



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:51 PM UTC

PDB ID : 2VOT / pdb_00002vot
Title : Structural and biochemical evidence for a boat-like transition state in beta-mannosidases
Authors : Tailford, L.N.; Offen, W.A.; Smith, N.; Dumon, C.; Moreland, C.; Gratien, J.; Heck, M.-P.; Stick, R.V.; Bleriot, Y.; Vasella, A.; Gilbert, H.J.; Davies, G.J.
Deposited on : 2008-02-20
Resolution : 1.95 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

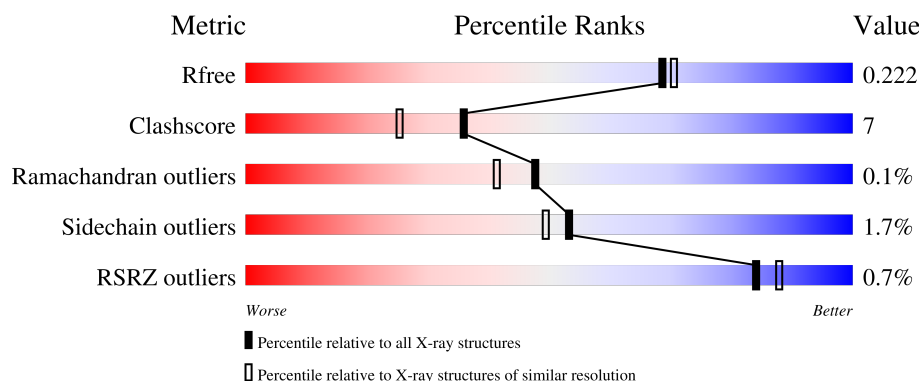
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	846	% 85% 13% ..
1	B	846	% 85% 13% ..

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
3	EDO	A	1867	-	-	X	-
3	EDO	B	1879	-	-	X	-
3	EDO	B	1898	-	-	X	-

2 Entry composition [i](#)

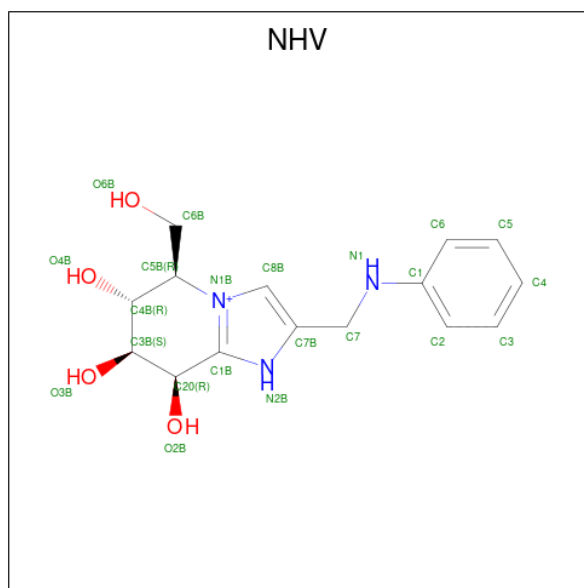
There are 6 unique types of molecules in this entry. The entry contains 15345 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BETA-MANNOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	837	Total	C	N	O	S	3	52	0
			7015	4502	1171	1308	34			
1	B	839	Total	C	N	O	S	0	37	0
			6916	4440	1149	1292	35			

- Molecule 2 is (5R,6R,7S,8R)-6,7,8-trihydroxy-5-(hydroxymethyl)-2-[(phenylamino)methyl]-5,6,7,8-tetrahydro-1H-imidazo[1,2-a]pyridin-4-ium (CCD ID: NHV) (formula: $C_{15}H_{20}N_3O_4$).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	A	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0
3	B	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0

- Molecule 4 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	5	Total Br 5 5	0	0
4	B	3	Total Br 3 3	0	0

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	5	Total 5	Cl 5	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	545	Total 545	O 545	0	0
6	B	604	Total 604	O 604	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.31Å 115.29Å 99.40Å 90.00° 113.14° 90.00°	Depositor
Resolution (Å)	20.03 – 1.95 20.03 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.03-1.95) 99.8 (20.03-1.95)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.76 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.4.0065	Depositor
R, R_{free}	0.166 , 0.221 0.166 , 0.222	Depositor DCC
R_{free} test set	6781 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	15345	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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: EDO, NHV, BR, CL

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	1.04	1/7247 (0.0%)	1.01	10/9852 (0.1%)
1	B	1.11	6/7161 (0.1%)	1.02	11/9734 (0.1%)
All	All	1.08	7/14408 (0.0%)	1.02	21/19586 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	254	ILE	CA-CB	6.69	1.59	1.54
1	B	357	ALA	CA-CB	5.97	1.62	1.53
1	B	395	TRP	CA-C	5.86	1.60	1.52
1	B	344	GLU	N-CA	5.51	1.53	1.46
1	A	207	SER	CA-C	-5.46	1.45	1.52
1	B	291	VAL	CA-CB	5.29	1.60	1.54
1	B	293	ASN	N-CA	5.17	1.50	1.46

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	425	THR	CA-C-N	-10.14	109.99	120.03
1	A	425	THR	C-N-CA	-10.14	109.99	120.03
1	A	75	GLY	N-CA-C	7.45	124.57	110.86
1	B	588[A]	LYS	N-CA-C	6.92	123.34	113.02
1	B	588[B]	LYS	N-CA-C	6.92	123.34	113.02
1	B	746	TRP	N-CA-C	-5.75	105.29	112.93
1	B	136	GLY	N-CA-C	-5.61	102.50	112.54
1	B	279	GLN	CA-C-N	-5.50	114.11	119.78
1	B	279	GLN	C-N-CA	-5.50	114.11	119.78
1	A	396	GLY	N-CA-C	5.47	122.09	115.31
1	A	521	ARG	CA-C-N	5.44	125.52	119.32
1	A	521	ARG	C-N-CA	5.44	125.52	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	200	TRP	CA-C-N	-5.39	117.46	121.61
1	A	200	TRP	C-N-CA	-5.39	117.46	121.61
1	A	38	THR	N-CA-C	5.33	117.51	109.41
1	B	425	THR	CA-C-N	-5.32	114.22	119.92
1	B	425	THR	C-N-CA	-5.32	114.22	119.92
1	B	208	GLY	N-CA-C	5.31	117.74	111.63
1	B	864	LYS	CA-C-N	-5.07	113.52	122.74
1	B	864	LYS	C-N-CA	-5.07	113.52	122.74
1	A	507	ARG	NE-CZ-NH1	-5.02	116.48	121.50

There are no chirality outliers.

There are no planarity outliers.

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	A	7015	0	6678	102	0
1	B	6916	0	6610	91	0
2	A	22	0	20	2	0
2	B	22	0	20	4	0
3	A	100	0	150	24	0
3	B	108	0	162	29	0
4	A	5	0	0	0	0
4	B	3	0	0	0	0
5	B	5	0	0	0	0
6	A	545	0	0	13	0
6	B	604	0	0	3	0
All	All	15345	0	13640	203	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 7.

All (203) 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:B:626[A]:ARG:HH22	1:B:674:ARG:HB2	1.04	1.14
1:A:575[A]:ASP:OD2	1:A:582[A]:VAL:HB	1.48	1.13
1:B:587:GLN:OE1	1:B:589[A]:SER:HB3	1.50	1.10
1:A:75:GLY:HA2	1:A:582[B]:VAL:HG13	1.44	1.00
1:B:587:GLN:NE2	1:B:593[B]:ASN:OD1	1.96	0.97
1:A:73:PHE:O	1:A:586[B]:HIS:NE2	2.00	0.94
1:B:298[A]:MET:HE3	1:B:298[A]:MET:HA	1.49	0.94
1:B:626[A]:ARG:NH2	1:B:674:ARG:HB2	1.82	0.93
1:B:202:ILE:O	3:B:1898:EDO:H11	1.70	0.92
1:A:575[A]:ASP:CG	1:A:582[A]:VAL:HB	1.96	0.90
1:B:201:GLY:CA	3:B:1898:EDO:H22	2.05	0.86
1:A:214[B]:THR:HG22	6:A:2142:HOH:O	1.76	0.85
1:B:626[A]:ARG:HH22	1:B:674:ARG:CB	1.88	0.83
1:B:695:ILE:HG23	3:B:1880:EDO:H22	1.62	0.80
1:A:525:TRP:CE2	3:A:1867:EDO:H22	2.19	0.78
1:B:341:LYS:HE3	3:B:1900:EDO:H21	1.66	0.78
1:A:575[A]:ASP:OD2	1:A:582[A]:VAL:CB	2.34	0.75
1:B:496:LEU:O	1:B:499[B]:THR:HG22	1.87	0.74
1:B:695:ILE:CG2	3:B:1880:EDO:H22	2.18	0.73
1:A:76:MET:SD	1:A:581[B]:GLU:O	2.47	0.73
1:A:328:ILE:HG23	3:A:1868:EDO:H11	1.72	0.71
1:B:201:GLY:HA2	3:B:1898:EDO:H22	1.72	0.71
1:B:464:LEU:HD22	1:B:511:HIS:HE1	1.56	0.69
1:B:787:THR:HA	1:B:808:MET:HE3	1.74	0.69
1:A:578[A]:ILE:HD13	1:A:597:ARG:HE	1.56	0.69
1:A:199:ASP:C	1:A:588[A]:LYS:HB2	2.17	0.69
1:B:525:TRP:CE2	3:B:1869:EDO:H22	2.27	0.69
1:A:76:MET:O	6:A:2026:HOH:O	2.10	0.69
1:A:284:HIS:HB3	3:A:1893:EDO:H21	1.73	0.69
1:A:470:TRP:CZ2	2:A:1865:NHV:H2	2.29	0.68
1:A:578[A]:ILE:CD1	1:A:597:ARG:HE	2.08	0.66
1:A:76:MET:SD	1:A:585[B]:ALA:HB2	2.37	0.65
1:B:201:GLY:HA3	3:B:1898:EDO:H22	1.79	0.65
1:B:470:TRP:CZ2	2:B:1868:NHV:H2	2.31	0.64
1:A:574[B]:GLU:H	1:A:574[B]:GLU:CD	2.04	0.64
1:B:35:LEU:HD23	1:B:214[B]:THR:HG22	1.80	0.64
1:A:199:ASP:HA	1:A:588[A]:LYS:HB2	1.80	0.64
1:B:33:VAL:CG1	1:B:214[A]:THR:HG23	2.27	0.64
1:B:298[A]:MET:HA	1:B:298[A]:MET:CE	2.26	0.64
1:A:570[A]:PHE:CD1	1:A:570[A]:PHE:C	2.77	0.62
1:B:33:VAL:HG13	1:B:214[A]:THR:CG2	2.28	0.62
1:A:591[B]:ILE:HG12	1:A:591[B]:ILE:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:442:VAL:HG12	3:A:1875:EDO:H11	1.82	0.62
1:B:33:VAL:HG13	1:B:214[A]:THR:HG23	1.81	0.61
1:B:406[A]:TYR:OH	1:B:418:GLN:OE1	2.15	0.61
1:A:77:ASN:HA	1:A:80:LYS:HG3	1.83	0.60
1:A:470:TRP:CH2	2:A:1865:NHV:H2	2.36	0.60
1:A:104:ARG:HG3	1:A:217:PHE:HB3	1.83	0.60
1:A:575[A]:ASP:OD2	1:A:580[A]:SER:OG	2.19	0.59
1:B:663:GLY:HA2	3:B:1901:EDO:H21	1.83	0.59
1:B:519:TRP:HE1	3:B:1879:EDO:H12	1.66	0.59
1:B:516[B]:LEU:HD13	1:B:527:THR:HG21	1.84	0.59
1:B:587:GLN:OE1	1:B:589[A]:SER:CB	2.40	0.59
1:A:464:LEU:HD22	1:A:511:HIS:HE1	1.67	0.58
1:A:102:LEU:O	1:A:142:LYS:HE3	2.04	0.58
1:A:532[A]:ASN:CG	3:A:1867:EDO:O2	2.46	0.58
1:A:70:PRO:HB2	1:A:77:ASN:ND2	2.19	0.58
1:A:578[B]:ILE:HG22	1:A:579[B]:GLU:HG2	1.86	0.58
1:A:536:TRP:CZ2	1:A:592[B]:GLY:HA2	2.39	0.57
1:B:470:TRP:CH2	2:B:1868:NHV:H2	2.38	0.57
1:A:532[A]:ASN:HB2	3:A:1867:EDO:O2	2.05	0.57
1:A:574[B]:GLU:CD	1:A:574[B]:GLU:N	2.62	0.57
1:A:575[B]:ASP:OD2	1:A:580[B]:SER:HB2	2.05	0.57
1:B:298[A]:MET:HE3	1:B:298[A]:MET:CA	2.23	0.56
1:B:85:GLU:HG3	1:B:160:ILE:HG13	1.88	0.56
1:A:708:GLU:HG2	3:A:1894:EDO:H12	1.86	0.56
1:B:695:ILE:HG23	3:B:1880:EDO:C2	2.34	0.56
2:B:1868:NHV:C6	3:B:1879:EDO:H11	2.36	0.55
3:A:1875:EDO:H22	6:A:2314:HOH:O	2.07	0.55
2:B:1868:NHV:C1	3:B:1879:EDO:H11	2.36	0.55
1:A:284:HIS:CB	3:A:1893:EDO:H21	2.37	0.55
1:A:570[A]:PHE:O	1:A:570[A]:PHE:HD1	1.90	0.55
1:A:28:ASN:HA	6:A:2002:HOH:O	2.07	0.54
1:B:519:TRP:HE1	3:B:1879:EDO:C1	2.19	0.54
3:B:1898:EDO:H11	6:B:2600:HOH:O	2.08	0.53
1:A:401:GLU:O	1:A:406[A]:TYR:HE2	1.91	0.53
1:A:740:ARG:HD2	6:A:2480:HOH:O	2.08	0.53
3:A:1868:EDO:H22	6:A:2534:HOH:O	2.08	0.53
1:A:754:ARG:CZ	1:A:783[A]:GLN:HE22	2.22	0.53
3:B:1898:EDO:C1	6:B:2600:HOH:O	2.57	0.53
1:B:536:TRP:CZ2	1:B:592[A]:GLY:HA2	2.44	0.52
1:B:126:LEU:HG	3:B:1872:EDO:H11	1.92	0.52
1:A:327:ARG:O	3:A:1881:EDO:H12	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:578:ILE:C	1:B:579[A]:GLU:HG2	2.35	0.52
1:B:424:CYS:SG	3:B:1877:EDO:H12	2.49	0.52
1:B:522:PRO:HA	1:B:525:TRP:CE2	2.45	0.51
1:B:104:ARG:HG3	1:B:217:PHE:HB3	1.93	0.51
3:A:1890:EDO:H12	6:A:2258:HOH:O	2.11	0.50
1:A:244:LEU:O	3:A:1893:EDO:H22	2.11	0.50
1:B:201:GLY:HA3	3:B:1898:EDO:C2	2.40	0.50
1:B:763:ASP:OD2	1:B:767:HIS:HB3	2.12	0.50
1:A:578[A]:ILE:HG13	1:A:579[A]:GLU:HG3	1.94	0.50
1:A:199:ASP:CA	1:A:588[A]:LYS:HB2	2.41	0.49
1:B:602:ARG:HD3	3:B:1889:EDO:H22	1.93	0.49
1:A:248[A]:GLN:OE1	1:A:278:LEU:HB2	2.13	0.49
1:A:501:LYS:O	1:A:505:SER:HB3	2.12	0.48
1:A:589[A]:SER:HB2	1:A:656:TRP:CH2	2.48	0.48
1:A:561:PHE:HE1	1:A:583[A]:MET:SD	2.37	0.48
1:B:525:TRP:CD2	3:B:1869:EDO:H22	2.48	0.48
1:A:525:TRP:CD2	3:A:1867:EDO:H22	2.48	0.48
1:A:424:CYS:SG	3:A:1878:EDO:H21	2.53	0.48
1:A:597:ARG:HD3	6:A:2397:HOH:O	2.14	0.48
1:B:380:LEU:HD13	3:B:1901:EDO:O2	2.13	0.48
1:B:464:LEU:HD22	1:B:511:HIS:CE1	2.44	0.48
1:B:731:PRO:HG2	1:B:736:LYS:HE2	1.94	0.48
1:B:199:ASP:OD2	3:B:1877:EDO:O2	2.28	0.47
1:A:570[B]:PHE:CD1	1:A:570[B]:PHE:N	2.82	0.47
1:A:33:VAL:HG13	1:A:214[B]:THR:OG1	2.14	0.47
1:B:753:ARG:NH1	1:B:783:GLN:OE1	2.48	0.47
1:A:71:ASN:OD1	1:A:71:ASN:C	2.57	0.47
1:B:536:TRP:HZ2	1:B:592[A]:GLY:HA2	1.80	0.46
1:A:211:ARG:HB3	1:A:212:PRO:CD	2.45	0.46
1:A:613:PHE:HB2	3:A:1871:EDO:H12	1.97	0.46
3:A:1890:EDO:C1	6:A:2258:HOH:O	2.63	0.46
1:A:536:TRP:HZ2	1:A:592[B]:GLY:HA2	1.78	0.46
1:B:105:ASP:HB2	1:B:219:ASP:HA	1.98	0.46
1:A:33:VAL:HG13	1:A:214[A]:THR:HG23	1.96	0.46
1:A:646:GLN:HG2	1:A:657:SER:HA	1.98	0.46
1:B:793:MET:SD	1:B:800[B]:CYS:SG	3.14	0.46
1:B:162:GLN:HA	3:B:1887:EDO:H21	1.98	0.46
1:A:75:GLY:HA2	1:A:582[B]:VAL:CG1	2.32	0.45
1:A:162:GLN:HA	3:A:1884:EDO:H22	1.97	0.45
1:B:394:ILE:HD11	1:B:405:PHE:CE2	2.51	0.45
1:B:589[B]:SER:HB2	1:B:656:TRP:CH2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:665:TRP:CZ2	3:B:1897:EDO:H21	2.52	0.45
1:A:532[A]:ASN:CG	3:A:1867:EDO:HO2	2.24	0.45
1:A:591[B]:ILE:O	1:A:591[B]:ILE:CG1	2.63	0.45
1:B:55:VAL:HA	1:B:56:PRO:C	2.41	0.45
1:B:298[A]:MET:HE3	1:B:299:PRO:HD2	1.98	0.45
1:A:270:THR:HG21	1:A:290:GLU:O	2.17	0.45
1:A:545:LEU:HD13	1:A:628:GLY:HA3	1.99	0.45
1:A:75:GLY:O	1:A:585[B]:ALA:HB3	2.16	0.45
1:B:591[B]:ILE:HG21	1:B:595:LEU:HD22	1.99	0.45
1:B:298[A]:MET:HE1	1:B:351:ASN:HD22	1.82	0.45
1:A:328:ILE:HG23	3:A:1868:EDO:C1	2.44	0.44
3:A:1866:EDO:H22	6:A:2259:HOH:O	2.16	0.44
1:B:33:VAL:CG1	1:B:214[A]:THR:CG2	2.91	0.44
1:B:194:TYR:CD1	3:B:1898:EDO:H12	2.52	0.44
1:B:199:ASP:HA	1:B:588[B]:LYS:HB3	1.99	0.44
1:A:28:ASN:HB3	6:A:2002:HOH:O	2.16	0.44
1:A:199:ASP:HA	1:A:588[A]:LYS:CB	2.48	0.44
1:A:522:PRO:HA	1:A:525:TRP:CD2	2.53	0.44
1:A:570[A]:PHE:CD1	1:A:570[A]:PHE:O	2.69	0.43
1:B:204[B]:MET:HE3	1:B:204[B]:MET:HB2	1.58	0.43
1:A:764:LYS:HD2	1:A:764:LYS:O	2.18	0.43
1:B:698:ARG:NH2	1:B:700[A]:ASP:OD2	2.49	0.43
1:A:485:MET:HE3	1:A:485:MET:HB3	1.93	0.43
1:A:701:THR:HG23	1:A:731:PRO:HA	2.01	0.43
1:B:626[A]:ARG:HG3	1:B:627:HIS:N	2.31	0.43
1:B:646:GLN:O	1:B:657:SER:HA	2.19	0.43
1:A:570[B]:PHE:CD2	1:A:586[B]:HIS:CG	3.07	0.43
1:B:626[A]:ARG:NH2	1:B:671:GLN:O	2.38	0.43
1:A:174:TYR:CD2	1:A:187:VAL:HG21	2.53	0.43
1:A:394:ILE:HD11	1:A:405:PHE:CE2	2.54	0.43
1:A:110:ILE:HB	1:A:214[B]:THR:HG23	2.01	0.42
1:B:504:ASP:OD2	1:B:507:ARG:HD2	2.19	0.42
1:B:586:HIS:HD2	1:B:652:PRO:O	2.02	0.42
1:A:80:LYS:HE2	1:A:80:LYS:HB3	1.79	0.42
1:A:99:GLU:O	1:A:103[A]:ASN:ND2	2.52	0.42
1:A:625:MET:HB2	1:A:668:LEU:HD13	2.02	0.42
1:A:105:ASP:HB2	1:A:219:ASP:HA	2.00	0.42
1:B:441:ALA:HB1	1:B:499[B]:THR:HG21	2.02	0.42
1:A:586[B]:HIS:HD2	1:A:652:PRO:O	2.02	0.42
1:B:516[B]:LEU:HD22	1:B:527:THR:HB	2.02	0.42
1:B:865:GLU:H	1:B:865:GLU:HG2	1.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:573[B]:PRO:HD2	1:A:574[B]:GLU:OE1	2.20	0.42
1:A:597:ARG:HG3	1:A:606:ILE:HD13	2.02	0.42
1:B:200:TRP:HZ3	3:B:1898:EDO:H21	1.85	0.41
1:B:519:TRP:NE1	3:B:1879:EDO:H12	2.34	0.41
1:B:95:PHE:CE1	1:B:215:LEU:HD11	2.55	0.41
1:B:625:MET:HB2	1:B:668:LEU:HD13	2.01	0.41
1:A:559:GLN:OE1	1:A:587[A]:GLN:NE2	2.29	0.41
1:A:760:ILE:HG23	3:A:1894:EDO:H22	2.01	0.41
1:B:44:GLN:HB2	1:B:89:TRP:CZ3	2.54	0.41
1:A:115:ASP:OD1	1:A:191:LYS:NZ	2.53	0.41
1:A:199:ASP:O	1:A:588[A]:LYS:HB2	2.19	0.41
1:A:798:GLY:HA2	1:A:843:SER:O	2.20	0.41
1:A:546:ASP:OD2	1:A:627:HIS:NE2	2.43	0.41
1:A:626:ARG:HD3	6:A:2445:HOH:O	2.19	0.41
1:B:118:ALA:HA	1:B:155:TYR:O	2.20	0.41
1:A:636:ARG:HG2	1:A:637:PRO:HA	2.03	0.41
1:A:682:ASN:HA	1:A:683:PRO:HD2	1.97	0.41
1:B:578:ILE:O	1:B:593[A]:ASN:ND2	2.53	0.41
1:B:591[B]:ILE:CG2	1:B:595:LEU:HD22	2.51	0.41
1:A:469:TYR:CE2	3:A:1870:EDO:H22	2.56	0.41
1:B:353:ILE:HD13	1:B:699:LEU:HD22	2.03	0.41
1:B:791:TYR:HA	1:B:803:THR:O	2.21	0.41
1:A:588[B]:LYS:NZ	6:A:2392:HOH:O	2.45	0.40
1:A:646:GLN:O	1:A:657:SER:HA	2.21	0.40
1:A:786:PRO:O	1:A:808:MET:HE1	2.21	0.40
1:B:437:VAL:HG11	1:B:491:LEU:HD11	2.03	0.40
1:B:541:PRO:HA	6:B:2409:HOH:O	2.21	0.40
1:A:100:GLU:H	1:A:100:GLU:CD	2.29	0.40
1:A:575[A]:ASP:CG	1:A:582[A]:VAL:CB	2.81	0.40
1:A:577[A]:GLN:HE21	1:A:579[A]:GLU:H	1.68	0.40
1:B:298[A]:MET:CE	1:B:298[A]:MET:CA	2.94	0.40
1:B:660:ASP:C	1:B:660:ASP:OD1	2.64	0.40
1:B:390:ASN:HD22	3:B:1884:EDO:C1	2.35	0.40
1:B:732:ALA:O	1:B:734:THR:HG23	2.22	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	874/846 (103%)	852 (98%)	21 (2%)	1 (0%)	48	41
1	B	862/846 (102%)	833 (97%)	28 (3%)	1 (0%)	48	41
All	All	1736/1692 (103%)	1685 (97%)	49 (3%)	2 (0%)	48	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	536	TRP
1	A	536	TRP

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	757/754 (100%)	742 (98%)	15 (2%)	48	43
1	B	752/754 (100%)	736 (98%)	16 (2%)	47	41
All	All	1509/1508 (100%)	1478 (98%)	31 (2%)	53	41

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	THR
1	A	80	LYS
1	A	144	VAL
1	A	319	ARG

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Mol	Chain	Res	Type
1	A	395	TRP
1	A	417	TRP
1	A	550	PRO
1	A	570[A]	PHE
1	A	570[B]	PHE
1	A	577[A]	GLN
1	A	577[B]	GLN
1	A	591[A]	ILE
1	A	591[B]	ILE
1	A	764	LYS
1	A	819	LEU
1	B	144	VAL
1	B	269	VAL
1	B	287	LEU
1	B	417	TRP
1	B	440	GLU
1	B	516[A]	LEU
1	B	516[B]	LEU
1	B	521	ARG
1	B	578	ILE
1	B	591[A]	ILE
1	B	591[B]	ILE
1	B	626[A]	ARG
1	B	626[B]	ARG
1	B	758	LYS
1	B	813	ILE
1	B	865	GLU

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

Mol	Chain	Res	Type
1	A	28	ASN
1	A	86	ASN
1	A	108	GLN
1	A	242	ASN
1	A	337	ASN
1	B	28	ASN
1	B	284	HIS
1	B	337	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 67 ligands modelled in this entry, 13 are monoatomic - leaving 54 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	$\# Z > 2$
3	EDO	B	1900	-	3,3,3	0.39	0	2,2,2	0.72	0
3	EDO	A	1894	-	3,3,3	0.50	0	2,2,2	0.34	0
3	EDO	A	1867	-	3,3,3	0.45	0	2,2,2	1.34	0
3	EDO	A	1869	-	3,3,3	0.83	0	2,2,2	0.28	0
3	EDO	A	1893	-	3,3,3	0.55	0	2,2,2	0.08	0
3	EDO	A	1880	-	3,3,3	0.40	0	2,2,2	0.70	0
3	EDO	B	1887	-	3,3,3	0.27	0	2,2,2	0.57	0
3	EDO	B	1902	-	3,3,3	0.32	0	2,2,2	1.01	0
3	EDO	B	1877	-	3,3,3	0.49	0	2,2,2	0.33	0
3	EDO	B	1889	-	3,3,3	0.65	0	2,2,2	0.13	0
3	EDO	A	1882	-	3,3,3	0.46	0	2,2,2	0.82	0
3	EDO	A	1875	-	3,3,3	0.61	0	2,2,2	0.09	0
3	EDO	A	1892	-	3,3,3	0.25	0	2,2,2	0.97	0
3	EDO	B	1872	-	3,3,3	0.34	0	2,2,2	0.54	0
3	EDO	B	1897	-	3,3,3	0.31	0	2,2,2	0.50	0
3	EDO	B	1884	-	3,3,3	0.60	0	2,2,2	0.57	0
2	NHV	B	1868	-	21,24,24	1.58	1 (4%)	23,34,34	1.77	4 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	A	1877	-	3,3,3	0.49	0	2,2,2	0.59	0
3	EDO	B	1871	-	3,3,3	0.39	0	2,2,2	0.52	0
3	EDO	B	1883	-	3,3,3	0.44	0	2,2,2	1.36	0
3	EDO	B	1882	-	3,3,3	0.51	0	2,2,2	0.19	0
3	EDO	A	1871	-	3,3,3	0.32	0	2,2,2	0.51	0
3	EDO	A	1884	-	3,3,3	0.64	0	2,2,2	0.51	0
3	EDO	A	1873	-	3,3,3	0.43	0	2,2,2	0.46	0
3	EDO	B	1870	-	3,3,3	0.48	0	2,2,2	0.43	0
3	EDO	B	1895	-	3,3,3	0.53	0	2,2,2	0.50	0
3	EDO	A	1883	-	3,3,3	0.54	0	2,2,2	0.45	0
3	EDO	B	1888	-	3,3,3	0.89	0	2,2,2	0.58	0
3	EDO	B	1869	-	3,3,3	0.44	0	2,2,2	1.02	0
3	EDO	B	1878	-	3,3,3	0.52	0	2,2,2	0.44	0
3	EDO	A	1891	-	3,3,3	0.41	0	2,2,2	0.69	0
3	EDO	B	1898	-	3,3,3	0.34	0	2,2,2	0.81	0
3	EDO	A	1879	-	3,3,3	0.63	0	2,2,2	0.37	0
3	EDO	A	1870	-	3,3,3	0.95	0	2,2,2	0.35	0
3	EDO	A	1878	-	3,3,3	0.38	0	2,2,2	0.70	0
3	EDO	A	1885	-	3,3,3	0.61	0	2,2,2	0.12	0
3	EDO	A	1881	-	3,3,3	1.13	0	2,2,2	0.50	0
3	EDO	B	1876	-	3,3,3	0.70	0	2,2,2	0.20	0
2	NHV	A	1865	-	21,24,24	1.37	2 (9%)	23,34,34	1.69	5 (21%)
3	EDO	B	1874	-	3,3,3	0.36	0	2,2,2	0.35	0
3	EDO	A	1876	-	3,3,3	0.42	0	2,2,2	0.88	0
3	EDO	A	1868	-	3,3,3	0.55	0	2,2,2	0.46	0
3	EDO	A	1890	-	3,3,3	0.57	0	2,2,2	0.48	0
3	EDO	A	1874	-	3,3,3	0.53	0	2,2,2	0.18	0
3	EDO	B	1879	-	3,3,3	0.61	0	2,2,2	0.38	0
3	EDO	A	1872	-	3,3,3	0.48	0	2,2,2	0.82	0
3	EDO	B	1886	-	3,3,3	0.55	0	2,2,2	0.21	0
3	EDO	B	1881	-	3,3,3	0.78	0	2,2,2	0.30	0
3	EDO	B	1885	-	3,3,3	0.43	0	2,2,2	0.48	0
3	EDO	B	1901	-	3,3,3	0.50	0	2,2,2	0.19	0
3	EDO	B	1880	-	3,3,3	0.72	0	2,2,2	0.43	0
3	EDO	B	1873	-	3,3,3	0.41	0	2,2,2	0.56	0
3	EDO	B	1875	-	3,3,3	0.70	0	2,2,2	0.32	0
3	EDO	A	1866	-	3,3,3	0.64	0	2,2,2	0.16	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	EDO	B	1900	-	-	1/1/1/1	-
3	EDO	A	1894	-	-	1/1/1/1	-
3	EDO	A	1867	-	-	0/1/1/1	-
3	EDO	A	1869	-	-	0/1/1/1	-
3	EDO	A	1893	-	-	1/1/1/1	-
3	EDO	A	1880	-	-	0/1/1/1	-
3	EDO	B	1887	-	-	1/1/1/1	-
3	EDO	B	1902	-	-	0/1/1/1	-
3	EDO	B	1877	-	-	0/1/1/1	-
3	EDO	B	1889	-	-	1/1/1/1	-
3	EDO	A	1882	-	-	0/1/1/1	-
3	EDO	A	1875	-	-	0/1/1/1	-
3	EDO	A	1892	-	-	0/1/1/1	-
3	EDO	B	1872	-	-	1/1/1/1	-
3	EDO	B	1897	-	-	1/1/1/1	-
3	EDO	B	1884	-	-	1/1/1/1	-
2	NHV	B	1868	-	-	0/7/27/27	0/3/3/3
3	EDO	A	1877	-	-	0/1/1/1	-
3	EDO	B	1871	-	-	1/1/1/1	-
3	EDO	B	1883	-	-	1/1/1/1	-
3	EDO	B	1882	-	-	0/1/1/1	-
3	EDO	A	1871	-	-	0/1/1/1	-
3	EDO	A	1884	-	-	0/1/1/1	-
3	EDO	A	1873	-	-	0/1/1/1	-
3	EDO	B	1870	-	-	0/1/1/1	-
3	EDO	B	1895	-	-	0/1/1/1	-
3	EDO	A	1883	-	-	0/1/1/1	-
3	EDO	B	1888	-	-	0/1/1/1	-
3	EDO	B	1869	-	-	0/1/1/1	-
3	EDO	B	1878	-	-	1/1/1/1	-
3	EDO	A	1891	-	-	0/1/1/1	-
3	EDO	B	1898	-	-	0/1/1/1	-
3	EDO	A	1879	-	-	1/1/1/1	-
3	EDO	A	1870	-	-	1/1/1/1	-
3	EDO	A	1878	-	-	0/1/1/1	-
3	EDO	A	1885	-	-	1/1/1/1	-
3	EDO	A	1881	-	-	0/1/1/1	-
3	EDO	B	1876	-	-	0/1/1/1	-
2	NHV	A	1865	-	-	0/7/27/27	0/3/3/3
3	EDO	B	1874	-	-	0/1/1/1	-
3	EDO	A	1876	-	-	0/1/1/1	-
3	EDO	A	1868	-	-	0/1/1/1	-
3	EDO	A	1890	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	1874	-	-	0/1/1/1	-
3	EDO	B	1879	-	-	0/1/1/1	-
3	EDO	A	1872	-	-	0/1/1/1	-
3	EDO	B	1886	-	-	1/1/1/1	-
3	EDO	B	1881	-	-	0/1/1/1	-
3	EDO	B	1885	-	-	0/1/1/1	-
3	EDO	B	1901	-	-	0/1/1/1	-
3	EDO	B	1880	-	-	0/1/1/1	-
3	EDO	B	1873	-	-	0/1/1/1	-
3	EDO	B	1875	-	-	0/1/1/1	-
3	EDO	A	1866	-	-	1/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1868	NHV	C8B-N1B	-5.36	1.33	1.39
2	A	1865	NHV	C8B-N1B	-4.07	1.35	1.39
2	A	1865	NHV	C5-C6	2.08	1.42	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1868	NHV	C7-C7B-N2B	5.36	126.74	121.10
2	A	1865	NHV	C7-C7B-N2B	4.67	126.01	121.10
2	B	1868	NHV	O2B-C20-C3B	4.12	119.10	110.53
2	A	1865	NHV	O2B-C20-C3B	3.89	118.61	110.53
2	A	1865	NHV	C8B-C7B-N2B	2.50	108.92	105.39
2	A	1865	NHV	C3-C2-C1	2.45	122.56	119.73
2	B	1868	NHV	C4B-C3B-C20	2.38	113.98	110.22
2	B	1868	NHV	C3-C2-C1	2.32	122.40	119.73
2	A	1865	NHV	C4B-C5B-N1B	-2.09	102.34	110.20

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1884	EDO	O1-C1-C2-O2
3	B	1900	EDO	O1-C1-C2-O2
3	B	1878	EDO	O1-C1-C2-O2
3	B	1883	EDO	O1-C1-C2-O2
3	B	1887	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	B	1889	EDO	O1-C1-C2-O2
3	A	1866	EDO	O1-C1-C2-O2
3	A	1890	EDO	O1-C1-C2-O2
3	B	1886	EDO	O1-C1-C2-O2
3	B	1897	EDO	O1-C1-C2-O2
3	A	1885	EDO	O1-C1-C2-O2
3	A	1894	EDO	O1-C1-C2-O2
3	B	1872	EDO	O1-C1-C2-O2
3	A	1870	EDO	O1-C1-C2-O2
3	B	1871	EDO	O1-C1-C2-O2
3	A	1879	EDO	O1-C1-C2-O2
3	A	1893	EDO	O1-C1-C2-O2

There are no ring outliers.

26 monomers are involved in 57 short contacts:

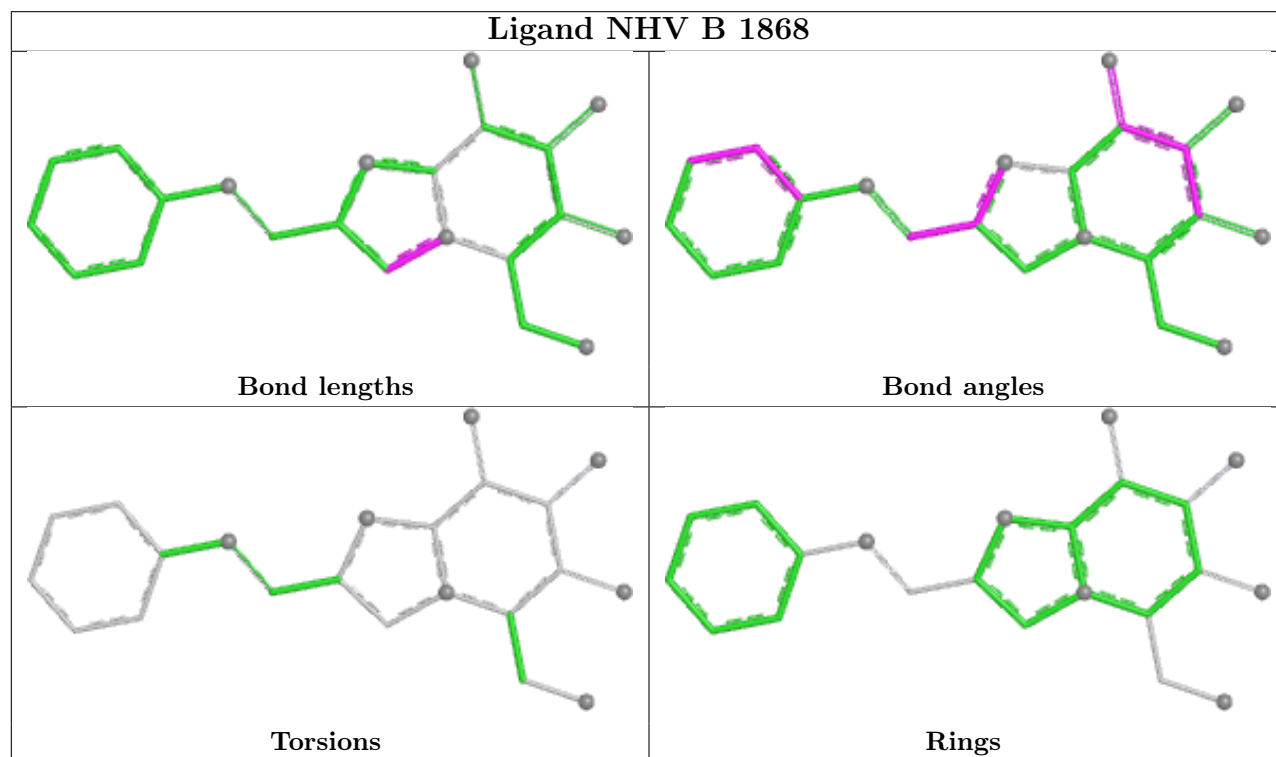
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1900	EDO	1	0
3	A	1894	EDO	2	0
3	A	1867	EDO	6	0
3	A	1893	EDO	3	0
3	B	1887	EDO	1	0
3	B	1877	EDO	2	0
3	B	1889	EDO	1	0
3	A	1875	EDO	2	0
3	B	1872	EDO	1	0
3	B	1897	EDO	1	0
3	B	1884	EDO	1	0
2	B	1868	NHV	4	0
3	A	1871	EDO	1	0
3	A	1884	EDO	1	0
3	B	1869	EDO	2	0
3	B	1898	EDO	9	0
3	A	1870	EDO	1	0
3	A	1878	EDO	1	0
3	A	1881	EDO	1	0
2	A	1865	NHV	2	0
3	A	1868	EDO	3	0
3	A	1890	EDO	2	0
3	B	1879	EDO	5	0
3	B	1901	EDO	2	0
3	B	1880	EDO	3	0

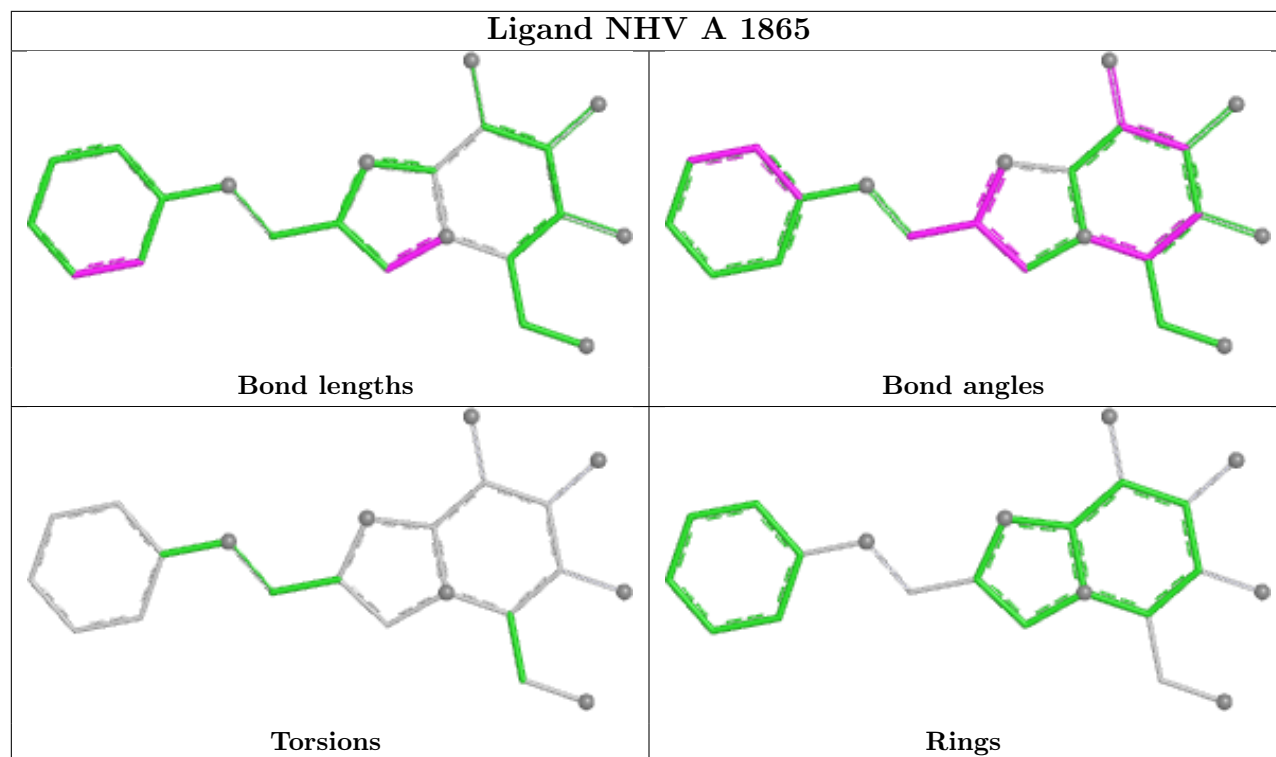
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1866	EDO	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	A	837/846 (98%)	-0.29	6 (0%) 84 88	9, 25, 44, 57	51 (6%)
1	B	839/846 (99%)	-0.37	6 (0%) 84 88	7, 24, 42, 52	35 (4%)
All	All	1676/1692 (99%)	-0.33	12 (0%) 84 88	7, 25, 43, 57	86 (5%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	74[A]	TYR	5.0
1	B	592[A]	GLY	2.7
1	B	580	SER	2.6
1	A	797	ASP	2.4
1	A	471	GLY	2.4
1	B	590[A]	SER	2.4
1	A	590[A]	SER	2.4
1	B	578	ILE	2.3
1	B	849	GLY	2.3
1	A	848	LYS	2.2
1	B	797	ASP	2.1
1	A	796	THR	2.1

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 oligosaccharides 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. 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
3	EDO	B	1883	4/4	0.75	0.17	38,39,41,42	0
3	EDO	B	1882	4/4	0.76	0.20	54,55,55,59	0
3	EDO	B	1886	4/4	0.79	0.13	48,50,50,51	0
3	EDO	A	1879	4/4	0.80	0.14	47,50,50,51	0
3	EDO	B	1879	4/4	0.82	0.13	40,44,45,48	0
3	EDO	B	1872	4/4	0.82	0.13	46,47,47,49	0
3	EDO	B	1884	4/4	0.84	0.16	36,37,37,37	0
3	EDO	A	1891	4/4	0.84	0.13	54,55,56,56	0
3	EDO	A	1884	4/4	0.85	0.18	35,39,41,47	0
3	EDO	A	1870	4/4	0.85	0.17	34,36,40,40	0
3	EDO	A	1892	4/4	0.85	0.13	44,45,45,48	0
3	EDO	B	1871	4/4	0.85	0.13	45,46,47,51	0
3	EDO	A	1867	4/4	0.85	0.15	34,36,36,37	0
3	EDO	B	1900	4/4	0.85	0.17	55,55,55,56	0
3	EDO	A	1885	4/4	0.86	0.11	45,47,48,50	0
3	EDO	A	1883	4/4	0.87	0.12	42,43,44,48	0
3	EDO	A	1893	4/4	0.87	0.21	45,46,47,48	0
3	EDO	B	1878	4/4	0.87	0.11	29,40,45,49	0
3	EDO	A	1881	4/4	0.88	0.14	25,30,32,33	0
3	EDO	A	1873	4/4	0.88	0.11	41,48,48,51	0
3	EDO	A	1890	4/4	0.88	0.10	33,33,35,38	0
3	EDO	A	1894	4/4	0.88	0.10	48,49,50,51	0
3	EDO	B	1880	4/4	0.88	0.14	31,32,33,35	0
3	EDO	B	1885	4/4	0.89	0.12	53,54,55,57	0
3	EDO	A	1876	4/4	0.90	0.09	40,44,45,47	0
3	EDO	B	1887	4/4	0.90	0.27	37,43,44,48	0
3	EDO	B	1889	4/4	0.90	0.10	38,41,42,44	0
3	EDO	A	1875	4/4	0.90	0.16	30,31,35,44	0
5	CL	B	1903	1/1	0.90	0.15	48,48,48,48	1
3	EDO	A	1880	4/4	0.91	0.09	47,48,48,49	0
3	EDO	A	1877	4/4	0.91	0.12	43,44,47,49	0
3	EDO	A	1878	4/4	0.91	0.13	31,33,36,40	0
3	EDO	A	1874	4/4	0.91	0.09	36,39,41,42	0
3	EDO	B	1897	4/4	0.92	0.09	27,33,33,35	0
3	EDO	A	1871	4/4	0.93	0.09	38,39,41,43	0
5	CL	B	1896	1/1	0.93	0.18	34,34,34,34	1
3	EDO	B	1888	4/4	0.93	0.08	27,29,29,34	0

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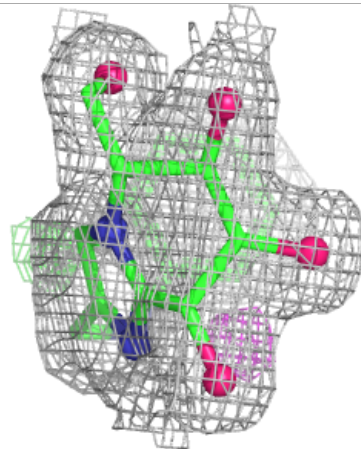
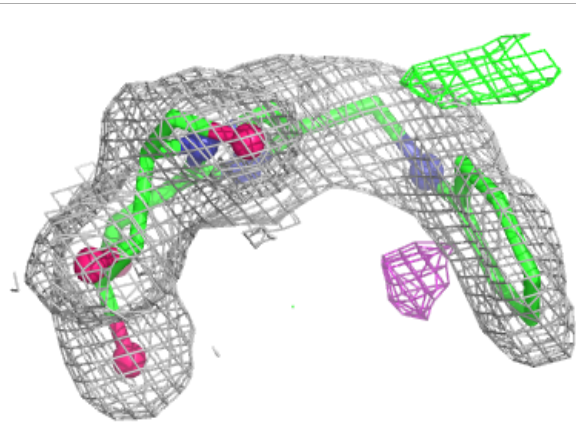
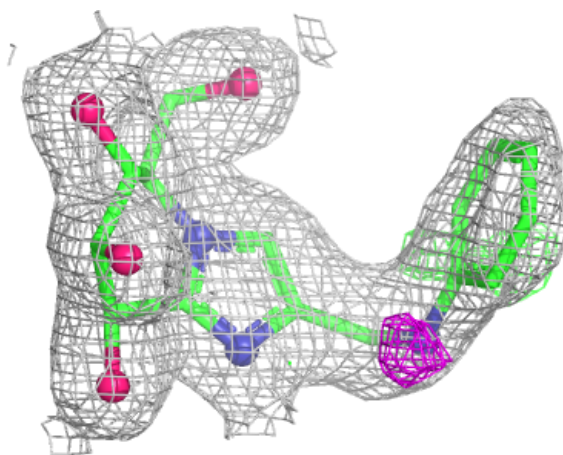
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	B	1901	4/4	0.94	0.13	35,41,44,44	0
3	EDO	B	1902	4/4	0.94	0.08	43,44,45,46	0
3	EDO	B	1874	4/4	0.94	0.09	47,48,50,51	0
5	CL	B	1899	1/1	0.94	0.08	30,30,30,30	1
3	EDO	A	1868	4/4	0.94	0.08	22,31,34,42	0
3	EDO	B	1869	4/4	0.95	0.07	27,28,31,37	0
3	EDO	B	1877	4/4	0.95	0.08	28,30,32,34	0
3	EDO	B	1870	4/4	0.95	0.06	23,26,26,28	0
3	EDO	A	1869	4/4	0.95	0.06	23,26,29,33	0
3	EDO	A	1866	4/4	0.95	0.08	26,28,30,31	0
3	EDO	B	1881	4/4	0.95	0.09	17,28,30,33	0
3	EDO	B	1873	4/4	0.95	0.07	32,34,37,39	0
3	EDO	A	1872	4/4	0.96	0.06	24,24,24,29	0
3	EDO	B	1895	4/4	0.96	0.06	28,28,29,31	0
2	NHV	A	1865	22/22	0.96	0.07	19,22,39,39	0
3	EDO	B	1898	4/4	0.96	0.12	25,26,26,26	0
2	NHV	B	1868	22/22	0.96	0.07	17,20,41,42	0
4	BR	A	1887	1/1	0.97	0.08	22,22,22,22	1
4	BR	A	1889	1/1	0.97	0.10	28,28,28,28	1
3	EDO	B	1875	4/4	0.97	0.07	25,27,30,32	0
3	EDO	B	1876	4/4	0.97	0.05	18,20,21,22	0
3	EDO	A	1882	4/4	0.97	0.10	12,32,37,45	0
4	BR	B	1891	1/1	0.98	0.13	23,23,23,23	1
4	BR	A	1888	1/1	0.98	0.07	28,28,28,28	1
4	BR	B	1892	1/1	0.99	0.11	20,20,20,20	1
5	CL	B	1893	1/1	0.99	0.06	31,31,31,31	0
5	CL	B	1894	1/1	0.99	0.05	33,33,33,33	1
4	BR	A	1900	1/1	0.99	0.10	21,21,21,21	1
4	BR	B	1890	1/1	0.99	0.04	28,28,28,28	1
4	BR	A	1886	1/1	0.99	0.06	26,26,26,26	1

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

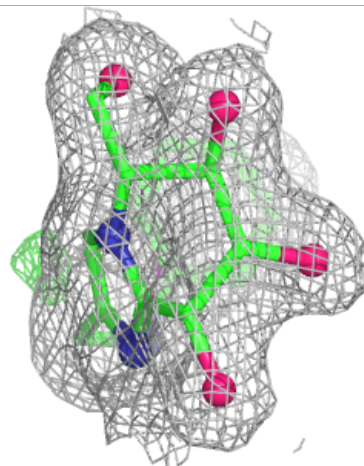
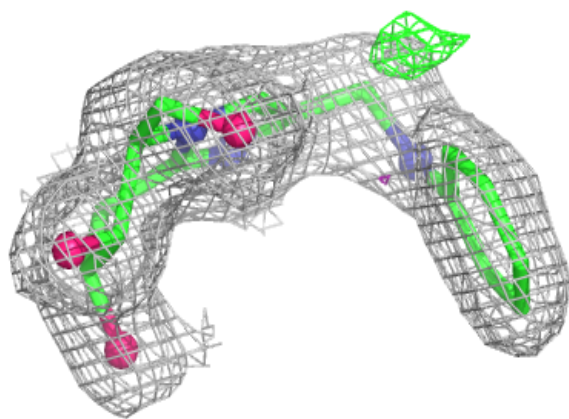
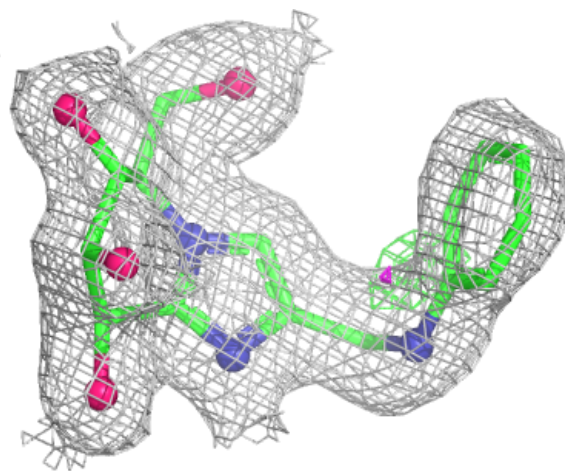
Electron density around NHV A 1865:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around NHV B 1868:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.