



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:01 PM UTC

PDB ID : 2HXC / pdb_00002hxc
Title : Crystal structure of the benzylamine complex of aromatic amine dehydrogenase in N-semiquinone form
Authors : Roujeinikova, A.; Leys, D.
Deposited on : 2006-08-03
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

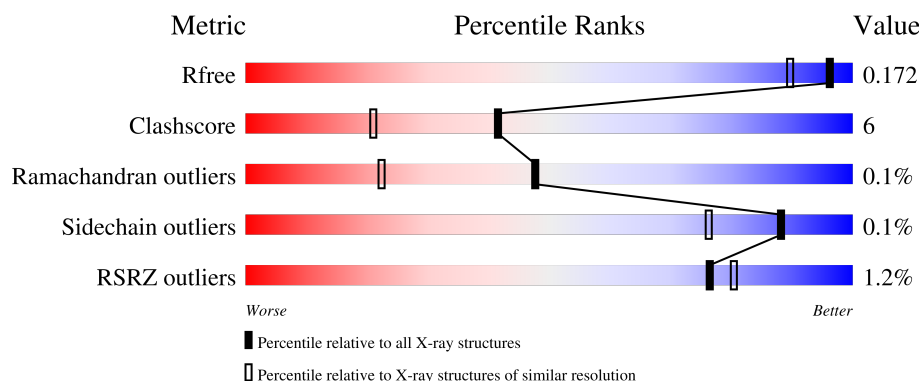
1 Overall quality at a glance

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	135	69% 12% 19%
1	H	135	73% 8% 19%
2	A	361	89% 10% 1%
2	B	361	92% 8%

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
1	TTQ	D	109[B]	-	-	X	-
3	ABN	D	1370[B]	-	-	X	-
3	ABN	H	1301	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8488 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.

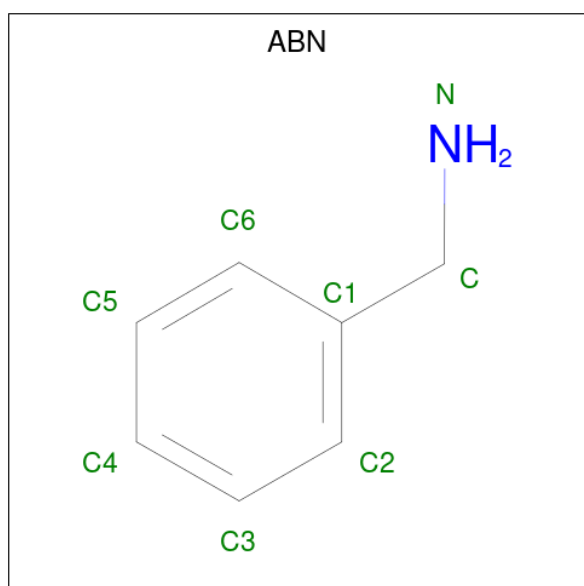
- Molecule 1 is a protein called Aromatic amine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	110	Total	C	N	O	S	0	1	0
			851	522	149	165	15			
1	H	110	Total	C	N	O	S	0	1	0
			845	516	149	165	15			

- Molecule 2 is a protein called Aromatic amine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	359	Total	C	N	O	S	0	1	0
			2797	1764	486	534	13			
2	B	360	Total	C	N	O	S	0	1	0
			2787	1758	482	534	13			

- Molecule 3 is BENZYLAMINE (CCD ID: ABN) (formula: C₇H₉N).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	127	Total O 127 127	0	0
4	H	123	Total O 123 123	0	0
4	A	504	Total O 504 504	0	0
4	B	430	Total O 430 430	0	0

- Molecule 1: Aromatic amine dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.79Å 88.59Å 79.99Å 90.00° 90.30° 90.00°	Depositor
Resolution (Å)	15.00 – 1.45 15.00 – 1.45	Depositor EDS
% Data completeness (in resolution range)	99.8 (15.00-1.45) 99.6 (15.00-1.45)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.99 (at 1.45Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.147 , 0.172 0.148 , 0.172	Depositor DCC
R_{free} test set	8689 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	11.8	Xtriage
Anisotropy	0.465	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 59.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8488	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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: TTQ, ABN

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	0.77	0/846	0.73	0/1151
1	H	0.71	0/854	0.74	0/1163
2	A	0.76	1/2864 (0.0%)	0.71	0/3884
2	B	0.74	0/2856	0.71	0/3874
All	All	0.75	1/7420 (0.0%)	0.72	0/10072

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	302	GLY	N-CA	5.69	1.49	1.44

There are no bond angle outliers.

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	D	851	0	744	26	0
1	H	845	0	746	16	0
2	A	2797	0	2734	31	0
2	B	2787	0	2704	26	0
3	D	16	0	17	8	0
3	H	8	0	9	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	504	0	0	8	0
4	B	430	0	0	5	0
4	D	127	0	0	9	0
4	H	123	0	0	4	0
All	All	8488	0	6954	91	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 (91) 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)
2:B:217:GLU:HB3	4:B:597:HOH:O	1.42	1.17
1:D:84:ASP:O	3:D:1369[A]:ABN:H2A	1.54	1.05
1:H:95[A]:THR:HG22	4:H:1322:HOH:O	1.76	0.84
2:A:124:ASN:HD21	2:A:178:GLY:H	1.30	0.79
1:D:95:THR:HG22	4:D:1391:HOH:O	1.83	0.79
2:B:124:ASN:HD21	2:B:178:GLY:H	1.30	0.77
2:B:104:ARG:HH11	2:B:106:HIS:HE1	1.32	0.76
2:A:104:ARG:HH11	2:A:106:HIS:HE1	1.34	0.74
4:H:1348:HOH:O	2:B:118:MET:SD	2.48	0.71
3:D:1370[B]:ABN:H1	4:D:1481:HOH:O	1.91	0.69
1:H:105:SER:HA	1:H:164:ASN:HD21	1.58	0.69
1:D:95:THR:HG23	4:D:1410:HOH:O	1.92	0.68
2:B:91:TYR:OH	2:B:428:HIS:HD2	1.75	0.68
2:A:91:TYR:OH	2:A:428:HIS:HD2	1.78	0.67
1:H:109:TTQ:N2	3:H:1301:ABN:N	2.42	0.67
1:D:105:SER:HA	1:D:164:ASN:HD21	1.62	0.65
1:D:109[B]:TTQ:HB2	1:D:160:TRP:NE1	2.11	0.64
1:D:84:ASP:O	3:D:1369[A]:ABN:C	2.41	0.64
1:H:109:TTQ:HB2	1:H:160:TRP:NE1	2.13	0.64
2:B:197:GLN:NE2	2:B:227:SER:H	1.97	0.63
1:D:109[A]:TTQ:HB2	1:D:160:TRP:NE1	2.14	0.61
1:H:109:TTQ:HN22	3:H:1301:ABN:H2A	1.64	0.61
2:A:263:ARG:HG3	4:A:784:HOH:O	2.01	0.61
2:A:197:GLN:NE2	2:A:227:SER:H	1.99	0.60
1:D:109[B]:TTQ:N2	3:D:1370[B]:ABN:N	2.51	0.59
2:A:308:ASP:HB2	4:A:666:HOH:O	2.03	0.58
1:D:167:SER:HB2	2:B:177:GLN:HE22	1.70	0.56
2:A:91:TYR:OH	2:A:428:HIS:CD2	2.58	0.56
2:B:335:ARG:H	2:B:386:GLN:NE2	2.03	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:335:ARG:H	2:B:386:GLN:HE22	1.51	0.56
1:H:109:TTQ:N2	3:H:1301:ABN:C	2.68	0.56
2:A:335:ARG:H	2:A:386:GLN:HE22	1.51	0.56
1:H:167:SER:HB2	2:A:177:GLN:HE22	1.70	0.56
1:H:109:TTQ:HN22	3:H:1301:ABN:C	2.19	0.55
1:H:95[A]:THR:HG23	4:H:1321:HOH:O	2.05	0.55
2:A:197:GLN:HE21	2:A:227:SER:H	1.52	0.55
2:A:74:GLU:N	4:A:912:HOH:O	2.40	0.54
1:H:71:GLU:N	4:H:1383:HOH:O	2.40	0.54
2:B:197:GLN:HE21	2:B:227:SER:H	1.55	0.54
1:D:127:HIS:HB2	2:A:76:LEU:HD22	1.89	0.54
2:A:335:ARG:H	2:A:386:GLN:NE2	2.05	0.54
1:D:117:HIS:HE1	4:D:1468:HOH:O	1.89	0.53
1:D:85:GLY:HA2	4:D:1481:HOH:O	2.08	0.53
2:B:106:HIS:HD2	2:B:418:ALA:O	1.93	0.52
2:A:106:HIS:HD2	2:A:418:ALA:O	1.94	0.51
1:D:109[B]:TTQ:CH2	3:D:1370[B]:ABN:HN2	2.23	0.50
1:H:159:ASN:HD22	2:A:177:GLN:NE2	2.10	0.49
2:A:398:ASN:HD22	2:A:416:GLU:HA	1.78	0.49
1:D:155:HIS:HE1	4:B:454:HOH:O	1.96	0.48
2:A:133:GLY:O	2:A:428:HIS:HE1	1.97	0.48
1:D:146:GLU:CD	2:B:354:LYS:HZ1	2.22	0.48
2:A:308:ASP:HB2	4:A:828:HOH:O	2.14	0.47
2:B:74:GLU:N	4:B:683:HOH:O	2.48	0.47
2:B:398:ASN:HD22	2:B:417:GLY:H	1.63	0.47
2:A:128:GLN:HE22	2:A:186:ARG:C	2.23	0.47
1:H:109:TTQ:HB2	1:H:160:TRP:HE1	1.79	0.46
1:H:155:HIS:HE1	4:A:455:HOH:O	1.97	0.46
1:H:146:GLU:CD	2:A:354:LYS:HZ1	2.25	0.45
2:B:128:GLN:HE22	2:B:186:ARG:C	2.24	0.45
2:A:143:HIS:HD2	4:A:570:HOH:O	1.99	0.45
1:D:109[B]:TTQ:HB2	1:D:160:TRP:HE1	1.79	0.45
2:B:398:ASN:HD22	2:B:416:GLU:HA	1.82	0.45
2:B:308:ASP:OD1	2:B:308:ASP:C	2.60	0.44
2:A:406[A]:GLN:H	2:A:406[A]:GLN:CD	2.26	0.44
2:A:131:ASN:H	2:A:187:GLN:HE22	1.65	0.44
2:B:216:VAL:O	2:B:217:GLU:HB2	2.16	0.44
2:B:236:ARG:HD3	4:B:700:HOH:O	2.18	0.43
2:B:275:ILE:HA	2:B:290:SER:HA	2.01	0.43
1:D:146:GLU:OE2	1:D:155:HIS:HD2	2.02	0.43
1:H:173:THR:HG21	2:B:76:LEU:HD13	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:GLY:CA	4:D:1481:HOH:O	2.66	0.42
2:A:200:SER:N	2:A:201:PRO:HA	2.34	0.42
2:A:261:GLN:NE2	4:A:628:HOH:O	2.44	0.42
1:D:109[B]:TTQ:HN22	3:D:1370[B]:ABN:C	2.32	0.42
2:A:398:ASN:HD22	2:A:417:GLY:H	1.68	0.42
1:D:159:ASN:HD22	2:B:177:GLN:NE2	2.16	0.42
4:D:1481:HOH:O	2:B:97:PHE:HZ	2.03	0.42
2:A:194:ILE:HB	2:A:208:VAL:HB	2.02	0.42
1:D:109[A]:TTQ:HB2	1:D:160:TRP:HE1	1.82	0.42
2:A:285:LYS:HD2	4:A:580:HOH:O	2.20	0.41
1:H:104:PRO:HB3	2:B:76:LEU:HD23	2.02	0.41
1:D:71:GLU:CA	4:D:1442:HOH:O	2.69	0.41
2:A:275:ILE:HA	2:A:290:SER:HA	2.02	0.41
2:A:124:ASN:ND2	2:A:178:GLY:H	2.09	0.41
1:D:109[B]:TTQ:HN22	3:D:1370[B]:ABN:H2A	1.85	0.41
2:B:416:GLU:HG3	4:B:710:HOH:O	2.19	0.41
2:B:398:ASN:ND2	2:B:417:GLY:H	2.19	0.40
1:D:81:CYS:O	1:D:142:THR:HG22	2.21	0.40
1:D:109[B]:TTQ:N2	3:D:1370[B]:ABN:C	2.85	0.40
1:D:139:GLN:NE2	4:D:1463:HOH:O	2.48	0.40
2:A:105:VAL:HG21	2:A:127:VAL:HG21	2.02	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	D	107/135 (79%)	105 (98%)	2 (2%)	0	100	100
1	H	108/135 (80%)	105 (97%)	3 (3%)	0	100	100
2	A	358/361 (99%)	346 (97%)	11 (3%)	1 (0%)	36	16
2	B	358/361 (99%)	345 (96%)	13 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	931/992 (94%)	901 (97%)	29 (3%)	1 (0%)	48 22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	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	94/112 (84%)	94 (100%)	0	100 100
1	H	96/112 (86%)	96 (100%)	0	100 100
2	A	303/304 (100%)	301 (99%)	2 (1%)	76 54
2	B	301/304 (99%)	301 (100%)	0	100 100
All	All	794/832 (95%)	792 (100%)	2 (0%)	88 74

All (2) residues with a non-rotameric sidechain are listed below:

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	117	HIS
1	D	139	GLN
1	D	143	GLN
1	D	155	HIS
1	D	164	ASN
1	H	143	GLN
1	H	155	HIS

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Mol	Chain	Res	Type
1	H	164	ASN
2	A	86	GLN
2	A	106	HIS
2	A	124	ASN
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	261	GLN
2	A	322	ASN
2	A	386	GLN
2	A	387	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	187	GLN
2	B	197	GLN
2	B	231	GLN
2	B	251	ASN
2	B	386	GLN
2	B	398	ASN
2	B	406	GLN
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 ⓘ

3 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	TTQ	D	109[A]	3,1	16,16,18	1.57	4 (25%)	16,22,26	3.46	2 (12%)
1	TTQ	H	109	1	16,17,18	1.43	3 (18%)	16,24,26	1.21	1 (6%)
1	TTQ	D	109[B]	1	16,17,18	1.37	3 (18%)	16,24,26	1.01	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
1	TTQ	D	109[A]	3,1	-	0/5/6/8	0/2/2/2
1	TTQ	H	109	1	-	0/5/6/8	0/2/2/2
1	TTQ	D	109[B]	1	-	0/5/6/8	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	109[A]	TTQ	CB-CG	3.79	1.57	1.50
1	D	109[B]	TTQ	CH2-CZ2	-3.25	1.36	1.40
1	H	109	TTQ	CD2-CG	2.92	1.48	1.44
1	H	109	TTQ	CH2-CZ2	-2.73	1.37	1.40
1	D	109[B]	TTQ	CE3-CD2	2.64	1.44	1.39
1	D	109[A]	TTQ	CD2-CG	2.60	1.48	1.44
1	H	109	TTQ	CE3-CD2	2.31	1.43	1.39
1	D	109[A]	TTQ	CE3-CD2	2.20	1.43	1.39
1	D	109[A]	TTQ	CA-N	-2.04	1.42	1.48
1	D	109[B]	TTQ	CA-N	-2.04	1.42	1.48

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	109[A]	TTQ	CZ2-CH2-N2	-11.68	101.08	120.36
1	D	109[A]	TTQ	CZ3-CH2-N2	6.45	132.74	120.11
1	H	109	TTQ	CZ3-CH2-CZ2	2.00	121.53	118.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	109[A]	TTQ	2	0
1	H	109	TTQ	6	0
1	D	109[B]	TTQ	7	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
3	ABN	D	1370[B]	-	8,8,8	0.59	0	9,9,9	0.35	0
3	ABN	H	1301	-	8,8,8	0.48	0	9,9,9	0.85	0
3	ABN	D	1369[A]	1	8,8,8	0.58	0	9,9,9	0.63	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	ABN	D	1370[B]	-	-	0/2/2/2	0/1/1/1
3	ABN	H	1301	-	-	0/2/2/2	0/1/1/1
3	ABN	D	1369[A]	1	-	0/2/2/2	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1370[B]	ABN	6	0
3	H	1301	ABN	4	0
3	D	1369[A]	ABN	2	0

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	109/135 (80%)	-0.23	1 (0%) 81 84	8, 14, 25, 35	0
1	H	109/135 (80%)	-0.06	1 (0%) 81 84	8, 15, 30, 43	1 (0%)
2	A	359/361 (99%)	-0.53	6 (1%) 69 72	6, 11, 23, 35	1 (0%)
2	B	360/361 (99%)	-0.17	3 (0%) 82 86	7, 16, 28, 38	1 (0%)
All	All	937/992 (94%)	-0.30	11 (1%) 76 80	6, 13, 27, 43	3 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	303	ASP	4.0
2	A	302	GLY	3.2
2	A	431	GLY	2.9
1	D	180	ALA	2.8
2	A	432	GLY	2.5
2	A	75	VAL	2.5
2	B	216	VAL	2.4
2	B	308	ASP	2.4
2	A	308	ASP	2.2
2	B	211	ALA	2.1
1	H	180	ALA	2.0

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	TTQ	H	109	16/17	0.95	0.06	11,13,18,32	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	TTQ	D	109[B]	16/17	0.96	0.07	12,14,17,28	12
1	TTQ	D	109[A]	15/17	0.96	0.07	11,12,14,14	11

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	ABN	H	1301	8/8	0.75	0.16	14,21,24,27	8
3	ABN	D	1370[B]	8/8	0.93	0.06	11,17,19,20	8
3	ABN	D	1369[A]	8/8	0.93	0.07	10,16,19,19	8

6.5 Other polymers [i](#)

There are no such residues in this entry.