



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 13, 2023 – 05:52 am GMT

PDB ID : 2X4D
Title : Crystal structure of human phospholysine phosphohistidine inorganic pyrophosphate phosphatase LHPP
Authors : Vollmar, M.; Gileadi, C.; Guo, K.; Savitsky, P.; Muniz, J.R.C.; Yue, W.; Allerton, C.; von Delft, F.; Bountra, C.; Arrowsmith, C.H.; Weigelt, J.; Edwards, A.; Kavanagh, K.L.; Oppermann, U.
Deposited on : 2010-01-29
Resolution : 1.92 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

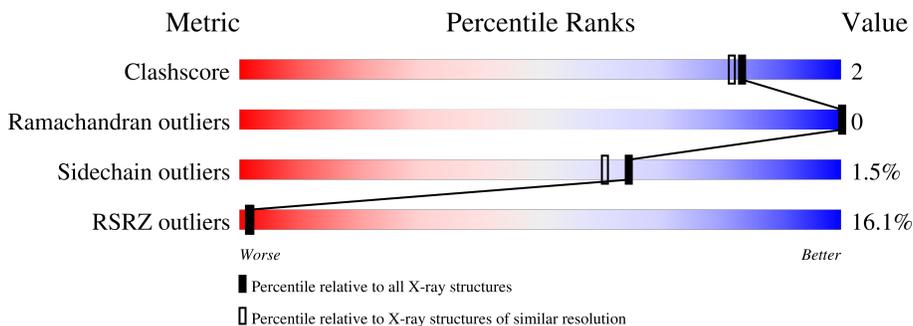
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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))
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	271	
1	B	271	

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
2	POP	A	500	-	X	-	-
2	POP	B	500	-	X	-	-

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 3704 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.

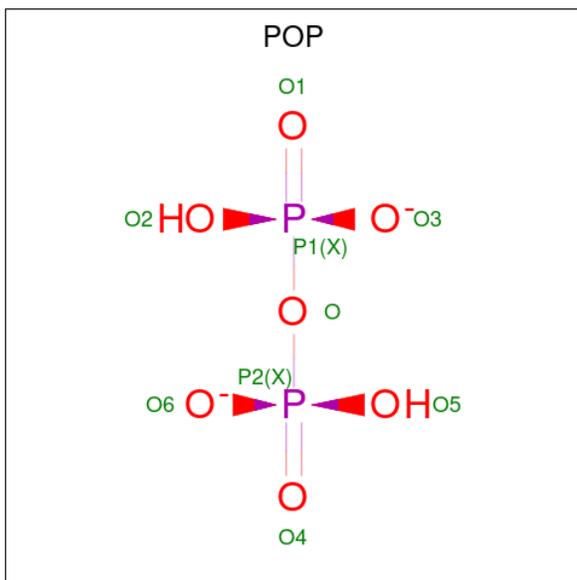
- Molecule 1 is a protein called PHOSPHOLYSINE PHOSPHOHISTIDINE INORGANIC PYROPHOSPHATE PHOSPHATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	270	1986	1249	348	377	12	0	2	0
1	B	213	1478	930	252	285	11	0	2	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	ARG	GLN	SEE REMARK 999	UNP Q9H008
A	160	ALA	LYS	engineered mutation	UNP Q9H008
A	161	ALA	GLU	engineered mutation	UNP Q9H008
B	94	ARG	GLN	SEE REMARK 999	UNP Q9H008
B	160	ALA	LYS	engineered mutation	UNP Q9H008
B	161	ALA	GLU	engineered mutation	UNP Q9H008

- Molecule 2 is PYROPHOSPHATE 2- (three-letter code: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).

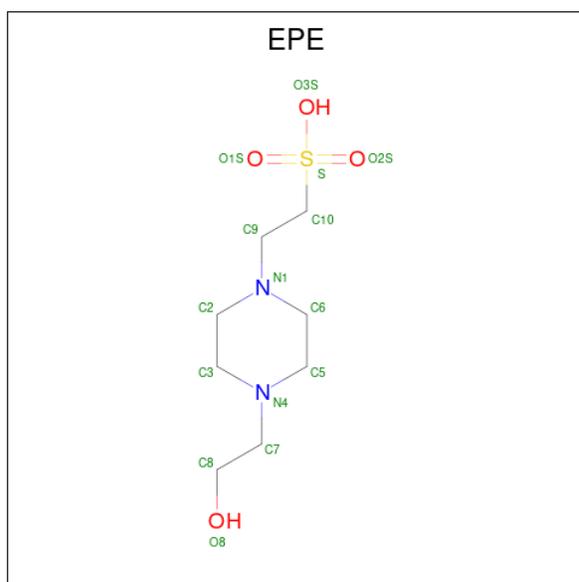


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O S 5 1 3 1	0	0
4	B	1	Total C O S 5 1 3 1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 6 3 3	0	0

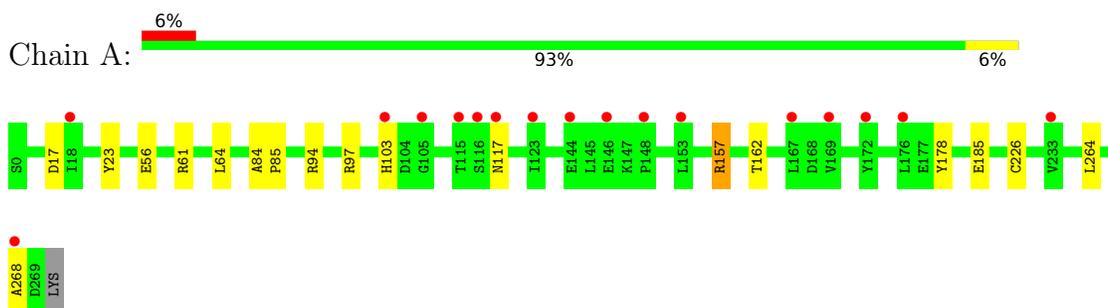
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	185	Total O 185 185	0	0
6	B	27	Total O 27 27	0	0

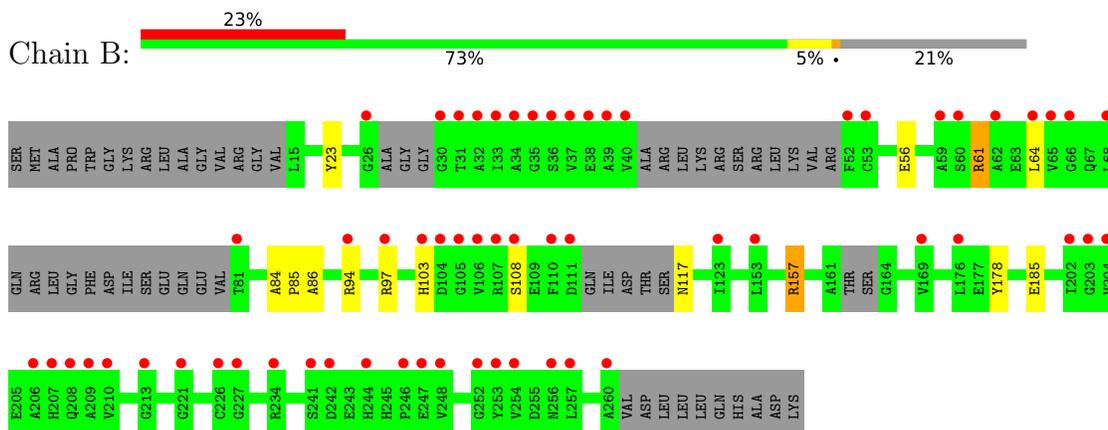
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: PHOSPHOLYSINE PHOSPHOHISTIDINE INORGANIC PYROPHOSPHATE PHOSPHATASE



- Molecule 1: PHOSPHOLYSINE PHOSPHOHISTIDINE INORGANIC PYROPHOSPHATE PHOSPHATASE



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.21Å 39.41Å 112.92Å 90.00° 91.79° 90.00°	Depositor
Resolution (Å)	40.00 – 1.92 31.35 – 1.92	Depositor EDS
% Data completeness (in resolution range)	94.2 (40.00-1.92) 94.2 (31.35-1.92)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 1.92Å)	Xtrriage
Refinement program	REFMAC 5.5.0089	Depositor
R, R_{free}	0.227 , 0.258 0.253 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtrriage
Anisotropy	0.554	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.018 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3704	wwPDB-VP
Average B, all atoms (Å ²)	10.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.86% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, POP, MG, EPE

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	0.83	1/2030 (0.0%)	0.84	3/2749 (0.1%)
1	B	0.59	0/1511	0.73	4/2049 (0.2%)
All	All	0.73	1/3541 (0.0%)	0.80	7/4798 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	226	CYS	CB-SG	-5.05	1.73	1.81

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	157	ARG	NE-CZ-NH2	-9.70	115.45	120.30
1	B	157	ARG	NE-CZ-NH2	-7.37	116.61	120.30
1	A	157	ARG	NE-CZ-NH1	6.83	123.72	120.30
1	A	61	ARG	NE-CZ-NH2	-6.31	117.15	120.30
1	B	157	ARG	NE-CZ-NH1	5.57	123.08	120.30
1	B	61[A]	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	B	61[B]	ARG	NE-CZ-NH2	-5.56	117.52	120.30

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	1986	0	1947	9	0
1	B	1478	0	1357	9	0
2	A	5	0	0	1	0
2	B	5	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	5	0	1	0	0
4	B	5	0	1	0	0
5	A	6	0	8	0	0
6	A	185	0	0	1	1
6	B	27	0	0	1	0
All	All	3704	0	3314	16	1

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 2.

All (16) 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:A:157:ARG:HD3	1:B:178:TYR:O	1.80	0.80
1:A:178:TYR:O	1:B:157:ARG:HD3	1.91	0.70
1:B:86:ALA:HB2	6:B:2002:HOH:O	1.91	0.69
1:A:162:THR:HG22	6:A:2014:HOH:O	1.98	0.63
1:B:61[B]:ARG:CZ	1:B:85:PRO:HG2	2.33	0.58
1:A:56:GLU:HG2	1:A:64:LEU:HD22	1.87	0.55
1:A:84:ALA:HB3	1:A:85:PRO:HD3	1.89	0.55
1:B:61[B]:ARG:NH1	1:B:85:PRO:HG2	2.22	0.54
1:B:56:GLU:HG2	1:B:64:LEU:HD22	1.92	0.52
1:A:264:LEU:O	1:A:268:ALA:HB3	2.14	0.47
1:B:94:ARG:NH2	1:B:185:GLU:OE1	2.48	0.46
1:A:94:ARG:NH2	1:A:185:GLU:OE1	2.49	0.45
1:B:84:ALA:HB3	1:B:85:PRO:HD3	1.99	0.43
1:B:97:ARG:NH1	1:B:117:ASN:O	2.43	0.43
1:A:17:ASP:OD1	2:A:500:POP:P1	2.77	0.43
1:A:97:ARG:NH1	1:A:117:ASN:O	2.50	0.42

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	Interatomic distance (Å)	Clash overlap (Å)
6:A:2028:HOH:O	6:A:2141:HOH:O[1_545]	2.04	0.16

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	270/271 (100%)	263 (97%)	7 (3%)	0	100	100
1	B	203/271 (75%)	197 (97%)	6 (3%)	0	100	100
All	All	473/542 (87%)	460 (97%)	13 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	198/211 (94%)	196 (99%)	2 (1%)	76	75
1	B	135/211 (64%)	131 (97%)	4 (3%)	41	31
All	All	333/422 (79%)	327 (98%)	6 (2%)	65	53

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	TYR
1	A	103	HIS
1	B	23	TYR
1	B	103	HIS
1	B	108[A]	SER
1	B	108[B]	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are

no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	POP	B	500	3	4,4,8	2.14	2 (50%)	6,6,13	2.44	3 (50%)
2	POP	A	500	3	4,4,8	2.21	2 (50%)	6,6,13	1.97	4 (66%)
4	EPE	A	502	-	4,4,15	1.84	1 (25%)	5,6,20	2.24	1 (20%)
5	GOL	A	503	-	5,5,5	0.61	0	5,5,5	1.00	0
4	EPE	B	502	-	4,4,15	1.47	0	5,6,20	2.86	2 (40%)

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
5	GOL	A	503	-	-	2/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	POP	P1-O	3.78	1.66	1.54
2	B	500	POP	P1-O	3.24	1.64	1.54
2	B	500	POP	P1-O1	2.78	1.57	1.50
4	A	502	EPE	O1S-S	2.67	1.54	1.43
2	A	500	POP	P1-O1	2.23	1.56	1.50

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	502	EPE	O2S-S-O1S	-4.91	104.77	118.02
4	A	502	EPE	O2S-S-C10	-4.14	98.91	107.54
2	B	500	POP	O2-P1-O1	3.90	125.16	110.89
4	B	502	EPE	O1S-S-C10	3.83	115.53	107.54
2	B	500	POP	O3-P1-O2	-2.97	98.44	107.97
2	A	500	POP	O3-P1-O2	2.81	116.98	107.97
2	B	500	POP	O-P1-O2	-2.71	99.26	107.97
2	A	500	POP	O3-P1-O1	-2.62	101.31	110.89
2	A	500	POP	O-P1-O3	2.07	114.62	107.97
2	A	500	POP	O-P1-O2	-2.05	101.38	107.97

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	503	GOL	O1-C1-C2-C3
5	A	503	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	POP	1	0

5.7 Other polymers [i](#)

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	270/271 (99%)	0.57	17 (6%) 20 22	4, 9, 15, 27	0
1	B	213/271 (78%)	1.47	61 (28%) 0 0	7, 9, 12, 14	0
All	All	483/542 (89%)	0.97	78 (16%) 1 2	4, 9, 13, 27	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	260	ALA	8.5
1	B	105	GLY	7.2
1	B	40	VAL	6.7
1	B	252	GLY	6.0
1	B	68	LEU	5.6
1	B	206	ALA	5.5
1	B	227	GLY	5.1
1	B	33	ILE	4.8
1	B	52	PHE	4.8
1	B	35	GLY	4.5
1	B	203	GLY	4.4
1	B	65	VAL	4.4
1	B	209	ALA	4.1
1	B	34	ALA	4.0
1	B	207	HIS	3.9
1	B	60	SER	3.9
1	B	248	VAL	3.8
1	B	247	GLU	3.8
1	A	176	LEU	3.8
1	B	107	ARG	3.8
1	A	116	SER	3.7
1	B	246	PRO	3.7
1	B	53	CYS	3.7
1	B	62	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	257	LEU	3.4
1	A	172	TYR	3.4
1	B	110	PHE	3.4
1	A	167	LEU	3.3
1	A	117	ASN	3.2
1	B	39	ALA	3.2
1	B	37	VAL	3.2
1	B	213	GLY	3.1
1	B	208	GLN	3.1
1	A	169	VAL	3.1
1	B	256	ASN	3.1
1	B	153	LEU	3.1
1	B	221	GLY	3.1
1	B	204	VAL	3.0
1	B	244	HIS	2.9
1	A	153	LEU	2.9
1	B	108[A]	SER	2.8
1	A	115	THR	2.8
1	A	18	ILE	2.8
1	B	202	ILE	2.8
1	B	32	ALA	2.8
1	A	268	ALA	2.7
1	B	111	ASP	2.7
1	B	66	GLY	2.7
1	B	36	SER	2.7
1	B	226	CYS	2.6
1	B	169	VAL	2.6
1	B	123	ILE	2.6
1	B	59	ALA	2.5
1	A	146	GLU	2.5
1	A	144	GLU	2.4
1	B	234	ARG	2.4
1	B	242	ASP	2.4
1	B	103	HIS	2.4
1	A	148	PRO	2.3
1	B	97	ARG	2.3
1	B	210	VAL	2.3
1	B	31	THR	2.3
1	B	94	ARG	2.3
1	B	106	VAL	2.2
1	B	26	GLY	2.2
1	A	105	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	81	THR	2.2
1	B	38	GLU	2.2
1	A	123	ILE	2.2
1	B	241	SER	2.2
1	B	30	GLY	2.2
1	B	64	LEU	2.1
1	A	103	HIS	2.1
1	B	253	TYR	2.1
1	A	233	VAL	2.1
1	B	104	ASP	2.1
1	B	176	LEU	2.0
1	B	254	VAL	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 monosaccharides 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. 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	B-factors(Å ²)	Q<0.9
4	EPE	A	502	5/15	0.81	0.39	34,35,39,43	0
2	POP	B	500	5/9	0.85	0.33	28,30,34,35	0
5	GOL	A	503	6/6	0.89	0.14	25,28,29,32	0
3	MG	B	501	1/1	0.92	0.12	27,27,27,27	0
4	EPE	B	502	5/15	0.93	0.51	33,35,38,42	0
2	POP	A	500	5/9	0.96	0.12	18,23,26,28	0
3	MG	A	501	1/1	0.99	0.21	9,9,9,9	0

6.5 Other polymers [i](#)

There are no such residues in this entry.