



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:00 AM UTC

PDB ID : 2OK6 / pdb_00002ok6
Title : Crystal structure of aromatic amine dehydrogenase TTQ-formamide adduct oxidized with ferricyanide.
Authors : Roujeinikova, A.; Leys, D.
Deposited on : 2007-01-16
Resolution : 1.45 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

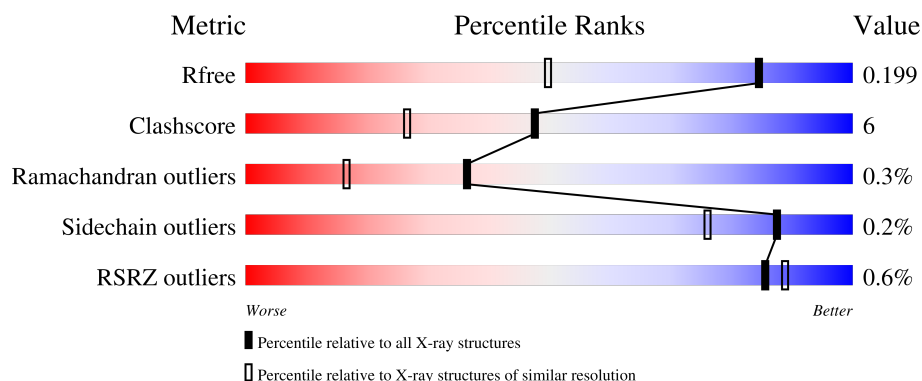
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.45 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}	180053	1756 (1.46-1.46)
Clashscore	190562	1795 (1.46-1.46)
Ramachandran outliers	187476	1776 (1.46-1.46)
Sidechain outliers	187428	1776 (1.46-1.46)
RSRZ outliers	180081	1756 (1.46-1.46)

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	D	136	73% 7% 18%
1	H	136	75% 12% 11%
2	A	361	90% 10%
2	B	361	90% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9130 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 Aromatic amine dehydrogenase, small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	110	Total	C	N	O	S	0	2	0
			869	535	152	167	15			
1	H	120	Total	C	N	O	S	0	5	0
			952	585	166	186	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	109	TQQ	TRP	modified residue	UNP Q0VKG6
D	109	TQQ	TRP	modified residue	UNP Q0VKG6

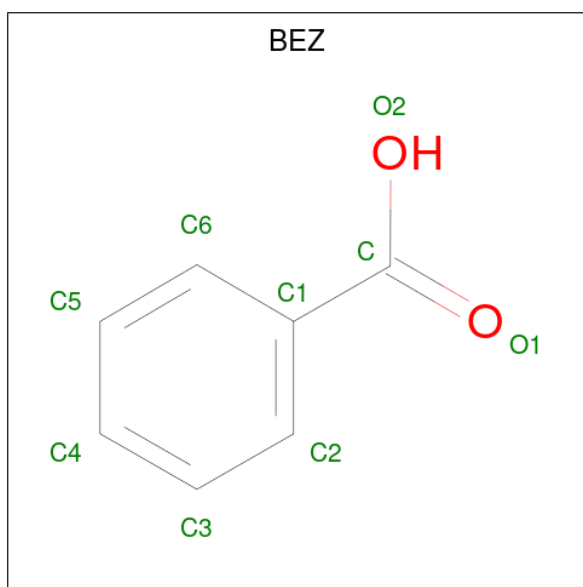
- Molecule 2 is a protein called Aromatic amine dehydrogenase, large subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	359	Total	C	N	O	S	0	2	0
			2805	1770	487	534	14			
2	B	360	Total	C	N	O	S	0	2	0
			2803	1768	487	534	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	433	THR	-	insertion	UNP Q0VKG7
B	433	THR	-	insertion	UNP Q0VKG7

- Molecule 3 is BENZOIC ACID (CCD ID: BEZ) (formula: C₇H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			8	7	1		
3	H	1	Total	C	O	0	0
			8	7	1		

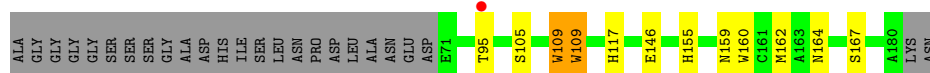
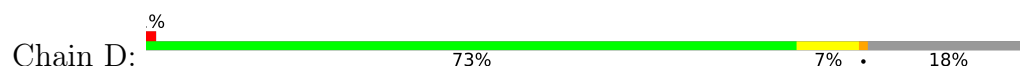
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	180	Total	O	0	1
			180	180		
4	H	179	Total	O	0	1
			179	179		
4	A	701	Total	O	0	0
			701	701		
4	B	625	Total	O	0	0
			625	625		

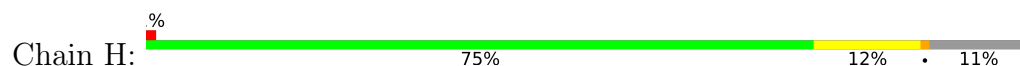
3 Residue-property plots [i](#)

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.

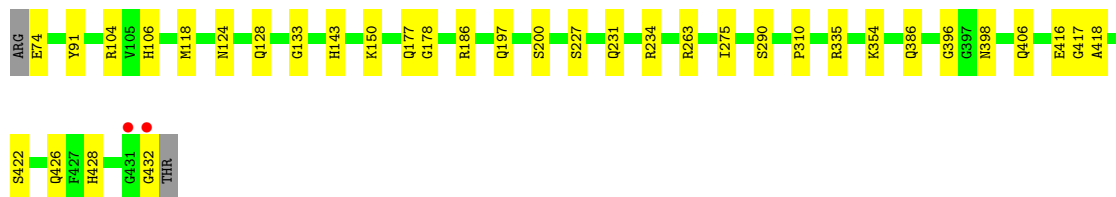
- Molecule 1: Aromatic amine dehydrogenase, small subunit



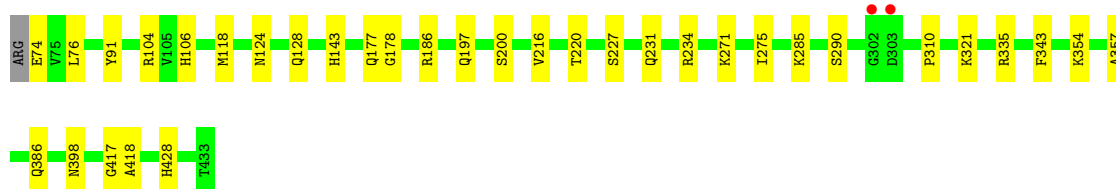
- Molecule 1: Aromatic amine dehydrogenase, small subunit



- Molecule 2: Aromatic amine dehydrogenase, large subunit



- Molecule 2: Aromatic amine dehydrogenase, large subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.34Å 89.27Å 80.49Å 90.00° 90.59° 90.00°	Depositor
Resolution (Å)	15.00 – 1.45 15.00 – 1.45	Depositor EDS
% Data completeness (in resolution range)	100.0 (15.00-1.45) 97.6 (15.00-1.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 1.45Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.161 , 0.192 0.170 , 0.199	Depositor DCC
R_{free} test set	8530 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	14.4	Xtriage
Anisotropy	0.446	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 61.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.009 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9130	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BEZ, 1TQ, TQQ

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	D	0.79	0/862	0.88	0/1174
1	H	0.79	0/953	0.84	0/1302
2	A	0.82	0/2876	0.87	0/3898
2	B	0.75	0/2876	0.86	0/3897
All	All	0.79	0/7567	0.86	0/10271

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
2	A	0	1
2	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	310	PRO	Peptide
2	B	310	PRO	Peptide

5.2 Too-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 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	D	869	0	755	11	0
1	H	952	0	825	19	0
2	A	2805	0	2747	28	0
2	B	2803	0	2735	25	0
3	D	8	0	5	0	0
3	H	8	0	5	0	0
4	A	701	0	0	5	0
4	B	625	0	0	6	0
4	D	180	0	0	3	0
4	H	179	0	0	8	0
All	All	9130	0	7072	76	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 6.

All (76) 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:H:173:THR:HA	4:H:2170:HOH:O	1.71	0.91
2:B:74:GLU:N	4:B:878:HOH:O	2.09	0.86
1:H:80:HIS:O	1:H:142[A]:THR:HG21	1.79	0.82
1:H:95[A]:THR:HG22	4:H:2065:HOH:O	1.82	0.79
2:B:104:ARG:HH11	2:B:106:HIS:HE1	1.31	0.78
2:A:104:ARG:HH11	2:A:106:HIS:HE1	1.30	0.77
4:D:2161:HOH:O	2:A:118[A]:MET:SD	2.44	0.75
4:H:2084:HOH:O	2:B:118[B]:MET:SD	2.45	0.74
1:D:105:SER:HA	1:D:164:ASN:HD21	1.54	0.73
1:H:105:SER:HA	1:H:164:ASN:HD21	1.59	0.68
1:H:95[A]:THR:HG23	4:H:2038:HOH:O	1.93	0.68
2:A:74:GLU:N	4:A:733:HOH:O	2.26	0.67
2:B:91:TYR:OH	2:B:428:HIS:HD2	1.76	0.67
2:A:91:TYR:OH	2:A:428:HIS:HD2	1.78	0.66
2:A:124:ASN:HD21	2:A:178:GLY:H	1.43	0.66
2:B:124:ASN:HD21	2:B:178:GLY:H	1.40	0.66
1:D:160[B]:TRP:CE3	1:D:162:MET:HE3	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:197:GLN:NE2	2:B:227:SER:H	1.95	0.64
2:A:197:GLN:NE2	2:A:227:SER:H	1.98	0.61
2:B:197:GLN:HE21	2:B:227:SER:H	1.49	0.60
1:H:109[A]:TQQ:HB2	1:H:160[A]:TRP:NE1	2.17	0.59
2:A:335:ARG:H	2:A:386:GLN:HE22	1.49	0.58
2:A:91:TYR:OH	2:A:428:HIS:CD2	2.57	0.57
1:H:80:HIS:O	1:H:142[A]:THR:CG2	2.51	0.57
2:B:106:HIS:HD2	2:B:418:ALA:O	1.88	0.56
2:A:335:ARG:H	2:A:386:GLN:NE2	2.03	0.55
1:H:81:CYS:C	1:H:142[A]:THR:HG23	2.32	0.55
1:H:146:GLU:CD	2:A:354:LYS:HZ1	2.15	0.55
2:B:285:LYS:NZ	4:B:832:HOH:O	2.39	0.54
2:B:398:ASN:HD22	2:B:417:GLY:H	1.55	0.54
1:D:167:SER:HB2	2:B:177:GLN:HE22	1.71	0.54
2:A:197:GLN:HE21	2:A:227:SER:H	1.54	0.54
2:A:133:GLY:O	2:A:428:HIS:HE1	1.92	0.53
2:A:106:HIS:HD2	2:A:418:ALA:O	1.92	0.53
2:A:263:ARG:HG3	4:A:1007:HOH:O	2.10	0.52
1:H:173:THR:CA	4:H:2170:HOH:O	2.43	0.52
2:A:128:GLN:HE22	2:A:186:ARG:C	2.19	0.51
1:D:146:GLU:CD	2:B:354:LYS:HZ1	2.19	0.51
2:B:321:LYS:HA	4:B:723:HOH:O	2.11	0.51
1:H:167:SER:HB2	2:A:177:GLN:HE22	1.76	0.50
1:D:117:HIS:HE1	4:D:2048:HOH:O	1.94	0.50
2:B:128:GLN:HE22	2:B:186:ARG:C	2.20	0.50
2:A:398:ASN:HD22	2:A:417:GLY:H	1.60	0.50
1:H:77:TYR:CE1	1:H:79:ARG:HG2	2.47	0.49
1:H:159:ASN:HD22	2:A:177:GLN:NE2	2.11	0.48
2:B:143:HIS:HD2	4:B:566:HOH:O	1.97	0.48
2:B:275:ILE:HA	2:B:290:SER:HA	1.96	0.47
1:H:109[A]:TQQ:HB2	1:H:160[A]:TRP:HE1	1.80	0.47
1:H:160[B]:TRP:CE3	1:H:162:MET:HE3	2.50	0.47
2:B:220:THR:HG22	4:B:803:HOH:O	2.14	0.47
2:B:335:ARG:H	2:B:386:GLN:NE2	2.13	0.47
2:A:275:ILE:HA	2:A:290:SER:HA	1.97	0.46
1:H:154:LEU:HD22	4:H:2125:HOH:O	2.15	0.46
2:A:432:GLY:C	4:A:1048:HOH:O	2.59	0.46
2:B:398:ASN:ND2	2:B:417:GLY:H	2.13	0.46
1:D:160[B]:TRP:CZ3	1:D:162:MET:HE3	2.51	0.45
1:D:155:HIS:HE1	4:B:442:HOH:O	2.00	0.45
2:B:335:ARG:H	2:B:386:GLN:HE22	1.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:231:GLN:NE2	2:A:234:ARG:HH11	2.15	0.43
1:D:159:ASN:HD22	2:B:177:GLN:NE2	2.16	0.43
4:H:2037:HOH:O	2:A:143:HIS:HE1	2.01	0.43
2:A:150:LYS:HE3	4:A:933:HOH:O	2.18	0.43
2:B:124:ASN:ND2	2:B:178:GLY:H	2.12	0.43
2:A:396:GLY:HA2	2:A:422:SER:O	2.18	0.42
2:A:398:ASN:HD22	2:A:416:GLU:HA	1.84	0.42
1:H:155:HIS:HE1	4:A:446:HOH:O	2.01	0.42
2:A:398:ASN:ND2	2:A:417:GLY:H	2.16	0.42
1:D:95:THR:HG22	4:D:2091:HOH:O	2.19	0.42
2:A:124:ASN:ND2	2:A:178:GLY:H	2.15	0.42
1:D:146:GLU:OE2	1:D:155:HIS:HD2	2.03	0.42
2:A:186:ARG:HH22	2:A:426:GLN:NE2	2.17	0.41
2:B:343:PHE:HB3	2:B:357:ALA:HB2	2.02	0.41
1:D:109[A]:TQQ:HB2	1:D:160[A]:TRP:NE1	2.35	0.41
1:H:174:SER:N	4:H:2170:HOH:O	2.49	0.41
2:B:231:GLN:NE2	2:B:234:ARG:HH11	2.19	0.41
1:H:173:THR:HG21	2:B:76:LEU:HD13	2.03	0.41

There are no symmetry-related clashes.

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	D	108/136 (79%)	105 (97%)	3 (3%)	0	100	100
1	H	121/136 (89%)	118 (98%)	3 (2%)	0	100	100
2	A	359/361 (99%)	346 (96%)	12 (3%)	1 (0%)	36	16
2	B	359/361 (99%)	347 (97%)	10 (3%)	2 (1%)	21	6
All	All	947/994 (95%)	916 (97%)	28 (3%)	3 (0%)	36	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	216	VAL
2	A	200	SER
2	B	200	SER

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	D	95/112 (85%)	95 (100%)	0	100	100
1	H	105/112 (94%)	105 (100%)	0	100	100
2	A	305/305 (100%)	303 (99%)	2 (1%)	76	54
2	B	305/305 (100%)	304 (100%)	1 (0%)	86	74
All	All	810/834 (97%)	807 (100%)	3 (0%)	87	70

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	406[A]	GLN
2	A	406[B]	GLN
2	B	271	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	117	HIS
1	D	143	GLN
1	D	155	HIS
1	D	164	ASN
1	H	117	HIS
1	H	139	GLN
1	H	143	GLN
1	H	155	HIS
1	H	164	ASN
2	A	106	HIS
2	A	124	ASN

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Mol	Chain	Res	Type
2	A	128	GLN
2	A	143	HIS
2	A	177	GLN
2	A	180	ASN
2	A	187	GLN
2	A	197	GLN
2	A	231	GLN
2	A	251	ASN
2	A	369	GLN
2	A	386	GLN
2	A	398	ASN
2	A	424	GLN
2	A	426	GLN
2	A	428	HIS
2	B	106	HIS
2	B	124	ASN
2	B	128	GLN
2	B	143	HIS
2	B	177	GLN
2	B	180	ASN
2	B	197	GLN
2	B	231	GLN
2	B	251	ASN
2	B	322	ASN
2	B	386	GLN
2	B	398	ASN
2	B	424	GLN
2	B	426	GLN
2	B	428	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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
1	TQQ	D	109[A]	-	16,17,18	2.13	6 (37%)	14,24,26	0.82	0
1	1TQ	D	109[B]	-	18,19,20	2.91	4 (22%)	17,26,28	1.69	3 (17%)
1	TQQ	H	109[A]	-	16,17,18	2.14	5 (31%)	14,24,26	0.84	0
1	1TQ	H	109[B]	-	18,19,20	3.16	5 (27%)	17,26,28	1.74	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
1	TQQ	D	109[A]	-	-	0/5/19/21	0/2/2/2
1	1TQ	D	109[B]	-	-	3/8/9/11	0/2/2/2
1	TQQ	H	109[A]	-	-	0/5/19/21	0/2/2/2
1	1TQ	H	109[B]	-	-	3/8/9/11	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	109[B]	1TQ	CH3-N2	11.51	1.49	1.34
1	D	109[B]	1TQ	CH3-N2	10.80	1.48	1.34
1	D	109[A]	TQQ	CH2-N2	4.82	1.36	1.28
1	H	109[A]	TQQ	CH2-N2	4.45	1.35	1.28
1	D	109[A]	TQQ	CH2-CZ2	-4.33	1.36	1.48
1	H	109[B]	1TQ	CH2-N2	4.06	1.47	1.41
1	H	109[A]	TQQ	CH2-CZ2	-3.93	1.38	1.48
1	D	109[B]	1TQ	CH2-N2	3.70	1.46	1.41
1	H	109[A]	TQQ	CE2-CZ2	-3.11	1.39	1.50
1	D	109[B]	1TQ	O2-CZ2	3.01	1.43	1.36
1	D	109[A]	TQQ	CZ3-CE3	2.98	1.42	1.35
1	H	109[B]	1TQ	O2-CZ2	2.79	1.43	1.36
1	D	109[A]	TQQ	CE2-CZ2	-2.79	1.40	1.50
1	H	109[A]	TQQ	CD1-CG	-2.73	1.34	1.38
1	H	109[B]	1TQ	CB-CG	2.68	1.55	1.50
1	H	109[A]	TQQ	CZ3-CE3	2.39	1.41	1.35
1	D	109[A]	TQQ	CD2-CG	2.30	1.47	1.42
1	H	109[B]	1TQ	CH2-CZ2	-2.29	1.36	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	109[A]	TQQ	CD1-CG	-2.16	1.35	1.38
1	D	109[B]	1TQ	CH2-CZ2	-2.03	1.37	1.40

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	109[B]	1TQ	CH2-N2-CH3	-5.00	116.86	126.87
1	D	109[B]	1TQ	CH2-N2-CH3	-4.41	118.04	126.87
1	D	109[B]	1TQ	O1-CH3-N2	-3.55	120.84	125.72
1	H	109[B]	1TQ	O1-CH3-N2	-3.45	120.99	125.72
1	H	109[B]	1TQ	CZ3-CH2-CZ2	2.50	121.37	119.75
1	D	109[B]	1TQ	CZ3-CH2-CZ2	2.41	121.31	119.75

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	D	109[B]	1TQ	CZ2-CH2-N2-CH3
1	D	109[B]	1TQ	O1-CH3-N2-CH2
1	H	109[B]	1TQ	CZ2-CH2-N2-CH3
1	H	109[B]	1TQ	O1-CH3-N2-CH2
1	D	109[B]	1TQ	CZ3-CH2-N2-CH3
1	H	109[B]	1TQ	CZ3-CH2-N2-CH3

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	109[A]	TQQ	1	0
1	H	109[A]	TQQ	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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$
3	BEZ	H	2001	-	8,8,9	1.33	1 (12%)	9,9,11	0.60	0
3	BEZ	D	2002	-	8,8,9	0.97	0	9,9,11	0.60	0

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	BEZ	H	2001	-	-	0/2/2/4	0/1/1/1
3	BEZ	D	2002	-	-	0/2/2/4	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	2001	BEZ	C6-C1	2.04	1.43	1.39

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	D	109/136 (80%)	-0.11	1 (0%) 81 84	7, 16, 26, 36	1 (0%)
1	H	119/136 (87%)	0.03	1 (0%) 82 86	6, 16, 38, 51	4 (3%)
2	A	359/361 (99%)	-0.47	2 (0%) 85 89	7, 13, 25, 35	2 (0%)
2	B	360/361 (99%)	-0.01	2 (0%) 85 89	7, 18, 31, 50	2 (0%)
All	All	947/994 (95%)	-0.19	6 (0%) 85 89	6, 15, 30, 51	9 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	431	GLY	3.3
2	B	302	GLY	2.7
2	A	432	GLY	2.4
1	H	61	SER	2.3
2	B	303	ASP	2.2
1	D	95	THR	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	TQQ	D	109[A]	16/17	0.94	0.07	13,15,22,23	16
1	1TQ	D	109[B]	18/19	-	-	10,15,19,20	18
1	TQQ	H	109[A]	16/17	0.94	0.07	8,13,18,18	16
1	1TQ	H	109[B]	18/19	-	-	9,12,18,23	18

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	BEZ	H	2001	8/9	0.90	0.08	15,18,20,20	0
3	BEZ	D	2002	8/9	0.91	0.07	16,18,21,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.