



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:29 PM GMT

PDB ID : 4II4
Title : The Phenylacetyl-CoA monooxygenase - mutant PaaA E49Q K68Q - PaaC wild type subcomplex with benzoyl-CoA
Authors : Cygler, M.; Grishin, A.M.; Montreal-Kingston Bacterial Structural Genomics Initiative (BSGI)
Deposited on : 2012-12-19
Resolution : 2.80 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 (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

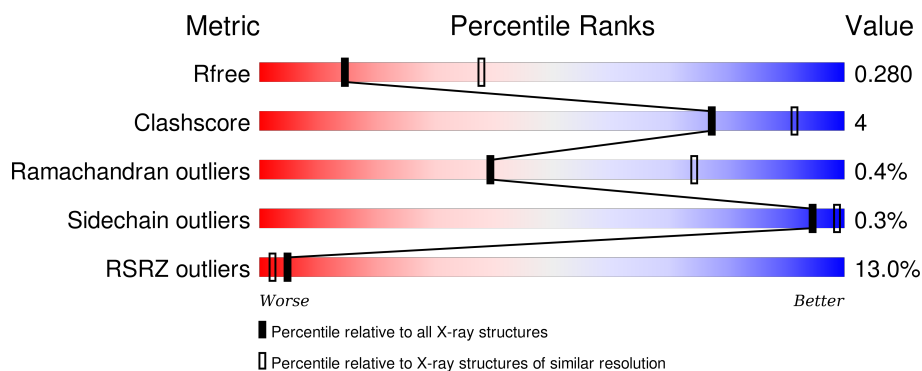
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	311	<div> <div>6%</div> <div>93%</div> <div>5%</div> </div>
2	B	259	<div> <div>20%</div> <div>81%</div> <div>10%</div> <div>8%</div> </div>
2	C	259	<div> <div>12%</div> <div>84%</div> <div>12%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6266 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 1,2-phenylacetyl-CoA epoxidase, subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2403	1511	429	447	16			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	EXPRESSION TAG	UNP P76077
A	0	ARG	-	EXPRESSION TAG	UNP P76077
A	1	SER	-	EXPRESSION TAG	UNP P76077
A	49	GLN	GLU	ENGINEERED MUTATION	UNP P76077
A	68	GLN	LYS	ENGINEERED MUTATION	UNP P76077

- Molecule 2 is a protein called 1,2-phenylacetyl-CoA epoxidase, subunit C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	239	Total	C	N	O	S	0	0	0
			1841	1160	328	347	6			
2	C	248	Total	C	N	O	S	0	0	0
			1951	1229	345	371	6			

There are 24 discrepancies between the modelled and reference sequences:

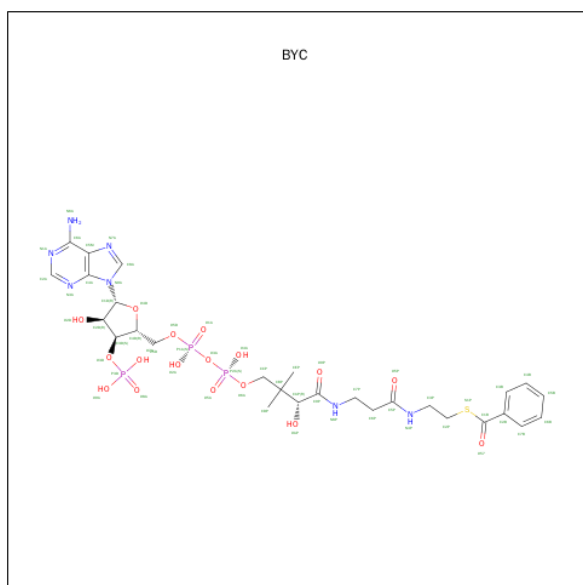
Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	MET	-	EXPRESSION TAG	UNP P76079
B	-9	GLY	-	EXPRESSION TAG	UNP P76079
B	-8	SER	-	EXPRESSION TAG	UNP P76079
B	-7	SER	-	EXPRESSION TAG	UNP P76079
B	-6	HIS	-	EXPRESSION TAG	UNP P76079
B	-5	HIS	-	EXPRESSION TAG	UNP P76079
B	-4	HIS	-	EXPRESSION TAG	UNP P76079
B	-3	HIS	-	EXPRESSION TAG	UNP P76079
B	-2	HIS	-	EXPRESSION TAG	UNP P76079

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	HIS	-	EXPRESSION TAG	UNP P76079
B	0	GLY	-	EXPRESSION TAG	UNP P76079
B	1	SER	-	EXPRESSION TAG	UNP P76079
C	-10	MET	-	EXPRESSION TAG	UNP P76079
C	-9	GLY	-	EXPRESSION TAG	UNP P76079
C	-8	SER	-	EXPRESSION TAG	UNP P76079
C	-7	SER	-	EXPRESSION TAG	UNP P76079
C	-6	HIS	-	EXPRESSION TAG	UNP P76079
C	-5	HIS	-	EXPRESSION TAG	UNP P76079
C	-4	HIS	-	EXPRESSION TAG	UNP P76079
C	-3	HIS	-	EXPRESSION TAG	UNP P76079
C	-2	HIS	-	EXPRESSION TAG	UNP P76079
C	-1	HIS	-	EXPRESSION TAG	UNP P76079
C	0	GLY	-	EXPRESSION TAG	UNP P76079
C	1	SER	-	EXPRESSION TAG	UNP P76079

- Molecule 3 is BENZOYL COENZYME A (three-letter code: BYC) (formula: $C_{28}H_{40}N_7O_{17}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			46	20	6	16	3	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	4	Total 4	O 4	0	0
4	C	12	Total 12	O 12	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	77.59Å 77.59Å 304.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.64 – 2.80 48.27 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (26.64-2.80) 91.3 (48.27-2.80)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.218 , 0.276 0.233 , 0.280	Depositor DCC
R_{free} test set	1225 reflections (5.89%)	DCC
Wilson B-factor (Å ²)	75.4	Xtriage
Anisotropy	0.102	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 50.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 23945 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6266	wwPDB-VP
Average B, all atoms (Å ²)	119.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.10% 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, 0.375 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: BYC

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	0.38	0/2460	0.59	0/3336
2	B	0.39	0/1877	0.64	1/2547 (0.0%)
2	C	0.37	0/1990	0.61	0/2697
All	All	0.38	0/6327	0.61	1/8580 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	149	ASP	CB-CG-OD1	5.69	123.42	118.30

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	2403	0	2285	10	0
2	B	1841	0	1760	20	0
2	C	1951	0	1890	16	0
3	A	46	0	25	1	0
4	A	9	0	0	1	0
4	B	4	0	0	0	0
4	C	12	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6266	0	5960	45	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 4.

All (45) 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:B:149:ASP:OD1	2:B:150:VAL:N	2.11	0.84
2:C:156:GLN:HE22	2:C:209:ASN:H	1.24	0.84
2:C:10:ARG:NH2	2:C:173:ASP:OD2	2.23	0.72
2:B:80:LEU:HD22	2:B:169:LEU:HD12	1.76	0.68
2:C:148:THR:HG23	2:C:151:SER:H	1.61	0.66
2:C:111:LEU:HD13	2:C:186:VAL:HG23	1.79	0.64
2:C:127:ILE:O	2:C:131:ARG:HG3	1.99	0.62
2:B:10:ARG:NH1	2:B:173:ASP:OD2	2.27	0.61
1:A:30:THR:HG23	1:A:33:ARG:HE	1.64	0.60
1:A:91:GLU:H	1:A:91:GLU:CD	2.04	0.59
1:A:0:ARG:NH2	1:A:5:GLU:OE1	2.27	0.59
2:B:6:ALA:O	2:B:10:ARG:HG3	2.04	0.58
2:B:80:LEU:O	2:B:83:VAL:HG22	2.06	0.55
2:B:161:LYS:HG3	2:B:162:LEU:HG	1.87	0.55
2:C:81:LEU:O	2:C:84:GLU:HG2	2.07	0.55
2:C:48:GLN:NE2	2:C:129:GLU:OE1	2.38	0.54
2:B:20:GLN:HG3	2:B:236:MET:SD	2.49	0.53
2:B:233:LEU:O	2:B:237:GLN:HG3	2.09	0.53
2:C:25:TRP:HE1	2:C:93:THR:HG1	1.57	0.52
1:A:185:ASN:HA	1:A:241:PRO:HG3	1.92	0.51
1:A:17:ILE:HD12	1:A:27:TYR:OH	2.10	0.51
2:B:116:ASP:HB3	2:B:119:LEU:HB2	1.92	0.50
1:A:123:TRP:HH2	1:A:233:VAL:HG22	1.77	0.48
2:B:18:LEU:HD21	2:B:100:ILE:HG21	1.96	0.47
1:A:6:ARG:NH1	4:A:502:HOH:O	2.48	0.47
2:B:67:THR:O	2:B:71:THR:OG1	2.18	0.46
1:A:62:LYS:HE2	2:C:20:GLN:OE1	2.15	0.46
2:B:236:MET:O	2:B:237:GLN:C	2.53	0.46
3:A:401:BYC:O3A	3:A:401:BYC:H2D	2.16	0.45
2:B:160:ASN:OD1	2:B:211:PRO:HG3	2.17	0.44
2:C:18:LEU:HB2	2:C:83:VAL:HB	1.98	0.44
1:A:59:LEU:HD12	2:C:65:GLU:HB2	1.99	0.44
2:B:18:LEU:CD2	2:B:100:ILE:HG21	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:236:MET:O	2:B:238:TYR:N	2.52	0.43
2:B:111:LEU:HD13	2:B:186:VAL:HG23	2.01	0.43
2:B:236:MET:HE3	2:B:236:MET:HB3	1.92	0.42
1:A:90:ARG:HE	1:A:90:ARG:HB2	1.57	0.42
2:B:153:GLN:O	2:B:157:GLN:N	2.39	0.42
2:C:37:ALA:O	2:C:41:ILE:HG13	2.20	0.41
2:C:110:ARG:HH22	2:C:190:THR:CB	2.33	0.41
2:B:31:GLU:HB2	2:B:34:ILE:HD12	2.02	0.41
2:C:153:GLN:O	2:C:157:GLN:HG2	2.20	0.41
2:B:110:ARG:HA	2:B:110:ARG:HD3	1.83	0.41
2:C:97:GLN:HG3	2:C:101:ASP:OD2	2.21	0.40
2:C:103:TRP:HE3	2:C:104:HIS:HD2	1.69	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	302/311 (97%)	300 (99%)	2 (1%)	0	100	100
2	B	237/259 (92%)	229 (97%)	5 (2%)	3 (1%)	15	44
2	C	246/259 (95%)	241 (98%)	5 (2%)	0	100	100
All	All	785/829 (95%)	770 (98%)	12 (2%)	3 (0%)	39	74

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	149	ASP
2	B	237	GLN
2	B	236	MET

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	242/259 (93%)	241 (100%)	1 (0%)	93	98
2	B	178/208 (86%)	177 (99%)	1 (1%)	90	98
2	C	196/208 (94%)	196 (100%)	0	100	100
All	All	616/675 (91%)	614 (100%)	2 (0%)	94	99

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	265	ASN
2	B	110	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	GLN
2	C	104	HIS
2	C	156	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 ⓘ

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$
3	BYC	A	401	-	36,47,59	2.37	10 (27%)	45,70,87	2.27	8 (17%)

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	BYC	A	401	-	-	0/38/59/71	0/3/3/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	BYC	O4D-C1D	-6.64	1.32	1.41
3	A	401	BYC	CEP-CBP	-2.99	1.47	1.53
3	A	401	BYC	O3D-C3D	-2.91	1.35	1.44
3	A	401	BYC	O2D-C2D	-2.82	1.36	1.43
3	A	401	BYC	P1A-O2A	-2.11	1.45	1.54
3	A	401	BYC	O5P-C5P	2.62	1.36	1.19
3	A	401	BYC	C7P-N8P	3.32	1.53	1.46
3	A	401	BYC	O9P-C9P	5.45	1.34	1.23
3	A	401	BYC	O4D-C4D	5.50	1.57	1.45
3	A	401	BYC	C7P-C6P	5.65	1.59	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	BYC	N3A-C2A-N1A	-9.02	121.99	128.89
3	A	401	BYC	C4D-O4D-C1D	-5.59	103.58	109.72
3	A	401	BYC	P2A-O3A-P1A	-4.95	118.83	132.73
3	A	401	BYC	OAP-CAP-C9P	-2.02	105.76	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	BYC	O4D-C4D-C5D	2.03	116.57	109.32
3	A	401	BYC	C2D-C1D-N9A	2.17	117.60	114.29
3	A	401	BYC	O4D-C1D-N9A	4.21	116.91	108.10
3	A	401	BYC	O9P-C9P-N8P	5.26	133.63	123.08

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	BYC	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	304/311 (97%)	0.60	20 (6%) 22 13	56, 110, 168, 209	0
2	B	239/259 (92%)	1.02	51 (21%) 1 1	86, 135, 179, 198	0
2	C	248/259 (95%)	0.75	32 (12%) 5 2	58, 109, 158, 189	0
All	All	791/829 (95%)	0.77	103 (13%) 5 2	56, 118, 173, 209	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-1	MET	7.3
2	C	206	ALA	6.7
2	B	144	LEU	6.6
2	B	184	ILE	6.4
1	A	295	ALA	6.0
2	C	146	ASN	5.8
2	B	179	LEU	5.4
2	B	175	ILE	5.3
2	B	186	VAL	5.3
2	B	124	ALA	4.4
2	B	58	GLU	4.4
2	B	128	LYS	4.3
2	B	185	ALA	4.3
2	C	147	GLY	4.3
2	B	176	ASP	4.1
2	C	141	LEU	4.1
2	B	98	TYR	3.9
2	B	112	MET	3.9
2	B	151	SER	3.7
2	B	150	VAL	3.6
1	A	279	LEU	3.6
2	C	145	GLY	3.6
1	A	22	TRP	3.5

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Mol	Chain	Res	Type	RSRZ
2	B	25	TRP	3.5
2	B	145	GLY	3.3
1	A	17	ILE	3.3
2	B	53	LEU	3.2
2	C	184	ILE	3.2
1	A	211	TRP	3.2
2	C	179	LEU	3.1
2	B	108	PHE	3.1
2	B	119	LEU	3.1
2	B	183	GLY	3.1
2	C	207	THR	3.1
2	B	148	THR	3.0
2	B	68	LEU	3.0
2	B	153	GLN	3.0
2	B	211	PRO	3.0
2	B	121	ALA	3.0
2	C	4	LEU	3.0
2	B	147	GLY	3.0
2	B	146	ASN	2.9
2	B	159	ILE	2.8
2	B	56	ALA	2.7
2	B	149	ASP	2.7
2	B	217	ARG	2.7
2	B	154	LYS	2.7
2	B	152	GLY	2.7
2	B	239	LEU	2.7
2	B	52	PHE	2.7
1	A	131	VAL	2.7
1	A	292	VAL	2.7
2	C	40	ASN	2.6
1	A	104	TYR	2.6
2	B	99	PHE	2.6
1	A	79	LEU	2.6
2	B	191	LEU	2.6
1	A	291	TRP	2.5
2	B	155	MET	2.5
2	C	39	ALA	2.5
2	C	149	ASP	2.5
2	B	210	VAL	2.5
2	B	195	TRP	2.5
1	A	97	MET	2.4
2	C	25	TRP	2.4

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Mol	Chain	Res	Type	RSRZ
2	C	16	LEU	2.4
2	B	4	LEU	2.4
2	B	80	LEU	2.4
1	A	82	ALA	2.4
2	C	82	LEU	2.4
2	C	37	ALA	2.4
2	C	19	SER	2.4
2	C	183	GLY	2.3
2	B	122	ILE	2.3
2	B	178	ALA	2.3
2	C	150	VAL	2.3
1	A	294	GLU	2.2
2	C	144	LEU	2.2
2	B	165	PHE	2.2
2	C	154	LYS	2.2
2	B	141	LEU	2.2
2	B	30	PRO	2.2
2	C	119	LEU	2.2
2	C	153	GLN	2.2
2	C	208	LEU	2.2
2	C	155	MET	2.2
2	C	122	ILE	2.1
1	A	98	LEU	2.1
2	C	42	GLY	2.1
2	C	30	PRO	2.1
2	B	81	LEU	2.1
1	A	290	THR	2.1
2	C	112	MET	2.1
1	A	32	ILE	2.1
1	A	225	PHE	2.1
1	A	93	ILE	2.0
2	C	98	TYR	2.0
2	C	215	ALA	2.0
2	C	41	ILE	2.0
1	A	9	GLN	2.0
2	B	157	GLN	2.0
2	B	114	SER	2.0
2	B	55	TYR	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 [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(\AA^2)	Q<0.9
3	BYC	A	401	46/56	0.95	0.16	-0.81	110,140,159,167	0

6.5 Other polymers [i](#)

There are no such residues in this entry.