



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:52 PM GMT

PDB ID : 2OOL
Title : Crystal structure of the chromophore-binding domain of an unusual bacterio-
phytochrome RpBphP3 from *R. palustris*
Authors : Yang, X.; Stojkovic, E.A.; Kuk, J.; Moffat, K.
Deposited on : 2007-01-25
Resolution : 2.20 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>

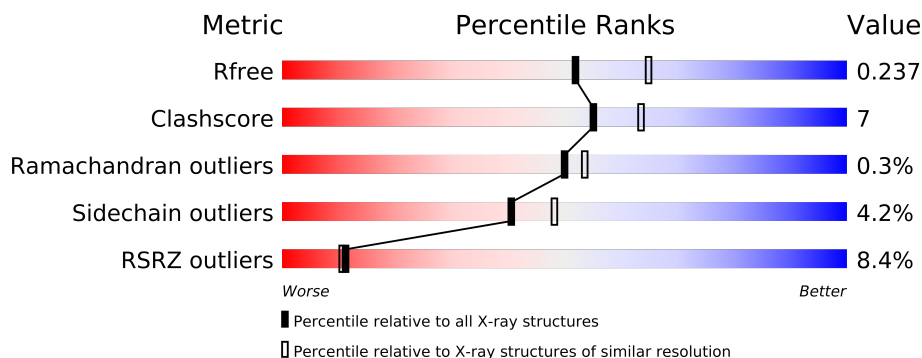
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.20 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	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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	337	
1	B	337	

2 Entry composition i

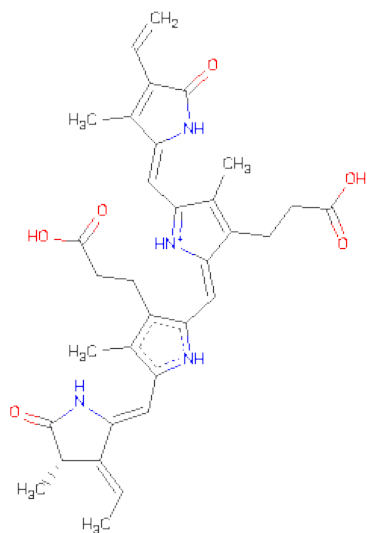
There are 3 unique types of molecules in this entry. The entry contains 5101 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Sensor protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	15	5	0
			2431	1542	435	445	9			
1	B	298	Total	C	N	O	S	5	3	0
			2368	1501	428	430	9			

- Molecule 2 is 3-[2-[(Z)-[3-(2-CARBOXYETHYL)-5-[(Z)-(4-ETHENYL-3-METHYL-5-OXIDANYLIDENE-PYRROL-2-YLIDENE)METHYL]-4-METHYL-PYRROL-1-IUM-2-YLIDENE]METHYL]-5-[(Z)-[(3E)-3-ETHYLIDENE-4-METHYL-5-OXIDANYLIDENE-PYRROLIDIN-2-YLIDENE]METHYL]-4-METHYL-1H-PYRROL-3-YL]PROPANOICACID (three-letter code: LBV) (formula: C₃₃H₃₇N₄O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	1
			44	34	4	6		
2	B	1	Total	C	N	O	0	0
			43	33	4	6		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	161	Total 161	O 161	0	0
3	B	54	Total 54	O 54	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	151.87Å 151.87Å 76.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.75 – 2.20 49.73 – 2.11	Depositor EDS
% Data completeness (in resolution range)	93.3 (49.75-2.20) 91.0 (49.73-2.11)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.188 , 0.231 0.202 , 0.237	Depositor DCC
R_{free} test set	4055 reflections (7.62%)	DCC
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.4	EDS
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 58049 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5101	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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: LBV

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	0.84	3/2492 (0.1%)	0.85	3/3402 (0.1%)
1	B	0.59	0/2428	0.68	0/3311
All	All	0.73	3/4920 (0.1%)	0.77	3/6713 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	144[A]	GLU	CB-CG	6.38	1.64	1.52
1	A	144[B]	GLU	CB-CG	6.38	1.64	1.52
1	A	28	CYS	CB-SG	-5.78	1.72	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	ARG	NE-CZ-NH2	-8.35	116.13	120.30
1	A	106	ARG	NE-CZ-NH1	7.34	123.97	120.30
1	A	222	ARG	NE-CZ-NH2	-5.20	117.70	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	143	ASN	Peptide

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	2431	0	2416	31	1
1	B	2368	0	2351	31	0
2	A	44	0	8	2	0
2	B	43	0	33	2	0
3	A	161	0	0	3	0
3	B	54	0	0	2	0
All	All	5101	0	4808	64	1

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 7.

All (64) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:73:LEU:O	1:B:76:VAL:HG12	1.63	0.98
1:A:51:THR:HG23	1:A:52:ASP:H	1.43	0.82
1:A:215:SER:O	2:A:400[B]:LBV:HMA3	1.86	0.76
1:A:51:THR:HG23	1:A:52:ASP:OD1	1.88	0.72
1:B:236:ILE:HD11	1:B:278:MET:O	1.88	0.72
1:B:103:ASN:HD21	1:B:122:LEU:H	1.42	0.67
1:B:236:ILE:HD11	1:B:273:MET:HB3	1.75	0.66
1:B:204:ILE:HD12	1:B:301:LEU:HD21	1.77	0.65
1:B:123[A]:HIS:CD2	3:B:402:HOH:O	2.49	0.65
1:B:250:ARG:HB2	1:B:251:LEU:HD13	1.80	0.63
1:B:236:ILE:CD1	1:B:278:MET:O	2.48	0.61
1:A:102:ILE:HG12	1:A:103:ASN:N	2.19	0.57
1:A:181:ARG:HE	1:A:183:LYS:NZ	2.04	0.56
1:B:204:ILE:HD12	1:B:301:LEU:CD2	2.35	0.56
1:B:123[A]:HIS:CE1	1:B:259:PHE:O	2.59	0.55
1:B:103:ASN:ND2	1:B:122:LEU:H	2.03	0.55
1:B:48:VAL:HG23	1:B:129:VAL:HB	1.88	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:181:ARG:HE	1:A:183:LYS:HZ3	1.53	0.54
1:B:245:PRO:HB2	1:B:247:ILE:HD12	1.89	0.54
1:A:121:ILE:HD12	1:A:308:TYR:CD1	2.44	0.53
1:A:26:THR:O	1:A:26:THR:HG22	2.11	0.51
1:B:123[A]:HIS:HD2	3:B:402:HOH:O	1.87	0.51
1:A:156:LEU:HD13	1:A:168:ILE:CG2	2.40	0.51
1:A:64:ASP:HB2	3:A:491:HOH:O	2.09	0.51
1:A:73:LEU:CD2	1:A:81:TYR:OH	2.61	0.49
1:B:183:LYS:HZ1	1:B:299:HIS:CE1	2.31	0.49
1:A:49:SER:O	1:A:53:LEU:HA	2.13	0.48
1:A:153:ILE:O	1:A:157:GLN:HG3	2.14	0.47
1:B:306:VAL:HB	1:B:310:VAL:HG21	1.95	0.47
2:B:400:LBV:N_D	2:B:400:LBV:HMC1	2.30	0.47
1:A:142:THR:O	1:A:145:PHE:HB3	2.15	0.47
1:A:106:ARG:HG3	1:A:136:ARG:NH1	2.30	0.46
1:B:153:ILE:HD11	1:B:316:LEU:HG	1.96	0.46
1:B:49:SER:O	1:B:53:LEU:HD23	2.15	0.46
1:A:50:GLU:OE1	1:A:124:ARG:NE	2.49	0.46
1:A:49:SER:O	1:A:53:LEU:HD23	2.16	0.45
1:A:47:VAL:HG21	1:A:255:ILE:HD13	1.98	0.45
1:B:218:PRO:O	1:B:221:SER:HB2	2.16	0.45
1:A:147:ARG:O	1:A:151:VAL:HG13	2.16	0.45
1:B:88:ALA:O	1:B:91:THR:HG22	2.17	0.45
1:B:47:VAL:HG13	1:B:128:ILE:HD12	1.99	0.45
1:B:230:VAL:HG13	1:B:284:ILE:HD13	1.98	0.45
1:A:128:ILE:HG21	1:A:249:PRO:HD2	1.98	0.45
1:A:192:SER:HB2	1:A:211[B]:HIS:CE1	2.53	0.43
1:A:121:ILE:HD12	1:A:308:TYR:CE1	2.53	0.43
1:A:268:THR:HG22	2:A:400[A]:LBV:HMA2	2.01	0.43
1:A:75:ASN:HD22	1:A:75:ASN:HA	1.66	0.43
2:B:400:LBV:O_D	2:B:400:LBV:HBD1	2.18	0.43
1:B:306:VAL:HB	1:B:310:VAL:CG2	2.48	0.42
1:A:151:VAL:HG22	3:A:515:HOH:O	2.19	0.42
1:A:323:TRP:O	1:A:327:VAL:HG23	2.20	0.42
1:B:168:ILE:O	1:B:172:GLU:HG2	2.19	0.42
1:A:279:HIS:HE1	3:A:499:HOH:O	2.03	0.41
1:B:119:ASN:HA	1:B:119:ASN:HD22	1.67	0.41
1:B:162:LEU:HB3	1:B:163:PRO:HD3	2.01	0.41
1:B:236:ILE:HG21	1:B:274:VAL:HG22	2.03	0.41
1:B:200:ARG:NE	1:B:204:ILE:O	2.54	0.41
1:B:76:VAL:HG11	1:B:81:TYR:HE1	1.85	0.41
1:A:269:HIS:O	1:A:272:TYR:HB3	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:156:LEU:HD13	1:A:168:ILE:HG22	2.03	0.41
1:A:53:LEU:HD23	1:A:53:LEU:HA	1.85	0.41
1:A:162:LEU:HB3	1:A:163:PRO:HD3	2.02	0.41
1:B:172:GLU:O	1:B:176:ILE:HG23	2.20	0.41
1:B:90:LEU:O	1:B:94:LEU:HD12	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:106:ARG:NH2	1:A:106:ARG:NH2[5.556]	1.99	0.21

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	305/337 (90%)	298 (98%)	5 (2%)	2 (1%)	30	28
1	B	295/337 (88%)	288 (98%)	7 (2%)	0	100	100
All	All	600/674 (89%)	586 (98%)	12 (2%)	2 (0%)	50	53

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	51	THR
1	A	95	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	267/288 (93%)	256 (96%)	11 (4%)	41	49
1	B	259/288 (90%)	248 (96%)	11 (4%)	40	48
All	All	526/576 (91%)	504 (96%)	22 (4%)	40	48

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	ASP
1	A	58	VAL
1	A	75	ASN
1	A	102	ILE
1	A	119	ASN
1	A	127	SER
1	A	151	VAL
1	A	156	LEU
1	A	231	ARG
1	A	260	SER
1	A	276	MET
1	B	29	ASP
1	B	50	GLU
1	B	91	THR
1	B	102	ILE
1	B	200	ARG
1	B	202	SER
1	B	207	LEU
1	B	208	LEU
1	B	227	ILE
1	B	231	ARG
1	B	251	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	ASN
1	A	119	ASN
1	A	186	GLN
1	B	75	ASN
1	B	103	ASN
1	B	119	ASN

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 ⓘ

Of 3 ligands modelled in this entry, 2 are modelled with single atom - leaving 1 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	$\# Z > 2$
2	LBV	B	400	1	46,46,46	1.85	11 (23%)	65,67,67	1.83	14 (21%)

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	LBV	B	400	1	1/1/14/22	2/26/74/74	0/4/4/4

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	400	LBV	CAA-C3A	5.79	1.40	1.33
2	B	400	LBV	CHC-C1C	4.54	1.40	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	400	LBV	CHB-C4A	3.54	1.42	1.34
2	B	400	LBV	C3D-C4D	-2.92	1.38	1.47
2	B	400	LBV	C1C-N_C	-2.42	1.34	1.38
2	B	400	LBV	CBC-CGC	2.40	1.56	1.50
2	B	400	LBV	CAD-C3D	-2.36	1.39	1.48
2	B	400	LBV	C4C-N_C	-2.35	1.33	1.37
2	B	400	LBV	C1D-C2D	-2.31	1.40	1.45
2	B	400	LBV	CBD-CAD	2.07	1.40	1.29
2	B	400	LBV	C1D-N_D	-2.05	1.34	1.37

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	LBV	C2A-C1A-N_A	5.14	112.22	107.82
2	B	400	LBV	CHB-C4A-C3A	-3.97	119.90	127.15
2	B	400	LBV	C1B-C2B-C3B	3.85	109.66	107.04
2	B	400	LBV	O_A-C1A-C2A	-3.78	122.72	126.60
2	B	400	LBV	CHD-C1D-N_D	3.24	133.22	126.19
2	B	400	LBV	CHD-C1D-C2D	-3.23	119.82	126.78
2	B	400	LBV	C4B-CHC-C1C	-3.20	124.60	128.71
2	B	400	LBV	C3D-C4D-N_D	2.96	109.33	106.66
2	B	400	LBV	CAC-C2C-C1C	2.42	128.28	125.52
2	B	400	LBV	CAB-C3B-C4B	2.39	129.85	124.73
2	B	400	LBV	CMC-C3C-C4C	2.39	127.94	124.87
2	B	400	LBV	CMD-C2D-C1D	2.38	127.48	124.23
2	B	400	LBV	C2A-C3A-CAA	-2.24	124.21	129.36
2	B	400	LBV	C2A-C3A-C4A	2.13	110.86	107.89

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	400	LBV	C2A

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	400	LBV	C1D-CHD-C4C-N_C
2	B	400	LBV	C1D-CHD-C4C-C3C

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	304/337 (90%)	0.61	15 (4%) 28 28	49, 56, 69, 78	2 (0%)
1	B	298/337 (88%)	0.68	36 (12%) 5 4	49, 59, 71, 87	1 (0%)
All	All	602/674 (89%)	0.64	51 (8%) 11 10	49, 57, 71, 87	3 (0%)

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	67	ARG	5.9
1	B	210	PHE	5.7
1	B	112	PRO	5.5
1	B	162	LEU	5.2
1	A	333	ILE	5.1
1	B	114	GLY	4.7
1	B	211	HIS	4.1
1	B	95	HIS	3.7
1	B	239	ARG	3.7
1	B	65	LEU	3.7
1	B	81	TYR	3.6
1	B	113	ASP	3.3
1	B	165	ALA	3.2
1	B	82	LEU	3.1
1	A	261	VAL	3.0
1	A	230	VAL	2.9
1	B	196	ILE	2.9
1	B	110	VAL	2.9
1	B	200	ARG	2.8
1	B	236	ILE	2.8
1	B	87	ALA	2.8
1	B	28	CYS	2.8
1	A	262	LEU	2.7
1	B	330	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	44	TYR	2.7
1	B	194	GLN	2.6
1	A	101	ALA	2.6
1	A	281	ALA	2.6
1	A	224	LEU	2.6
1	B	136	ARG	2.6
1	B	69	PRO	2.5
1	B	191	TRP	2.5
1	B	64	ASP	2.5
1	A	314	CYS	2.4
1	B	206	SER	2.4
1	B	203	GLY	2.4
1	B	66	LEU	2.4
1	A	282	MET	2.4
1	B	185	TYR	2.3
1	A	331	ALA	2.3
1	B	202	SER	2.3
1	B	109	VAL	2.2
1	B	212	PHE	2.2
1	B	116	ARG	2.1
1	A	105[A]	ILE	2.1
1	A	233	ILE	2.1
1	B	167	TRP	2.1
1	A	305	PHE	2.1
1	B	74	LEU	2.0
1	A	298	CYS	2.0
1	A	284	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	LBV	B	400	43/43	0.16	0.13	61,75,81,82	0
2	LBV	A	400[B]	1/43	0.12	-0.46	54,54,54,54	1
2	LBV	A	400[A]	1/43	0.12	-2.56	56,56,56,56	1

6.5 Other polymers ⓘ

There are no such residues in this entry.