



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:38 PM GMT

PDB ID : 3GWD
Title : Closed crystal structure of cyclohexanone monooxygenase
Authors : Mirza, I.A.; Yachnin, B.J.; Berghuis, A.M.
Deposited on : 2009-03-31
Resolution : 2.30 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>

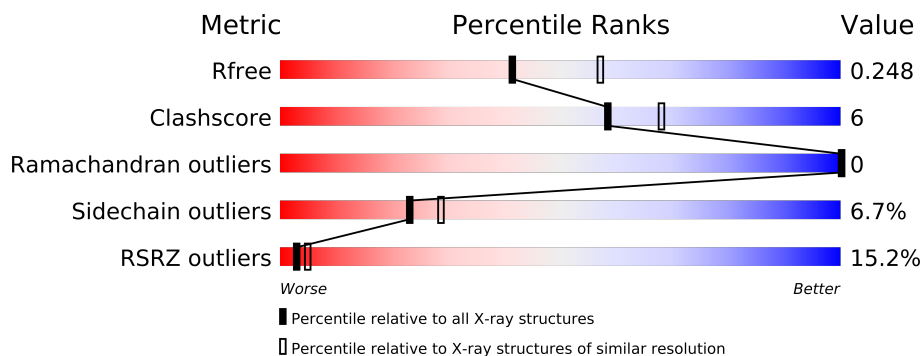
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.30 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	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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	540	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 4454 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.

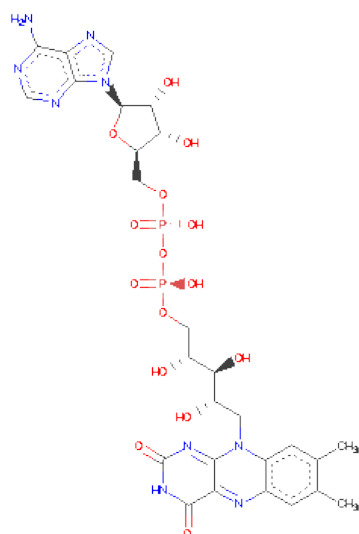
- Molecule 1 is a protein called Cyclohexanone monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	531	4178	2660	709	801	8	0	1	0

There are 4 discrepancies between the modelled and reference sequences:

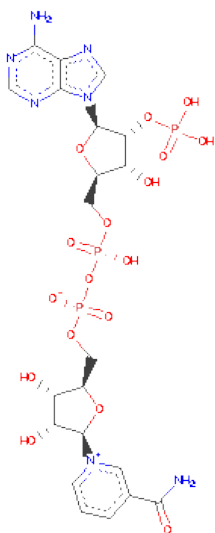
Chain	Residue	Modelled	Actual	Comment	Reference
A	6	THR	ILE	CONFLICT	UNP Q6RXW1
A	81	THR	SER	CONFLICT	UNP Q6RXW1
A	223	ASN	THR	CONFLICT	UNP Q6RXW1
A	323	LYS	THR	CONFLICT	UNP Q6RXW1

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	53	27	9	15	2	0	0

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	A	1	Total	C	N	O	P	0	0
			26	11	2	11	2		

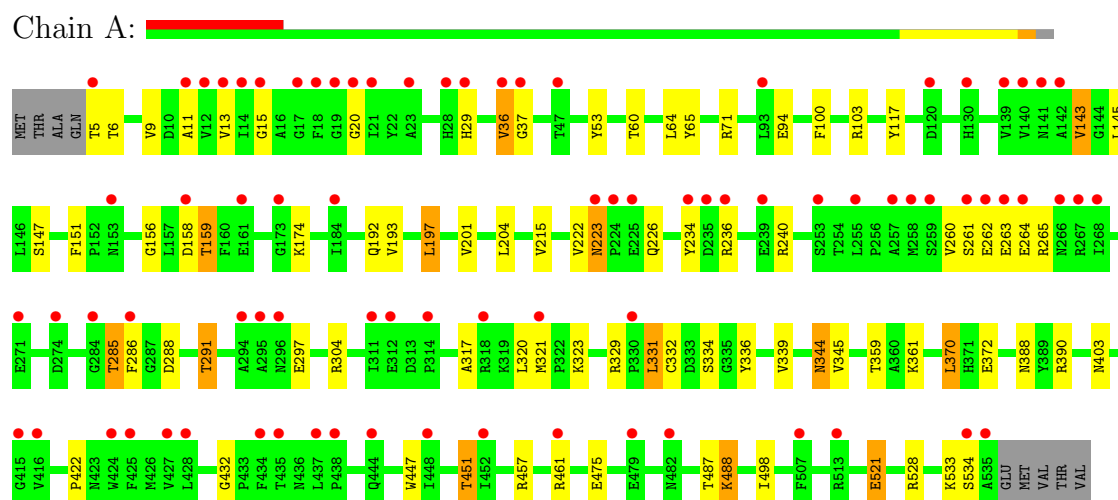
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	149	Total	O	0	0
			149	149		

3 Residue-property plots

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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.48Å 90.99Å 100.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.53 – 2.30 26.53 – 2.29	Depositor EDS
% Data completeness (in resolution range)	99.9 (26.53-2.30) 99.7 (26.53-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.183 , 0.245 0.184 , 0.248	Depositor DCC
R_{free} test set	2644 reflections (11.20%)	DCC
Wilson B-factor (Å ²)	40.7	Xtriage
Anisotropy	0.199	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 34.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 26246 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4454	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% 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: NAP, FAD

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	0.71	4/4292 (0.1%)	0.68	0/5848

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	263	GLU	CD-OE2	16.64	1.44	1.25
1	A	264	GLU	CD-OE1	10.06	1.36	1.25
1	A	264	GLU	CD-OE2	9.11	1.35	1.25
1	A	264	GLU	CG-CD	8.01	1.64	1.51

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	4178	0	4000	51	0
2	A	53	0	31	3	0
3	A	74	0	39	3	0
4	A	149	0	0	2	0
All	All	4454	0	4070	52	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 6.

All (52) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:147:SER:HB2	3:A:543:NAP:N7N	1.97	0.79
1:A:475:GLU:HG3	4:A:693:HOH:O	1.82	0.77
1:A:297:GLU:OE2	1:A:304:ARG:NH2	2.23	0.70
1:A:344:ASN:ND2	1:A:345:VAL:HG23	2.07	0.69
1:A:336:TYR:O	1:A:339:VAL:HG22	1.92	0.69
1:A:94:GLU:OE1	1:A:103:ARG:NH2	2.26	0.68
1:A:156:GLY:O	1:A:159:THR:HB	1.97	0.64
1:A:236:ARG:HE	1:A:240:ARG:HH12	1.45	0.63
1:A:260:VAL:HG22	1:A:261:SER:H	1.65	0.62
1:A:53:TYR:OH	1:A:192:GLN:NE2	2.34	0.60
1:A:344:ASN:HD22	1:A:344:ASN:C	2.05	0.59
1:A:223:ASN:ND2	1:A:226:GLN:H	2.00	0.59
1:A:215:VAL:O	1:A:332:CYS:HA	2.02	0.59
1:A:223:ASN:HD22	1:A:226:GLN:H	1.52	0.58
1:A:390:ARG:HH11	1:A:403:ASN:HD22	1.52	0.57
1:A:236:ARG:NE	1:A:240:ARG:HH12	2.02	0.57
2:A:541:FAD:HM73	3:A:542:NAP:C5N	2.35	0.56
1:A:223:ASN:HD22	1:A:223:ASN:C	2.11	0.54
1:A:447:TRP:CZ3	1:A:451:THR:HG21	2.42	0.54
1:A:329:ARG:O	1:A:331:LEU:HD13	2.08	0.52
1:A:390:ARG:HH11	1:A:403:ASN:ND2	2.07	0.52
1:A:222:VAL:HG13	1:A:226:GLN:HB3	1.92	0.51
1:A:147:SER:OG	1:A:388:ASN:ND2	2.45	0.50
1:A:262:GLU:OE1	1:A:265:ARG:NH1	2.45	0.49
1:A:9:VAL:HG11	1:A:36:VAL:HG23	1.95	0.49
1:A:432:GLY:HA3	4:A:593:HOH:O	2.13	0.48
1:A:288:ASP:HB2	1:A:291:THR:HG22	1.96	0.47
1:A:317:ALA:O	1:A:321[A]:MET:HG2	2.13	0.47
1:A:461:ARG:NH1	1:A:534:SER:O	2.47	0.47
1:A:60:THR:HB	1:A:65:TYR:HB2	1.97	0.46
1:A:143:VAL:HG22	2:A:541:FAD:C8A	2.46	0.46
1:A:15:GLY:O	1:A:20:GLY:HA3	2.15	0.46
1:A:521:GLU:OE2	1:A:528:ARG:NH1	2.41	0.46
1:A:204:LEU:HB3	1:A:345:VAL:HG22	1.98	0.45
1:A:447:TRP:O	1:A:451:THR:CG2	2.65	0.44
1:A:461:ARG:NH2	1:A:533:LYS:HE2	2.32	0.44
1:A:370:LEU:HD13	1:A:372:GLU:HG2	1.99	0.44
1:A:344:ASN:HD21	1:A:345:VAL:HG23	1.80	0.44
1:A:29:HIS:HE1	1:A:100:PHE:O	2.01	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:147:SER:HB2	3:A:543:NAP:H72N	1.77	0.43
1:A:344:ASN:ND2	1:A:344:ASN:C	2.71	0.43
1:A:143:VAL:HG22	2:A:541:FAD:N7A	2.34	0.43
1:A:285:THR:CG2	1:A:286:PHE:N	2.83	0.42
1:A:117:TYR:CZ	1:A:422:PRO:HG2	2.55	0.42
1:A:13:VAL:O	1:A:37:GLY:HA2	2.21	0.41
1:A:260:VAL:HG22	1:A:261:SER:N	2.34	0.41
1:A:288:ASP:HA	1:A:291:THR:HG22	2.02	0.41
1:A:9:VAL:HG12	1:A:11:ALA:H	1.86	0.41
1:A:193:VAL:HG22	1:A:197:LEU:HD22	2.01	0.41
1:A:390:ARG:CG	1:A:403:ASN:HD21	2.33	0.41
1:A:487:THR:HG23	1:A:488:LYS:HG2	2.02	0.40
1:A:159:THR:HG22	1:A:359:THR:C	2.42	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	530/540 (98%)	514 (97%)	16 (3%)	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	435/442 (98%)	406 (93%)	29 (7%)	23	29

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	6	THR
1	A	36	VAL
1	A	64	LEU
1	A	71	ARG
1	A	143	VAL
1	A	145	LEU
1	A	151	PHE
1	A	158	ASP
1	A	159	THR
1	A	174	LYS
1	A	197	LEU
1	A	201	VAL
1	A	223	ASN
1	A	234	TYR
1	A	285	THR
1	A	291	THR
1	A	320	LEU
1	A	323	LYS
1	A	331	LEU
1	A	334	SER
1	A	344	ASN
1	A	361	LYS
1	A	370	LEU
1	A	451	THR
1	A	457	ARG
1	A	488	LYS
1	A	498	ILE
1	A	521	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	141	ASN
1	A	192	GLN
1	A	223	ASN
1	A	353	ASN
1	A	388	ASN
1	A	403	ASN
1	A	431	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 ⓘ

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	FAD	A	541	-	58,58,58	0.98	6 (10%)	85,89,89	1.84	12 (14%)
3	NAP	A	542	-	52,52,52	1.32	3 (5%)	80,80,80	1.91	12 (15%)
3	NAP	A	543	-	27,27,52	1.68	2 (7%)	41,41,80	1.32	4 (9%)

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	FAD	A	541	-	-	0/34/50/50	0/1/6/6
3	NAP	A	542	-	-	0/35/67/67	0/3/5/5
3	NAP	A	543	-	-	0/20/36/67	0/2/2/5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	543	NAP	O7N-C7N	7.27	1.41	1.24
3	A	542	NAP	O7N-C7N	6.88	1.40	1.24
3	A	543	NAP	PA-O5B	3.48	1.62	1.50
3	A	542	NAP	C2A-N3A	2.96	1.38	1.32
2	A	541	FAD	C2A-N3A	2.48	1.37	1.32
3	A	542	NAP	C2N-N1N	2.38	1.38	1.35
2	A	541	FAD	C2A-N1A	2.30	1.38	1.33
2	A	541	FAD	C4X-N5	2.23	1.40	1.36
2	A	541	FAD	C1'-N10	2.16	1.50	1.48
2	A	541	FAD	C5X-N5	2.09	1.38	1.35
2	A	541	FAD	C6-C5X	-2.04	1.39	1.41

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	541	FAD	N3A-C2A-N1A	-11.04	119.48	128.71
3	A	542	NAP	N3A-C2A-N1A	-9.77	120.54	128.71
3	A	542	NAP	C2D-C1D-N1N	-5.99	103.71	113.86
3	A	542	NAP	O4B-C1B-N9A	5.84	113.87	108.44
2	A	541	FAD	C2-N1-C10	5.05	120.07	114.98
3	A	542	NAP	C3N-C7N-N7N	4.25	122.61	117.77
2	A	541	FAD	N3A-C4A-N9A	3.88	132.43	125.43
3	A	542	NAP	O4D-C1D-N1N	3.82	111.86	107.95
3	A	543	NAP	O4D-C1D-N1N	3.69	111.73	107.95
3	A	542	NAP	N3A-C4A-N9A	3.55	131.84	125.43
2	A	541	FAD	C5X-C9A-N10	3.49	120.24	116.80
3	A	543	NAP	C2D-C1D-N1N	-3.45	108.02	113.86
3	A	542	NAP	O4B-C1B-C2B	-3.38	103.79	106.95
2	A	541	FAD	C4-N3-C2	-3.21	118.81	125.39
3	A	542	NAP	O3X-P2B-O2B	3.10	116.01	107.09
3	A	543	NAP	PA-O3-PN	-3.02	123.41	133.69
2	A	541	FAD	C4X-C10-N1	-2.92	119.81	122.73
2	A	541	FAD	C4X-N5-C5X	2.89	119.94	116.69
3	A	542	NAP	C5A-C4A-N3A	-2.66	119.92	125.70
3	A	543	NAP	O2A-PA-O3	2.59	112.53	104.29
3	A	542	NAP	O7N-C7N-C3N	-2.48	116.78	119.58
2	A	541	FAD	N7A-C8A-N9A	-2.41	107.54	114.36
2	A	541	FAD	C2A-N3A-C4A	2.39	120.83	114.01
3	A	542	NAP	C2A-N3A-C4A	2.36	120.74	114.01
2	A	541	FAD	C5A-C4A-N3A	-2.27	120.76	125.70
3	A	542	NAP	C4A-C5A-N7A	-2.23	107.61	109.52
2	A	541	FAD	P-O3P-PA	-2.08	125.59	131.68
2	A	541	FAD	C4X-C4-N3	2.02	121.30	115.39

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	531/540 (98%)	0.77	81 (15%) 3 5	34, 44, 52, 61	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	13	VAL	5.3
1	A	12	VAL	5.1
1	A	257	ALA	4.7
1	A	140	VAL	4.6
1	A	139	VAL	4.6
1	A	263	GLU	4.6
1	A	18	PHE	4.2
1	A	261	SER	3.9
1	A	158	ASP	3.9
1	A	224	PRO	3.9
1	A	130	HIS	3.8
1	A	14	ILE	3.7
1	A	19	GLY	3.6
1	A	535	ALA	3.6
1	A	448	ILE	3.4
1	A	225	GLU	3.4
1	A	416	VAL	3.4
1	A	312	GLU	3.3
1	A	314	PRO	3.3
1	A	236	ARG	3.2
1	A	267	ARG	3.2
1	A	534	SER	3.2
1	A	255	LEU	3.2
1	A	235	ASP	3.2
1	A	20	GLY	3.1
1	A	434	PHE	3.1
1	A	415	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	437	LEU	3.1
1	A	5	THR	3.1
1	A	29	HIS	3.1
1	A	425	PHE	3.1
1	A	223	ASN	3.0
1	A	266	ASN	3.0
1	A	21	ILE	2.9
1	A	28	HIS	2.9
1	A	428	LEU	2.9
1	A	268	ILE	2.9
1	A	286	PHE	2.9
1	A	479	GLU	2.9
1	A	173	GLY	2.9
1	A	435	THR	2.8
1	A	284	GLY	2.8
1	A	153	ASN	2.8
1	A	239	GLU	2.8
1	A	184	ILE	2.8
1	A	507	PHE	2.7
1	A	294	ALA	2.7
1	A	253	SER	2.7
1	A	330	PRO	2.6
1	A	11	ALA	2.6
1	A	142	ALA	2.6
1	A	36	VAL	2.6
1	A	23	ALA	2.6
1	A	259	SER	2.5
1	A	264	GLU	2.5
1	A	120	ASP	2.5
1	A	427	VAL	2.5
1	A	141	ASN	2.4
1	A	234	TYR	2.4
1	A	262	GLU	2.4
1	A	438	PRO	2.4
1	A	47	THR	2.3
1	A	296	ASN	2.3
1	A	258	MET	2.3
1	A	424	TRP	2.3
1	A	461	ARG	2.3
1	A	295	ALA	2.3
1	A	444	GLN	2.2
1	A	321[A]	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	93	LEU	2.2
1	A	274	ASP	2.2
1	A	161	GLU	2.2
1	A	17	GLY	2.2
1	A	37	GLY	2.2
1	A	271	GLU	2.2
1	A	318	ARG	2.1
1	A	452	ILE	2.1
1	A	513	ARG	2.0
1	A	15	GLY	2.0
1	A	311	ILE	2.0
1	A	482	ASN	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	NAP	A	543	26/48	0.17	0.65	33,42,63,65	0
3	NAP	A	542	48/48	0.10	-1.49	17,26,29,30	0
2	FAD	A	541	53/53	0.11	-1.59	22,28,29,31	0

6.5 Other polymers ⓘ

There are no such residues in this entry.