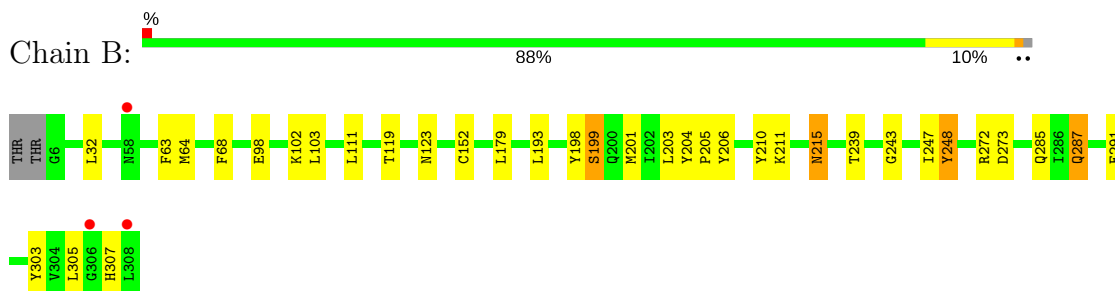
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

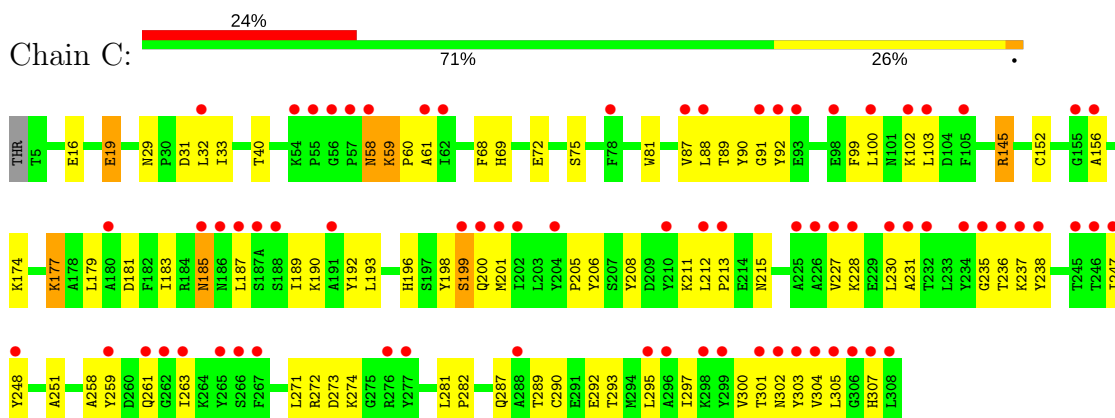
• Molecule 1: Carboxypeptidase B



• Molecule 1: Carboxypeptidase B



• Molecule 1: Carboxypeptidase B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.79Å 100.89Å 136.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 1.60 8.00 – 1.60	Depositor EDS
% Data completeness (in resolution range)	90.9 (8.00-1.60) 91.1 (8.00-1.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 1.60Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.232 , 0.264 0.225 , 0.259	Depositor DCC
R_{free} test set	4510 reflections (4.28%)	DCC
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.954	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 60.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7973	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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: ZN, 606

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	0.28	0/2512	0.47	0/3418
1	B	0.26	0/2505	0.46	0/3408
1	C	0.25	0/2512	0.45	0/3418
All	All	0.26	0/7529	0.46	0/10244

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

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	2443	0	2324	25	0
1	B	2436	0	2317	24	0
1	C	2443	0	2324	72	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	27	0	21	0	0
3	B	27	0	21	0	0
3	C	27	0	21	0	0
4	A	249	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	223	0	0	1	0
4	C	95	0	0	2	0
All	All	7973	0	7028	119	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 8.

All (119) 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:C:258:ALA:HA	1:C:261:GLN:OE1	1.76	0.84
1:C:145:ARG:CG	1:C:145:ARG:HH11	1.92	0.82
1:A:203:LEU:HD22	1:A:247:ILE:HD11	1.59	0.82
1:C:258:ALA:O	1:C:261:GLN:HG2	1.82	0.79
1:A:203:LEU:CD2	1:A:247:ILE:HD11	2.17	0.72
1:C:193:LEU:HD23	1:C:297:ILE:HD12	1.72	0.72
1:C:145:ARG:HG2	1:C:145:ARG:HH11	1.55	0.72
1:C:205:PRO:HB2	1:C:213:PRO:HG3	1.70	0.72
1:A:171:GLU:OE1	1:A:184:ARG:NH2	2.24	0.68
1:C:289:THR:HG23	4:C:651:HOH:O	1.93	0.67
1:C:192:TYR:C	1:C:193:LEU:HD12	2.15	0.66
1:C:247:ILE:HG23	1:C:248:TYR:H	1.59	0.66
1:C:200:GLN:HB2	1:C:274:LYS:NZ	2.10	0.65
1:A:215:ASN:HD22	1:A:215:ASN:H	1.41	0.65
1:B:215:ASN:HD22	1:B:215:ASN:H	1.46	0.62
1:B:272:ARG:HH11	1:B:285:GLN:HE21	1.48	0.62
1:C:183:ILE:HG21	1:C:261:GLN:HE22	1.64	0.62
1:C:99:PHE:O	1:C:103:LEU:HG	1.99	0.62
1:C:89:THR:O	1:C:91:GLY:N	2.33	0.62
1:C:227:VAL:O	1:C:231:ALA:HB2	2.01	0.61
1:C:200:GLN:HB2	1:C:274:LYS:HZ3	1.65	0.60
1:A:139:ILE:HD11	4:B:625:HOH:O	2.03	0.59
1:C:183:ILE:CG2	1:C:261:GLN:HE22	2.15	0.59
1:C:302:ASN:HA	1:C:305:LEU:HD12	1.85	0.58
1:C:213:PRO:HB3	1:C:259:TYR:CZ	2.38	0.57
1:C:208:TYR:HB3	1:C:251:ALA:HA	1.86	0.57
1:A:102:LYS:HD3	1:A:305:LEU:HD21	1.87	0.57
1:C:213:PRO:HB3	1:C:259:TYR:CE2	2.39	0.57
1:C:145:ARG:NH1	1:C:145:ARG:HG2	2.14	0.56
1:B:303:TYR:O	1:B:307:HIS:HD2	1.87	0.56
1:C:300:VAL:O	1:C:304:VAL:HG23	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:302:ASN:HA	1:C:305:LEU:CD1	2.36	0.56
1:C:88:LEU:O	1:C:92:TYR:CD1	2.59	0.56
1:C:200:GLN:HG3	1:C:273:ASP:O	2.05	0.56
1:C:81:TRP:CH2	1:C:287:GLN:HG3	2.41	0.56
1:A:276:ARG:HH11	1:A:276:ARG:HG3	1.71	0.55
1:A:247:ILE:N	1:A:247:ILE:HD12	2.21	0.55
1:A:233:LEU:HD21	1:A:291:GLU:HG2	1.88	0.55
1:C:200:GLN:O	1:C:238:TYR:HA	2.07	0.54
1:C:19:GLU:HG2	4:C:683:HOH:O	2.06	0.54
1:C:303:TYR:O	1:C:307:HIS:N	2.39	0.54
1:B:203:LEU:HD23	1:B:243:GLY:HA2	1.89	0.54
1:C:261:GLN:NE2	1:C:263:ILE:HD11	2.23	0.53
1:B:215:ASN:HD22	1:B:215:ASN:N	2.07	0.53
1:A:215:ASN:HD22	1:A:215:ASN:N	2.03	0.53
1:C:293:THR:O	1:C:297:ILE:HG12	2.08	0.53
1:C:145:ARG:HG3	1:C:145:ARG:HH11	1.72	0.52
1:C:201:MET:CE	1:C:247:ILE:HB	2.41	0.51
1:B:64:MET:HA	1:B:193:LEU:O	2.10	0.51
1:C:201:MET:HE3	1:C:247:ILE:HB	1.91	0.51
1:B:198:TYR:O	1:B:199:SER:CB	2.60	0.50
1:A:177:LYS:HD2	4:A:634:HOH:O	2.12	0.50
1:C:181:ASP:O	1:C:185:ASN:HB2	2.11	0.50
1:C:145:ARG:HD3	1:C:156:ALA:HB2	1.93	0.50
1:A:102:LYS:HE2	1:A:305:LEU:HD11	1.94	0.50
1:C:81:TRP:CD1	1:C:290:CYS:HB3	2.47	0.49
1:C:81:TRP:HH2	1:C:287:GLN:HG3	1.76	0.49
1:A:247:ILE:H	1:A:247:ILE:HD12	1.78	0.49
1:B:287:GLN:O	1:B:291:GLU:HG3	2.13	0.49
1:C:190:LYS:HB2	1:C:304:VAL:HG11	1.93	0.49
1:C:81:TRP:CG	1:C:290:CYS:HB3	2.48	0.49
1:C:198:TYR:O	1:C:199:SER:CB	2.61	0.48
1:A:198:TYR:O	1:A:199:SER:CB	2.61	0.48
1:C:230:LEU:HG	1:C:292:GLU:HB2	1.96	0.48
1:B:98:GLU:HG3	1:B:102:LYS:NZ	2.28	0.48
1:A:53:GLY:HA2	1:A:100:LEU:HG	1.96	0.48
1:C:31:ASP:C	1:C:32:LEU:HD12	2.35	0.47
1:C:177:LYS:C	1:C:177:LYS:HE2	2.34	0.47
1:B:272:ARG:HH11	1:B:285:GLN:NE2	2.12	0.47
1:C:301:THR:O	1:C:305:LEU:HG	2.14	0.47
1:C:69:HIS:CE1	1:C:196:HIS:HE1	2.32	0.47
1:A:11:LYS:HD3	4:A:621:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:LEU:O	1:C:92:TYR:HD1	1.97	0.47
1:C:193:LEU:HD12	1:C:193:LEU:N	2.30	0.47
1:C:29:ASN:OD1	1:C:87:VAL:HB	2.15	0.46
1:C:235:GLY:O	1:C:236:THR:C	2.54	0.46
1:B:203:LEU:HD21	1:B:247:ILE:CD1	2.46	0.46
1:C:238:TYR:OH	1:C:271:LEU:HA	2.15	0.46
1:B:204:TYR:HB2	1:B:205:PRO:CD	2.46	0.46
1:C:228:LYS:HA	1:C:231:ALA:HB3	1.98	0.46
1:A:24:GLN:O	1:A:28:GLU:HG3	2.17	0.45
1:C:58:ASN:O	1:C:59:LYS:HB2	2.17	0.45
1:A:203:LEU:HD13	1:A:247:ILE:HD11	1.98	0.45
1:B:203:LEU:HD21	1:B:247:ILE:HD12	1.98	0.45
1:A:203:LEU:CD1	1:A:247:ILE:HD11	2.47	0.45
1:B:63:PHE:CZ	1:B:179:LEU:HD11	2.52	0.45
1:C:190:LYS:HB2	1:C:304:VAL:CG1	2.47	0.45
1:A:204:TYR:HB2	1:A:205:PRO:CD	2.48	0.44
1:A:247:ILE:HB	1:A:248:TYR:H	1.65	0.44
1:A:64:MET:HA	1:A:193:LEU:O	2.17	0.44
1:B:103:LEU:C	1:B:103:LEU:HD12	2.38	0.44
1:B:111:LEU:HD13	1:B:111:LEU:C	2.38	0.44
1:A:215:ASN:ND2	1:A:215:ASN:H	2.11	0.44
1:C:61:ALA:O	1:C:189:ILE:HA	2.18	0.43
1:C:72:GLU:O	1:C:75:SER:HB3	2.18	0.43
1:C:179:LEU:HD23	1:C:179:LEU:C	2.38	0.43
1:C:272:ARG:HA	1:C:273:ASP:HA	1.78	0.43
1:C:69:HIS:CE1	1:C:196:HIS:CE1	3.07	0.43
1:B:247:ILE:HG22	1:B:248:TYR:CD2	2.54	0.43
1:B:210:TYR:CD2	1:C:102:LYS:HE3	2.53	0.43
1:A:303:TYR:O	1:A:307:HIS:HD2	2.01	0.43
1:C:59:LYS:HB3	1:C:59:LYS:HE2	1.87	0.43
1:B:272:ARG:HA	1:B:273:ASP:HA	1.83	0.43
1:C:237:LYS:HD3	1:C:237:LYS:HA	1.81	0.43
1:B:210:TYR:CG	1:C:102:LYS:HE3	2.55	0.42
1:B:203:LEU:N	1:B:203:LEU:HD12	2.35	0.42
1:C:281:LEU:HA	1:C:282:PRO:HD3	1.94	0.42
1:B:201:MET:HA	1:B:239:THR:O	2.19	0.42
1:C:33:ILE:HD11	1:C:87:VAL:HG11	2.00	0.42
1:B:205:PRO:HA	1:B:206:TYR:HA	1.78	0.41
1:C:211:LYS:HE2	1:C:211:LYS:HB3	1.70	0.41
1:C:60:PRO:HD2	1:C:102:LYS:O	2.21	0.41
1:C:205:PRO:HB2	1:C:213:PRO:CG	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:THR:HA	1:B:123:ASN:O	2.20	0.40
1:C:40:THR:O	1:C:174:LYS:HE2	2.21	0.40
1:C:100:LEU:N	1:C:100:LEU:HD12	2.37	0.40
1:C:205:PRO:HA	1:C:206:TYR:HA	1.72	0.40
1:C:212:LEU:HA	1:C:213:PRO:HD3	1.94	0.40
1:A:243:GLY:O	1:A:247:ILE:HD12	2.21	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	303/306 (99%)	295 (97%)	7 (2%)	1 (0%)	44	22
1	B	302/306 (99%)	290 (96%)	11 (4%)	1 (0%)	44	22
1	C	303/306 (99%)	285 (94%)	14 (5%)	4 (1%)	14	2
All	All	908/918 (99%)	870 (96%)	32 (4%)	6 (1%)	25	7

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	SER
1	B	199	SER
1	C	199	SER
1	C	90	TYR
1	C	59	LYS
1	C	58	ASN

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	258/259 (100%)	250 (97%)	8 (3%)	45	18
1	B	257/259 (99%)	249 (97%)	8 (3%)	45	18
1	C	258/259 (100%)	248 (96%)	10 (4%)	37	12
All	All	773/777 (100%)	747 (97%)	26 (3%)	42	15

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LEU
1	A	68	PHE
1	A	100	LEU
1	A	152	CYS
1	A	215	ASN
1	A	247	ILE
1	A	287	GLN
1	A	291	GLU
1	B	32	LEU
1	B	68	PHE
1	B	152	CYS
1	B	211	LYS
1	B	215	ASN
1	B	248	TYR
1	B	287	GLN
1	B	305	LEU
1	C	16	GLU
1	C	19	GLU
1	C	68	PHE
1	C	145	ARG
1	C	152	CYS
1	C	177	LYS
1	C	185	ASN
1	C	187	LEU
1	C	215	ASN
1	C	295	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	46	ASN
1	A	215	ASN
1	A	287	GLN
1	A	307	HIS
1	B	215	ASN
1	B	285	GLN
1	B	287	GLN
1	B	307	HIS
1	C	24	GLN
1	C	46	ASN
1	C	261	GLN
1	C	285	GLN
1	C	287	GLN
1	C	302	ASN

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](#)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 for Mogul analysis.

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
3	606	A	401	-	22,28,28	1.83	6 (27%)	25,38,38	0.77	0
3	606	B	501	-	22,28,28	1.79	5 (22%)	25,38,38	0.80	1 (4%)
3	606	C	601	-	22,28,28	1.89	5 (22%)	25,38,38	0.89	2 (8%)

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
3	606	A	401	-	-	0/20/24/24	0/2/2/2
3	606	B	501	-	-	0/20/24/24	0/2/2/2
3	606	C	601	-	-	0/20/24/24	0/2/2/2

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	601	606	C7-N9	-3.70	1.34	1.41
3	A	401	606	C7-N9	-3.46	1.34	1.41
3	B	501	606	C7-N9	-3.37	1.35	1.41
3	C	601	606	C10-N9	-2.38	1.34	1.37
3	A	401	606	C10-N9	-2.29	1.35	1.37
3	B	501	606	C10-N9	-2.15	1.35	1.37
3	A	401	606	C8-C3	2.09	1.42	1.39
3	B	501	606	C6-C7	2.16	1.42	1.39
3	A	401	606	P24-C23	2.40	1.81	1.79
3	B	501	606	P24-C23	2.47	1.82	1.79
3	C	601	606	C4-C3	2.49	1.43	1.39
3	A	401	606	P24-C25	2.61	1.82	1.79
3	A	401	606	C4-C3	2.61	1.43	1.39
3	B	501	606	C4-C3	2.74	1.43	1.39
3	C	601	606	P24-C25	2.92	1.82	1.79
3	C	601	606	P24-C23	3.28	1.82	1.79

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	601	606	C8-C3-C1	-2.63	117.33	120.81
3	B	501	606	C7-N9-C10	-2.32	122.47	128.03
3	C	601	606	C7-N9-C10	-2.14	122.92	128.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

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	305/306 (99%)	-0.28	2 (0%)  	16, 21, 28, 36	0
1	B	304/306 (99%)	-0.21	3 (0%)  	17, 23, 30, 39	0
1	C	286/306 (93%)	1.31	74 (25%)  	21, 32, 42, 45	0
All	All	895/918 (97%)	0.26	79 (8%)  	16, 24, 39, 45	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	303	TYR	10.3
1	C	306	GLY	9.8
1	C	225	ALA	8.7
1	C	308	LEU	7.7
1	C	304	VAL	5.2
1	C	204	TYR	5.0
1	C	247	ILE	5.0
1	C	305	LEU	4.9
1	C	299	TYR	4.8
1	C	246	THR	4.7
1	C	231	ALA	4.6
1	C	55	PRO	4.6
1	C	228	LYS	4.6
1	C	58	ASN	4.6
1	C	238	TYR	4.5
1	C	245	THR	4.4
1	C	265	TYR	4.3
1	C	191	ALA	4.1
1	C	188	SER	4.1
1	C	87	VAL	4.0
1	A	308	LEU	4.0
1	C	187	LEU	4.0
1	C	187(A)	SER	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	61	ALA	3.9
1	C	288	ALA	3.9
1	C	32	LEU	3.8
1	C	235	GLY	3.6
1	C	105	PHE	3.6
1	C	226	ALA	3.6
1	C	56	GLY	3.5
1	C	57	PRO	3.5
1	C	236	THR	3.5
1	C	92	TYR	3.4
1	C	261	GLN	3.3
1	C	103	LEU	3.3
1	C	100	LEU	3.3
1	C	302	ASN	3.2
1	C	267	PHE	3.2
1	C	248	TYR	3.1
1	C	102	LYS	3.1
1	B	58	ASN	3.0
1	C	93	GLU	2.9
1	C	185	ASN	2.9
1	C	200	GLN	2.8
1	C	155	GLY	2.8
1	C	180	ALA	2.8
1	C	276	ARG	2.7
1	C	88	LEU	2.7
1	C	237	LYS	2.7
1	C	262	GLY	2.7
1	C	232	THR	2.7
1	C	266	SER	2.6
1	C	298	LYS	2.6
1	C	234	TYR	2.6
1	C	227	VAL	2.6
1	C	210	TYR	2.6
1	C	62	ILE	2.5
1	C	201	MET	2.5
1	C	230	LEU	2.5
1	C	54	LYS	2.5
1	C	295	LEU	2.5
1	C	213	PRO	2.4
1	C	98	GLU	2.4
1	C	78	PHE	2.4
1	C	212	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	263	ILE	2.4
1	C	296	ALA	2.4
1	C	202	ILE	2.4
1	C	186	ASN	2.4
1	A	5	THR	2.3
1	C	277	TYR	2.3
1	C	307	HIS	2.3
1	C	259	TYR	2.2
1	C	91	GLY	2.2
1	B	308	LEU	2.2
1	C	156	ALA	2.2
1	C	199	SER	2.1
1	B	306	GLY	2.1
1	C	301	THR	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	606	B	501	27/27	0.91	0.13	2.26	22,28,36,36	0
3	606	A	401	27/27	0.92	0.11	1.27	17,25,33,34	0
3	606	C	601	27/27	0.86	0.14	-0.03	27,30,40,41	0
2	ZN	C	600	1/1	0.99	0.05	-1.79	26,26,26,26	0
2	ZN	B	500	1/1	1.00	0.02	-3.54	19,19,19,19	0
2	ZN	A	400	1/1	1.00	0.02	-4.44	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.