



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 01:26 PM EDT

PDB ID : 1TMO / pdb_00001tmo
Title : TRIMETHYLAMINE N-OXIDE REDUCTASE FROM SHEWANELLA MASSILIA
Authors : Czjzek, M.; Dos Santos, J.P.; Giordano, G.; Mejean, V.
Deposited on : 1998-08-03
Resolution : 2.50 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

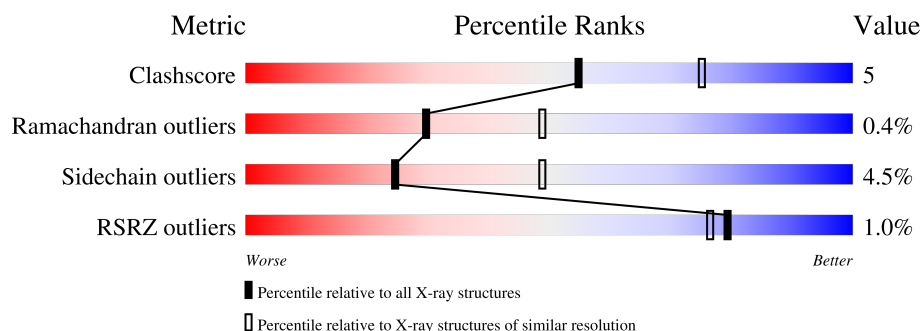
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	829	

2 Entry composition [i](#)

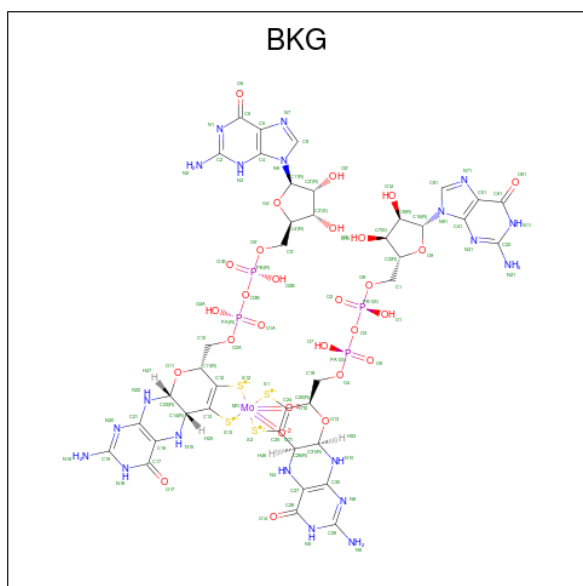
There are 3 unique types of molecules in this entry. The entry contains 6857 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 TRIMETHYLAMINE N-OXIDE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	794	Total	C	N	O	S	0	0	0
			6255	3972	1080	1167	36			

- Molecule 2 is bis-molybdopterin guanine dinucleotide dioxo molybdenum (CCD ID: BKG) (formula: $C_{40}H_{48}MoN_{20}O_{28}P_4S_4$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Mo	N	O	P	S	0
			97	40	1	20	28	4	4	0

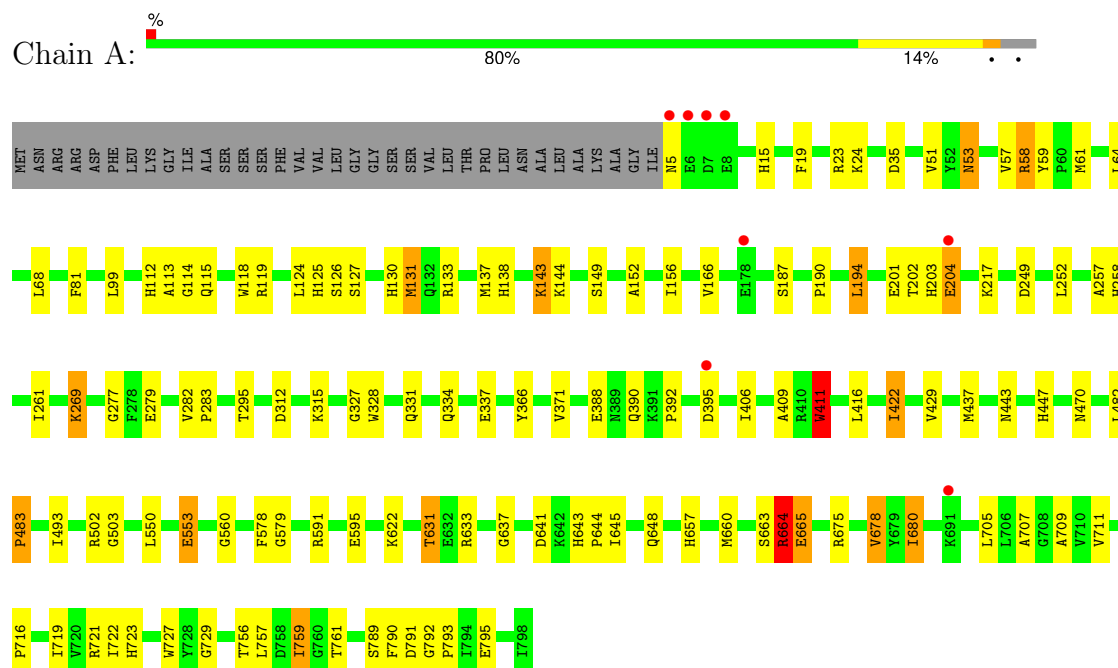
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	505	Total	O	0	0
			505	505		

3 Residue-property plots

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: TRIMETHYLAMINE N-OXIDE REDUCTASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	123.32Å 140.70Å 59.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.5 (20.00-2.50) 84.5 (20.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.41Å)	Xtriage
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.182 , 0.247 0.209 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	14.5	Xtriage
Anisotropy	0.533	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 66.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6857	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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: BKG

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	0.59	0/6422	1.03	42/8702 (0.5%)

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	127	SER	N-CA-C	10.40	122.20	111.07
1	A	789	SER	N-CA-C	7.97	120.05	111.36
1	A	644	PRO	N-CA-C	7.20	122.86	113.86
1	A	792	GLY	CA-C-N	7.09	127.05	120.03
1	A	792	GLY	C-N-CA	7.09	127.05	120.03
1	A	390	GLN	N-CA-C	7.08	119.93	109.24
1	A	791	ASP	N-CA-C	6.99	121.41	113.02
1	A	643	HIS	CA-C-N	6.71	126.34	119.56
1	A	643	HIS	C-N-CA	6.71	126.34	119.56
1	A	35	ASP	N-CA-C	6.69	119.43	111.33
1	A	756	THR	N-CA-C	6.36	119.14	110.55
1	A	201	GLU	N-CA-C	6.11	118.60	109.25
1	A	482	LEU	CA-C-N	6.11	127.03	120.13
1	A	482	LEU	C-N-CA	6.11	127.03	120.13
1	A	645	ILE	N-CA-C	5.93	117.88	108.87
1	A	125	HIS	N-CA-C	5.89	120.47	113.16
1	A	166	VAL	N-CA-C	5.89	118.42	111.05
1	A	58	ARG	N-CA-C	5.89	117.78	111.36
1	A	249	ASP	N-CA-C	5.73	118.26	111.33
1	A	447	HIS	N-CA-C	5.64	119.88	112.89
1	A	709	ALA	N-CA-C	5.63	118.41	109.96
1	A	118	TRP	N-CA-C	5.60	118.30	109.39
1	A	443	ASN	CA-C-N	5.56	125.66	119.32
1	A	443	ASN	C-N-CA	5.56	125.66	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	790	PHE	N-CA-C	5.55	120.16	113.17
1	A	680	ILE	N-CA-C	5.47	116.36	108.48
1	A	53	ASN	CA-C-N	5.46	125.55	119.87
1	A	53	ASN	C-N-CA	5.46	125.55	119.87
1	A	156	ILE	N-CA-C	5.42	117.33	111.58
1	A	793	PRO	N-CA-C	5.40	119.52	111.41
1	A	727	TRP	N-CA-C	5.27	117.99	110.24
1	A	579	GLY	N-CA-C	5.26	117.86	110.38
1	A	411	TRP	N-CA-C	5.26	117.70	111.33
1	A	483	PRO	N-CA-C	5.26	119.12	110.55
1	A	19	PHE	N-CA-C	5.20	116.56	109.18
1	A	729	GLY	CA-C-N	5.19	125.09	119.85
1	A	729	GLY	C-N-CA	5.19	125.09	119.85
1	A	560	GLY	N-CA-C	5.18	119.40	112.77
1	A	277	GLY	N-CA-C	5.15	122.42	115.32
1	A	366	TYR	N-CA-C	5.13	116.77	108.41
1	A	406	ILE	N-CA-C	5.07	113.47	107.84
1	A	409	ALA	N-CA-C	5.00	119.25	113.20

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	6255	0	6069	55	0
2	A	97	0	0	8	0
3	A	505	0	0	6	0
All	All	6857	0	6069	59	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 5.

All (59) 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:A:392:PRO:HG2	1:A:395:ASP:HB2	1.57	0.87
1:A:113:ALA:HB1	1:A:131:MET:HE1	1.57	0.85
1:A:5:ASN:N	1:A:24:LYS:HG3	1.92	0.84
2:A:799:BKG:OT2	2:A:799:BKG:MO	1.57	0.75
1:A:657:HIS:CE1	2:A:799:BKG:S2	2.81	0.73
2:A:799:BKG:OT2	2:A:799:BKG:OT1	2.07	0.72
1:A:663:SER:HB3	3:A:826:HOH:O	1.91	0.70
2:A:799:BKG:MO	2:A:799:BKG:OT1	1.64	0.69
1:A:422:ILE:HG22	1:A:429:VAL:HB	1.74	0.68
1:A:58:ARG:HG3	1:A:58:ARG:HH11	1.58	0.68
1:A:664:ARG:HH11	1:A:664:ARG:HB3	1.60	0.66
1:A:204:GLU:HG3	1:A:502:ARG:HH12	1.64	0.62
1:A:411:TRP:HE1	1:A:437:MET:HE1	1.70	0.57
1:A:190:PRO:O	1:A:194:LEU:HB2	2.06	0.56
1:A:550:LEU:O	1:A:553:GLU:HG3	2.05	0.56
1:A:112:HIS:HB3	1:A:437:MET:HG2	1.88	0.55
1:A:15:HIS:HE1	1:A:119:ARG:O	1.89	0.54
1:A:657:HIS:NE2	2:A:799:BKG:S2	2.81	0.54
1:A:633:ARG:HB3	1:A:637:GLY:HA3	1.90	0.54
1:A:187:SER:O	1:A:327:GLY:HA3	2.08	0.54
1:A:678:VAL:HG22	1:A:707:ALA:CB	2.39	0.53
1:A:114:GLY:HA3	1:A:143:LYS:HE3	1.88	0.53
1:A:665:GLU:H	1:A:665:GLU:CD	2.17	0.53
1:A:657:HIS:HE1	2:A:799:BKG:S2	2.32	0.53
1:A:57:VAL:HG12	1:A:483:PRO:HB3	1.89	0.52
1:A:61:MET:HE1	1:A:795:GLU:HG2	1.91	0.52
1:A:678:VAL:HG22	1:A:707:ALA:HB1	1.94	0.50
1:A:757:LEU:HD12	1:A:757:LEU:O	2.11	0.50
1:A:328:TRP:O	1:A:331:GLN:HG3	2.12	0.49
2:A:799:BKG:OT2	2:A:799:BKG:S2	2.70	0.49
1:A:723:HIS:HD2	3:A:1216:HOH:O	1.95	0.49
1:A:334:GLN:HE22	1:A:631:THR:H	1.61	0.48
1:A:124:LEU:O	1:A:130:HIS:HE1	1.96	0.48
1:A:152:ALA:HB1	1:A:337:GLU:HG3	1.97	0.47
1:A:648:GLN:NE2	1:A:721:ARG:HE	2.12	0.47
1:A:503:GLY:HA2	1:A:578:PHE:CD2	2.50	0.46
1:A:257:ALA:O	1:A:261:ILE:HG12	2.16	0.46
1:A:282:VAL:HB	1:A:283:PRO:HD3	1.97	0.46
1:A:759:ILE:HD13	1:A:759:ILE:HA	1.84	0.46
1:A:716:PRO:HG2	1:A:719:ILE:HD12	1.97	0.45
1:A:133:ARG:O	1:A:137:MET:HG3	2.16	0.45
1:A:149:SER:OG	2:A:799:BKG:OT1	2.36	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:GLU:HG3	1:A:502:ARG:HH22	1.82	0.44
1:A:502:ARG:NE	3:A:955:HOH:O	2.51	0.43
1:A:312:ASP:OD1	1:A:315:LYS:HE3	2.18	0.43
1:A:678:VAL:HG13	1:A:722:ILE:HG12	1.99	0.43
1:A:622:LYS:HG2	3:A:948:HOH:O	2.19	0.42
1:A:258:HIS:CE1	1:A:295:THR:HG22	2.54	0.42
1:A:58:ARG:NH1	3:A:833:HOH:O	2.52	0.42
1:A:680:ILE:O	1:A:711:VAL:HA	2.20	0.42
1:A:143:LYS:HE2	1:A:144:LYS:H	1.84	0.42
1:A:58:ARG:HH11	1:A:58:ARG:CG	2.28	0.42
1:A:202:THR:O	1:A:203:HIS:HB2	2.19	0.41
1:A:269:LYS:HB2	1:A:269:LYS:NZ	2.36	0.41
1:A:675:ARG:N	1:A:675:ARG:HD2	2.35	0.41
1:A:99:LEU:HB3	1:A:138:HIS:CE1	2.56	0.41
1:A:53:ASN:HD21	1:A:470:ASN:HD21	1.70	0.40
1:A:641:ASP:HB2	3:A:1029:HOH:O	2.21	0.40
1:A:705:LEU:C	1:A:705:LEU:HD12	2.47	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	792/829 (96%)	749 (95%)	40 (5%)	3 (0%)	30	49

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	664	ARG
1	A	660	MET
1	A	388	GLU

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	665/692 (96%)	635 (96%)	30 (4%)	24 49

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ARG
1	A	51	VAL
1	A	59	TYR
1	A	64	LEU
1	A	68	LEU
1	A	81	PHE
1	A	115	GLN
1	A	126	SER
1	A	131	MET
1	A	143	LYS
1	A	194	LEU
1	A	204	GLU
1	A	217	LYS
1	A	252	LEU
1	A	269	LYS
1	A	279	GLU
1	A	371	VAL
1	A	411	TRP
1	A	416	LEU
1	A	422	ILE
1	A	493	ILE
1	A	553	GLU
1	A	591	ARG
1	A	595	GLU
1	A	631	THR
1	A	664	ARG
1	A	665	GLU
1	A	678	VAL
1	A	759	ILE
1	A	761	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	HIS
1	A	71	HIS
1	A	77	GLN
1	A	125	HIS
1	A	130	HIS
1	A	199	ASN
1	A	203	HIS
1	A	234	GLN
1	A	245	ASN
1	A	334	GLN
1	A	425	ASN
1	A	442	ASN
1	A	443	ASN
1	A	446	ASN
1	A	452	ASN
1	A	470	ASN
1	A	597	ASN
1	A	635	HIS
1	A	643	HIS
1	A	648	GLN
1	A	651	HIS
1	A	723	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	BKG	A	799	1	100,110,110	2.43	23 (23%)	116,181,181	2.94	33 (28%)

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	BKG	A	799	1	-	4/44/160/160	0/14/14/14

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	799	BKG	OT2-MO	-9.60	1.57	1.73
2	A	799	BKG	O14-C28	8.26	1.39	1.23
2	A	799	BKG	O17-C17	8.09	1.38	1.23
2	A	799	BKG	PA-O3B	5.92	1.65	1.59
2	A	799	BKG	C23-C14	5.74	1.58	1.53
2	A	799	BKG	OT1-MO	-5.30	1.64	1.73
2	A	799	BKG	PB-O3B	5.03	1.64	1.59
2	A	799	BKG	C31-C26	5.02	1.57	1.53
2	A	799	BKG	C16-C21	4.93	1.47	1.38
2	A	799	BKG	C27-C30	4.56	1.46	1.38
2	A	799	BKG	C4-N9	4.09	1.41	1.37
2	A	799	BKG	C18-C20	3.56	1.56	1.51
2	A	799	BKG	PA1-O3	3.28	1.63	1.59
2	A	799	BKG	C2-N1	2.83	1.40	1.33
2	A	799	BKG	C10-C11	2.80	1.55	1.51
2	A	799	BKG	MO-S1	2.79	2.80	2.52
2	A	799	BKG	C6-N1	2.76	1.42	1.38
2	A	799	BKG	PB1-O3	2.46	1.62	1.59
2	A	799	BKG	O11-C23	-2.46	1.40	1.43
2	A	799	BKG	O13-C20	-2.16	1.39	1.43
2	A	799	BKG	O13-C31	-2.10	1.40	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	799	BKG	O9-C15	2.01	1.46	1.42
2	A	799	BKG	C26-N4	2.01	1.48	1.46

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	799	BKG	C24-C25-S2	-12.64	112.98	120.15
2	A	799	BKG	C13-C12-S12	-12.15	113.26	120.15
2	A	799	BKG	C25-C24-S1	-10.42	114.24	120.15
2	A	799	BKG	O11-C11-C12	8.78	121.49	109.09
2	A	799	BKG	O13-C31-N10	8.14	115.99	108.61
2	A	799	BKG	C23-C14-N15	7.43	115.17	107.87
2	A	799	BKG	C12-C13-S13	-7.28	116.02	120.15
2	A	799	BKG	O3A-C10-C11	5.67	122.57	108.58
2	A	799	BKG	C29-N8-C30	5.30	122.71	113.36
2	A	799	BKG	C19-N20-C21	5.18	122.49	113.36
2	A	799	BKG	O4-C18-C20	4.17	118.88	108.58
2	A	799	BKG	O14-C28-C27	-3.72	118.30	127.26
2	A	799	BKG	C31-C26-N4	3.64	111.44	107.87
2	A	799	BKG	O13-C31-C26	-3.60	106.56	108.96
2	A	799	BKG	O11-C23-C14	3.55	111.33	108.96
2	A	799	BKG	C27-C28-N5	3.26	121.08	112.13
2	A	799	BKG	C16-C17-N18	3.21	120.95	112.13
2	A	799	BKG	O17-C17-C16	-3.17	119.61	127.26
2	A	799	BKG	C19-N18-C17	-3.06	119.57	125.11
2	A	799	BKG	O7-PA1-O3	2.64	114.42	107.27
2	A	799	BKG	N6-C29-N5	2.49	122.02	116.76
2	A	799	BKG	O3-PB1-O2	-2.49	103.20	110.70
2	A	799	BKG	O9-C15-N91	2.47	113.96	108.36
2	A	799	BKG	O4'-C1'-N9	2.46	113.94	108.36
2	A	799	BKG	C28-C27-N4	2.43	122.89	116.27
2	A	799	BKG	O9-C15-C9	-2.33	101.64	106.62
2	A	799	BKG	N2-C2-N3	2.32	121.66	116.76
2	A	799	BKG	C29-N5-C28	-2.17	121.17	125.11
2	A	799	BKG	C8-N9-C4	-2.15	104.86	106.67
2	A	799	BKG	O4'-C1'-C2'	-2.14	102.03	106.62
2	A	799	BKG	C6-N1-C2	2.05	125.26	119.24
2	A	799	BKG	N5-C29-N8	-2.04	119.59	123.32
2	A	799	BKG	C3-O9-C15	-2.00	105.05	109.47

There are no chirality outliers.

All (4) torsion outliers are listed below:

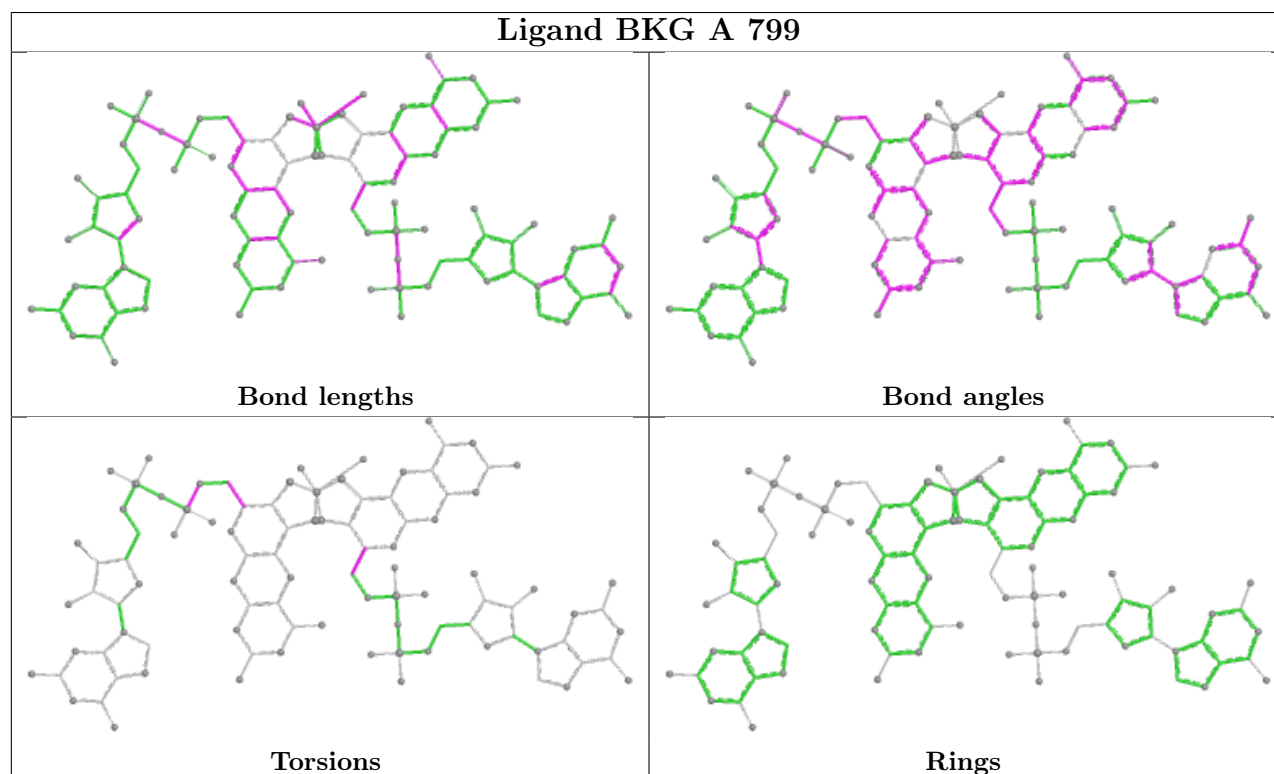
Mol	Chain	Res	Type	Atoms
2	A	799	BKG	O4-C18-C20-O13
2	A	799	BKG	O4-C18-C20-C24
2	A	799	BKG	C18-O4-PA1-O5
2	A	799	BKG	O3A-C10-C11-C12

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	799	BKG	8	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

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 [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	794/829 (95%)	0.17	8 (1%) 79 76	2, 11, 29, 64	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	ASN	4.0
1	A	395	ASP	3.8
1	A	6	GLU	2.5
1	A	204	GLU	2.4
1	A	178	GLU	2.1
1	A	691	LYS	2.1
1	A	8	GLU	2.1
1	A	7	ASP	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](#)

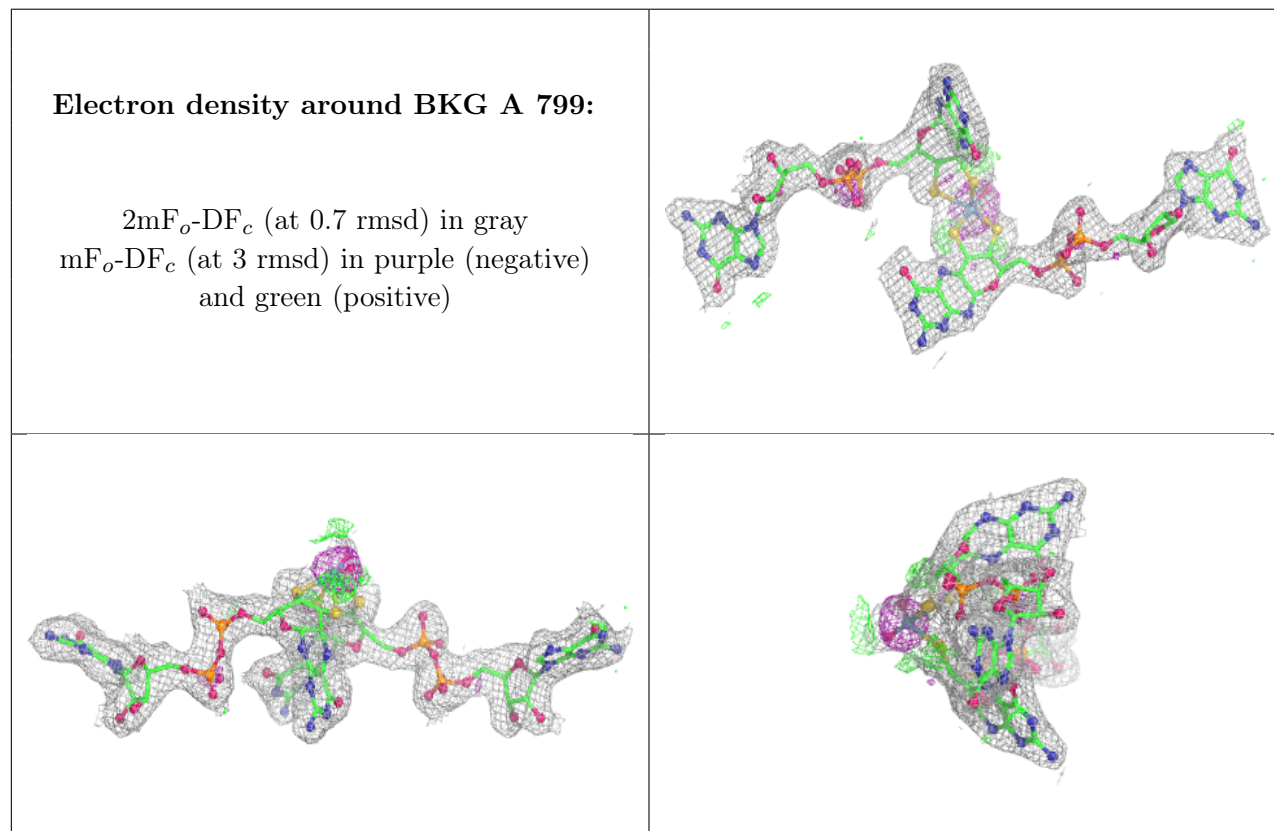
There are no oligosaccharides in this entry.

6.4 Ligands [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
2	BKG	A	799	97/97	0.92	0.10	1,11,16,24	0

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.



6.5 Other polymers [i](#)

There are no such residues in this entry.