



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 12:17 PM GMT

PDB ID : 3ALP
Title : Cell adhesion protein
Authors : Narita, H.; Nakagawa, A.; Suzuki, M.
Deposited on : 2010-08-05
Resolution : 2.80 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>

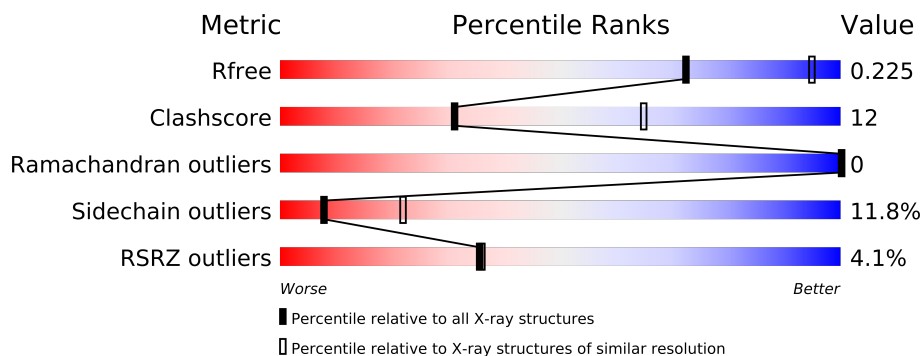
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.80 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}	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (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. 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	331	
1	B	331	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4727 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 Poliovirus receptor-related protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	0	0	0
			2299	1442	399	447	11			
1	B	304	Total	C	N	O	S	0	1	0
			2357	1479	409	458	11			

There are 50 discrepancies between the modelled and reference sequences:

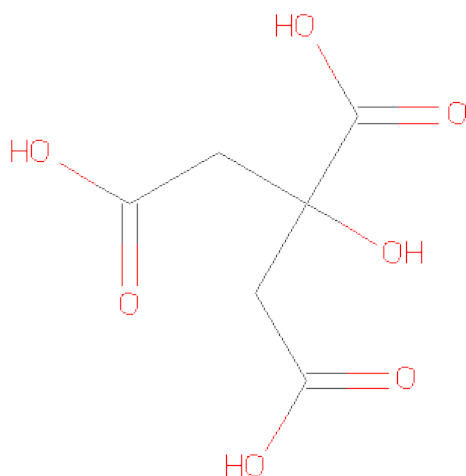
Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	EXPRESSION TAG	UNP Q15223
A	17	ALA	-	EXPRESSION TAG	UNP Q15223
A	18	SER	-	EXPRESSION TAG	UNP Q15223
A	19	MET	-	EXPRESSION TAG	UNP Q15223
A	20	THR	-	EXPRESSION TAG	UNP Q15223
A	21	GLY	-	EXPRESSION TAG	UNP Q15223
A	22	GLY	-	EXPRESSION TAG	UNP Q15223
A	23	GLN	-	EXPRESSION TAG	UNP Q15223
A	24	GLN	-	EXPRESSION TAG	UNP Q15223
A	25	MET	-	EXPRESSION TAG	UNP Q15223
A	26	GLY	-	EXPRESSION TAG	UNP Q15223
A	27	ARG	-	EXPRESSION TAG	UNP Q15223
A	28	ASP	-	EXPRESSION TAG	UNP Q15223
A	29	PRO	-	EXPRESSION TAG	UNP Q15223
A	336	ALA	-	EXPRESSION TAG	UNP Q15223
A	337	ALA	-	EXPRESSION TAG	UNP Q15223
A	338	ALA	-	EXPRESSION TAG	UNP Q15223
A	339	LEU	-	EXPRESSION TAG	UNP Q15223
A	340	GLU	-	EXPRESSION TAG	UNP Q15223
A	341	HIS	-	EXPRESSION TAG	UNP Q15223
A	342	HIS	-	EXPRESSION TAG	UNP Q15223
A	343	HIS	-	EXPRESSION TAG	UNP Q15223
A	344	HIS	-	EXPRESSION TAG	UNP Q15223
A	345	HIS	-	EXPRESSION TAG	UNP Q15223
A	346	HIS	-	EXPRESSION TAG	UNP Q15223

Continued on next page...

Continued from previous page...

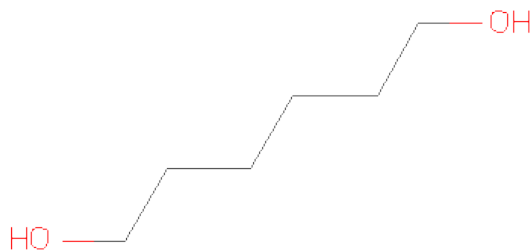
Chain	Residue	Modelled	Actual	Comment	Reference
B	16	MET	-	EXPRESSION TAG	UNP Q15223
B	17	ALA	-	EXPRESSION TAG	UNP Q15223
B	18	SER	-	EXPRESSION TAG	UNP Q15223
B	19	MET	-	EXPRESSION TAG	UNP Q15223
B	20	THR	-	EXPRESSION TAG	UNP Q15223
B	21	GLY	-	EXPRESSION TAG	UNP Q15223
B	22	GLY	-	EXPRESSION TAG	UNP Q15223
B	23	GLN	-	EXPRESSION TAG	UNP Q15223
B	24	GLN	-	EXPRESSION TAG	UNP Q15223
B	25	MET	-	EXPRESSION TAG	UNP Q15223
B	26	GLY	-	EXPRESSION TAG	UNP Q15223
B	27	ARG	-	EXPRESSION TAG	UNP Q15223
B	28	ASP	-	EXPRESSION TAG	UNP Q15223
B	29	PRO	-	EXPRESSION TAG	UNP Q15223
B	336	ALA	-	EXPRESSION TAG	UNP Q15223
B	337	ALA	-	EXPRESSION TAG	UNP Q15223
B	338	ALA	-	EXPRESSION TAG	UNP Q15223
B	339	LEU	-	EXPRESSION TAG	UNP Q15223
B	340	GLU	-	EXPRESSION TAG	UNP Q15223
B	341	HIS	-	EXPRESSION TAG	UNP Q15223
B	342	HIS	-	EXPRESSION TAG	UNP Q15223
B	343	HIS	-	EXPRESSION TAG	UNP Q15223
B	344	HIS	-	EXPRESSION TAG	UNP Q15223
B	345	HIS	-	EXPRESSION TAG	UNP Q15223
B	346	HIS	-	EXPRESSION TAG	UNP Q15223

- Molecule 2 is CITRIC ACID (three-letter code: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is HEXANE-1,6-DIOL (three-letter code: HEZ) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	15	Total	O	0	0
			15	15		
4	B	22	Total	O	0	0
			22	22		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	164.92Å 164.92Å 164.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.99 – 2.80 47.61 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.2 (35.99-2.80) 97.9 (47.61-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.30 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1.357)	Depositor
R, R_{free}	0.190 , 0.226 0.190 , 0.225	Depositor DCC
R_{free} test set	1810 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	71.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.0	EDS
Estimated twinning fraction	0.022 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 36079 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4727	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% 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: HEZ, CIT

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	0.48	0/2350	0.64	0/3212
1	B	0.48	0/2409	0.63	0/3289
All	All	0.48	0/4759	0.63	0/6501

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	A	0	9
1	B	0	3
All	All	0	12

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	259	TYR	Peptide
1	A	263	MET	Peptide
1	A	296	GLN	Peptide
1	A	303	LYS	Peptide
1	A	311	ALA	Peptide
1	A	322	ILE	Peptide
1	A	323	GLY	Peptide
1	A	59	SER	Peptide
1	A	71	THR	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	B	71	THR	Peptide
1	B	73	GLY	Peptide
1	B	74	SER	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	2299	0	0	38	0
1	B	2357	0	0	16	0
2	A	13	0	5	1	0
2	B	13	0	5	0	0
3	B	8	0	14	0	0
4	A	15	0	0	0	0
4	B	22	0	0	1	0
All	All	4727	0	24	55	0

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 12.

All (55) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:263:MET:SD	1:A:263:MET:N	2.30	1.03
1:A:261:GLN:O	1:A:261:GLN:CG	2.30	0.80
1:A:55:ASN:O	1:A:57:LEU:CD2	2.30	0.79
1:A:101:ARG:CG	1:A:101:ARG:O	2.30	0.78
1:A:56:PRO:O	1:A:57:LEU:CD2	2.33	0.77
1:B:259:TYR:CE1	1:B:262:ARG:CG	2.69	0.75
1:A:298:ARG:CG	1:A:298:ARG:NH1	2.51	0.73
1:A:53:PHE:CE1	1:A:134:ARG:NH2	2.59	0.71
1:A:71:THR:CG2	1:A:71:THR:O	2.39	0.71
1:B:259:TYR:CE1	1:B:262:ARG:CB	2.76	0.68
1:A:320:ASN:O	1:A:323:GLY:N	2.30	0.65
1:A:320:ASN:OD1	1:A:321:PRO:N	2.30	0.64
1:A:119:GLU:OE2	1:A:178:LYS:NZ	2.32	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:158:ARG:CG	1:B:158:ARG:NH1	2.59	0.63
1:A:297:ASN:N	1:A:297:ASN:OD1	2.30	0.63
1:A:320:ASN:C	1:A:320:ASN:OD1	2.37	0.63
1:A:53:PHE:CD1	1:A:53:PHE:C	2.71	0.62
1:B:101:ARG:O	1:B:101:ARG:CG	2.46	0.61
1:A:293:VAL:CG1	1:A:302:PHE:CD1	2.85	0.60
1:A:56:PRO:C	1:A:57:LEU:CD2	2.71	0.59
1:A:320:ASN:OD1	1:A:322:ILE:N	2.36	0.59
1:A:113:ARG:NH1	1:A:113:ARG:CG	2.65	0.58
1:A:259:TYR:CB	1:A:260:LEU:O	2.53	0.56
1:B:260:LEU:O	1:B:261:GLN:CB	2.54	0.56
1:A:53:PHE:CD1	1:A:54:ALA:N	2.75	0.54
1:B:53:PHE:C	1:B:53:PHE:CD1	2.81	0.54
1:A:265:VAL:CG1	1:A:265:VAL:O	2.56	0.52
1:B:113:ARG:NH1	1:B:113:ARG:CG	2.72	0.52
1:A:296:GLN:O	1:A:299:THR:OG1	2.26	0.52
1:A:134:ARG:CG	1:A:134:ARG:NH1	2.73	0.52
1:B:70:SER:CA	1:B:74:SER:O	2.59	0.50
1:B:71:THR:O	1:B:74:SER:OG	2.30	0.50
2:A:1:CIT:O4	2:A:1:CIT:H21	2.12	0.49
1:A:319:THR:OG1	1:A:324:THR:OG1	2.30	0.49
1:A:298:ARG:O	1:A:298:ARG:CG	2.61	0.48
1:B:70:SER:CB	1:B:74:SER:O	2.62	0.48
1:A:96:ARG:NH2	1:A:118:ASP:OD2	2.47	0.47
1:A:297:ASN:O	1:A:298:ARG:CB	2.62	0.47
1:B:128:THR:OG1	1:B:131:THR:CG2	2.63	0.46
1:A:220:HIS:NE2	1:A:274:ASN:ND2	2.62	0.46
1:A:262:ARG:O	1:A:304:GLY:O	2.34	0.46
1:A:309:SER:O	1:A:310:LEU:CB	2.64	0.45
1:A:94:ARG:N	1:A:94:ARG:NE	2.64	0.45
1:A:284:THR:CG2	1:A:288:SER:O	2.65	0.45
1:A:53:PHE:CD1	1:A:134:ARG:NH2	2.86	0.43
1:A:320:ASN:CB	1:A:321:PRO:CD	2.96	0.43
1:A:246:GLU:CA	1:A:322:ILE:CG2	2.97	0.43
1:B:119:GLU:OE2	1:B:178:LYS:NZ	2.52	0.42
1:A:231:HIS:CG	1:A:231:HIS:O	2.73	0.42
1:A:246:GLU:CB	1:A:322:ILE:CG2	2.98	0.42
1:B:96:ARG:NH2	1:B:118:ASP:OD2	2.52	0.42
1:B:259:TYR:C	1:B:259:TYR:CD1	2.93	0.42
1:B:261:GLN:NE2	4:B:351:HOH:O	2.52	0.41
1:B:196:GLN:OE1	1:B:210:ARG:NH2	2.53	0.41
1:A:222:GLN:NE2	1:A:222:GLN:CA	2.84	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	299/331 (90%)	286 (96%)	13 (4%)	0	100	100
1	B	303/331 (92%)	293 (97%)	10 (3%)	0	100	100
All	All	602/662 (91%)	579 (96%)	23 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	246/286 (86%)	214 (87%)	32 (13%)	6	17
1	B	256/286 (90%)	229 (90%)	27 (10%)	10	27
All	All	502/572 (88%)	443 (88%)	59 (12%)	8	22

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	PHE
1	A	57	LEU
1	A	71	THR
1	A	76	GLN
1	A	84	SER
1	A	94	ARG
1	A	101	ARG
1	A	103	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	111	LEU
1	A	113	ARG
1	A	114	LEU
1	A	116	LEU
1	A	134	ARG
1	A	140	LEU
1	A	153	THR
1	A	165	ASP
1	A	213	LEU
1	A	226	CYS
1	A	228	VAL
1	A	232	MET
1	A	261	GLN
1	A	263	MET
1	A	265	VAL
1	A	296	GLN
1	A	298	ARG
1	A	299	THR
1	A	300	LEU
1	A	307	ASN
1	A	319	THR
1	A	320	ASN
1	A	322	ILE
1	A	333	ILE
1	B	59	SER
1	B	70	SER
1	B	74	SER
1	B	111	LEU
1	B	113	ARG
1	B	116	LEU
1	B	130	PRO
1	B	131	THR
1	B	140	LEU
1	B	158	ARG
1	B	161	LYS
1	B	163	GLN
1	B	167	VAL
1	B	169	VAL
1	B	181	SER
1	B	213	LEU
1	B	226	CYS
1	B	232	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	259	TYR
1	B	260	LEU
1	B	262	ARG
1	B	263	MET
1	B	267	LEU
1	B	298	ARG
1	B	299	THR
1	B	307	ASN
1	B	309	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

3 ligands are 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	$\# Z > 2$
2	CIT	A	1	-	12,12,12	0.93	0	17,17,17	1.79	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HEZ	B	1	-	7,7,7	0.26	0	6,6,6	0.65	0
2	CIT	B	2	-	12,12,12	0.89	0	17,17,17	1.64	3 (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
2	CIT	A	1	-	-	0/16/16/16	0/0/0/0
3	HEZ	B	1	-	-	0/5/5/5	0/0/0/0
2	CIT	B	2	-	-	0/16/16/16	0/0/0/0

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	CIT	O6-C6-C3	5.19	120.44	112.89
2	B	2	CIT	O6-C6-C3	4.78	119.84	112.89
2	A	1	CIT	O2-C1-C2	2.41	123.14	114.63
2	B	2	CIT	C3-C4-C5	-2.27	108.27	113.77
2	A	1	CIT	O2-C1-O1	-2.19	117.73	123.30
2	B	2	CIT	O4-C5-C4	2.12	122.10	114.63

There are no chirality outliers.

There are no torsion outliers.

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	301/331 (90%)	0.18	16 (5%) 25 26	38, 76, 168, 230	0
1	B	304/331 (91%)	0.03	9 (2%) 48 49	40, 66, 123, 164	0
All	All	605/662 (91%)	0.11	25 (4%) 35 36	38, 69, 153, 230	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	258	TRP	5.0
1	A	256	GLY	4.3
1	B	33	VAL	4.1
1	A	257	ASN	3.7
1	A	302	PHE	3.6
1	A	336	ALA	3.2
1	A	310	LEU	3.2
1	B	293	VAL	3.1
1	A	267	LEU	2.9
1	B	296	GLN	2.8
1	A	313	THR	2.8
1	A	314	TYR	2.8
1	A	254	PHE	2.7
1	A	332	ASN	2.7
1	B	54	ALA	2.6
1	B	62	ILE	2.5
1	B	55	ASN	2.5
1	A	265	VAL	2.2
1	B	104	PHE	2.2
1	A	263	MET	2.2
1	B	302	PHE	2.2
1	A	294	GLU	2.1
1	A	331	VAL	2.1
1	B	264	ASP	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	277	ALA	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
3	HEZ	B	1	8/8	0.22	1.61	55,57,69,78	0
2	CIT	B	2	13/13	0.20	0.67	88,100,128,131	0
2	CIT	A	1	13/13	0.16	-0.35	91,114,124,125	0

6.5 Other polymers ⓘ

There are no such residues in this entry.