



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:13 PM UTC

PDB ID : 1NZY / pdb_00001nzy
Title : 4-CHLOROBENZOYL COENZYME A DEHALOGENASE FROM PSEUDOMONAS SP. STRAIN CBS-3
Authors : Benning, M.M.; Holden, H.M.
Deposited on : 1996-04-02
Resolution : 1.80 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

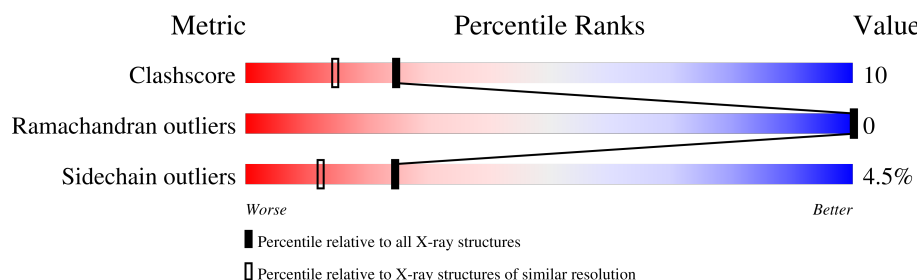
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	269	81% 17% .
1	C	269	72% 25% .
2	B	269	77% 19% . .

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7053 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 4-CHLOROBENZOYL COENZYME A DEHALOGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	1	0
			2097	1325	379	377	16			
1	C	268	Total	C	N	O	S	0	1	0
			2086	1319	375	376	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	85	ALA	GLY	conflict	PIR A42560
C	85	ALA	GLY	conflict	PIR A42560

- Molecule 2 is a protein called 4-CHLOROBENZOYL COENZYME A DEHALOGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	269	Total	C	N	O	S	0	1	0
			2097	1325	378	378	16			

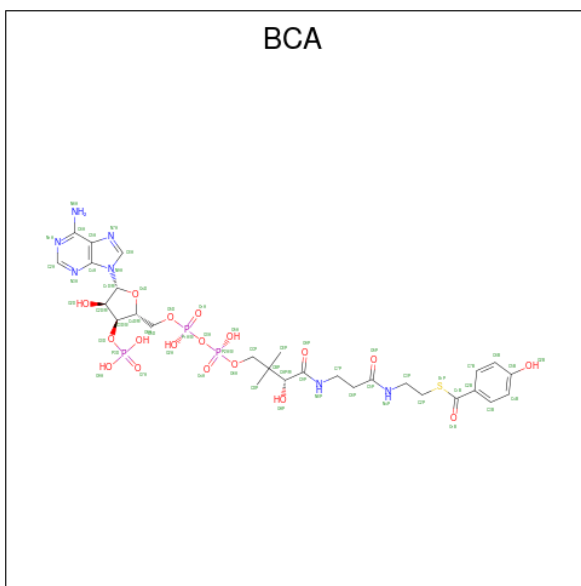
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	85	ALA	GLY	conflict	PIR A42560
B	168	ASP	ASN	conflict	PIR A42560

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

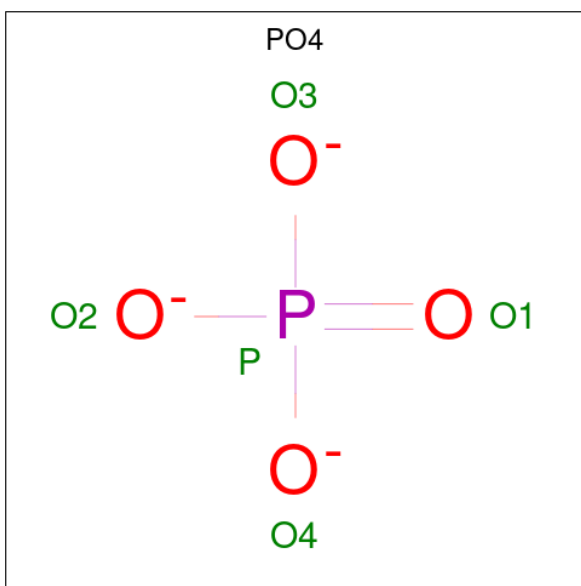
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		
3	C	1	Total	Ca	0	0
			1	1		

- Molecule 4 is 4-HYDROXYBENZOYL COENZYME A (CCD ID: BCA) (formula: $C_{28}H_{40}N_7O_{18}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			53	28	7	15	2	1		
4	B	1	Total	C	N	O	P	S	0	0
			53	28	7	15	2	1		
4	C	1	Total	C	N	O	P	S	0	0
			53	28	7	15	2	1		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	240	Total	O	0	0
			240	240		
7	B	201	Total	O	0	0
			201	201		
7	C	157	Total	O	0	0
			157	157		

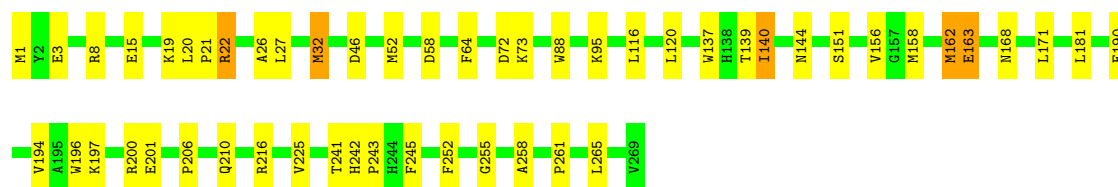
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

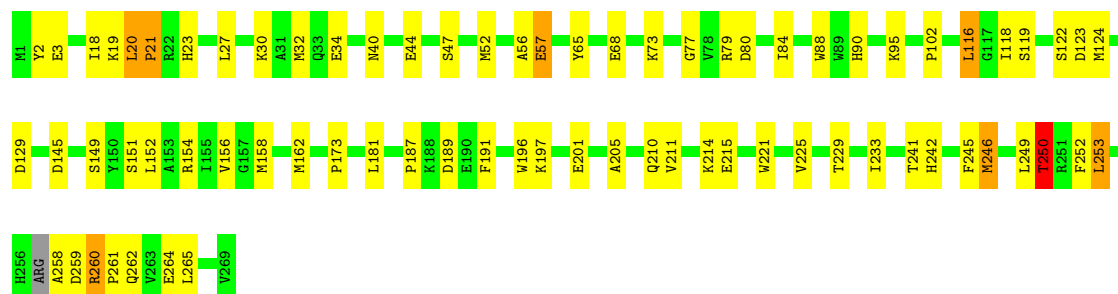
• Molecule 1: 4-CHLOROBENZOYL COENZYME A DEHALOGENASE

Chain A: 



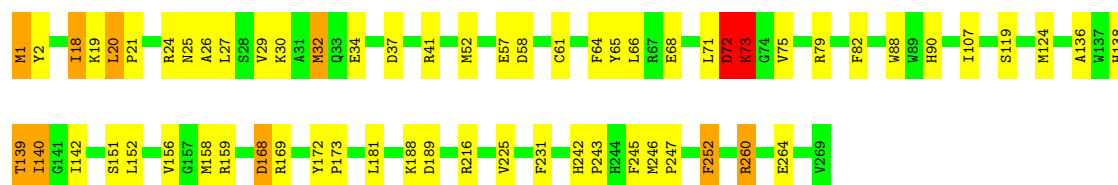
• Molecule 1: 4-CHLOROBENZOYL COENZYME A DEHALOGENASE

Chain C: 



• Molecule 2: 4-CHLOROBENZOYL COENZYME A DEHALOGENASE

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.80Å 102.40Å 90.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.188 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7053	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

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: EDO, BCA, CA, PO4

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	1.30	1/2149 (0.0%)	1.61	16/2911 (0.5%)
1	C	1.24	1/2137 (0.0%)	1.56	24/2894 (0.8%)
2	B	1.25	1/2149 (0.0%)	1.60	24/2911 (0.8%)
All	All	1.27	3/6435 (0.0%)	1.59	64/8716 (0.7%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	139	THR	C-N	-5.40	1.28	1.34
1	C	162	MET	N-CA	-5.39	1.39	1.46
2	B	189	ASP	CG-OD1	5.18	1.35	1.25

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	151	SER	N-CA-C	10.96	123.31	111.36
2	B	73	LYS	N-CA-C	10.68	123.95	110.61
1	A	245	PHE	N-CA-C	10.07	121.94	110.97
2	B	151	SER	N-CA-C	9.87	122.12	111.36
2	B	90	HIS	CA-CB-CG	-8.21	105.59	113.80
1	A	46	ASP	CA-CB-CG	-7.94	104.66	112.60
1	C	151	SER	N-CA-C	7.63	121.84	112.23
2	B	82	PHE	CA-CB-CG	-7.47	106.33	113.80
1	C	156	VAL	N-CA-C	7.36	119.97	112.90
2	B	158	MET	N-CA-C	7.15	119.16	111.36
2	B	72	ASP	CA-CB-CG	6.84	119.44	112.60
1	A	156	VAL	N-CA-C	6.82	119.58	112.83
1	C	158	MET	N-CA-C	6.71	118.59	111.28
2	B	225	VAL	N-CA-C	6.70	118.25	110.62
2	B	18	ILE	N-CA-C	-6.68	100.95	109.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	200	ARG	NE-CZ-NH2	-6.68	113.19	119.20
2	B	245	PHE	N-CA-C	6.58	124.81	110.80
2	B	216	ARG	NE-CZ-NH1	6.44	127.94	121.50
2	B	260	ARG	N-CA-C	6.38	115.65	108.25
1	C	116	LEU	N-CA-C	-6.38	104.41	111.36
1	C	225	VAL	N-CA-C	6.37	117.05	110.36
1	C	261	PRO	N-CA-CB	6.36	108.97	103.31
2	B	138	HIS	CA-CB-CG	-6.35	107.45	113.80
1	C	152	LEU	N-CA-C	6.31	117.85	110.97
1	C	242	HIS	CA-CB-CG	6.30	120.10	113.80
1	A	19	LYS	N-CA-CB	-6.02	99.68	110.82
1	C	259	ASP	CA-CB-CG	6.01	118.61	112.60
1	A	19	LYS	N-CA-C	5.97	121.15	107.48
1	A	168	ASN	CA-CB-CG	-5.92	106.68	112.60
1	C	19	LYS	N-CA-C	5.82	120.80	107.48
1	A	158	MET	N-CA-C	5.78	117.58	111.28
1	A	255	GLY	N-CA-C	5.71	122.96	115.36
1	A	252	PHE	CA-CB-CG	-5.66	108.14	113.80
1	C	173	PRO	N-CA-C	5.66	122.19	113.75
1	C	19	LYS	N-CA-CB	-5.63	100.40	110.82
1	C	21	PRO	CA-C-N	-5.58	112.24	120.28
1	C	21	PRO	C-N-CA	-5.58	112.24	120.28
2	B	152	LEU	N-CA-C	5.57	117.03	111.07
1	A	26	ALA	N-CA-C	-5.57	102.64	110.50
1	A	58	ASP	N-CA-C	5.52	120.00	112.88
2	B	52	MET	N-CA-CB	5.46	119.17	110.65
1	C	191	PHE	N-CA-C	5.41	122.33	110.80
2	B	26	ALA	N-CA-C	-5.39	103.37	110.43
1	C	245	PHE	N-CA-C	5.31	122.11	110.80
1	A	261	PRO	CB-CA-C	-5.30	104.03	110.98
2	B	168	ASP	N-CA-CB	5.29	120.61	112.41
1	A	265	LEU	CB-CA-C	-5.27	104.95	110.17
2	B	169	ARG	N-CA-CB	-5.21	102.91	110.36
1	C	252	PHE	O-C-N	5.19	127.42	122.07
2	B	252	PHE	CA-CB-CG	-5.18	108.62	113.80
1	C	18	ILE	N-CA-C	-5.17	102.01	108.84
2	B	19	LYS	N-CA-CB	-5.17	101.25	110.82
2	B	231	PHE	CA-CB-CG	-5.14	108.66	113.80
1	C	250	THR	CB-CA-C	5.13	119.26	110.74
1	C	189	ASP	N-CA-C	5.11	116.85	111.28
1	C	245	PHE	CA-CB-CG	-5.10	108.70	113.80
2	B	57	GLU	N-CA-C	5.10	117.18	108.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	VAL	N-CA-C	5.09	115.70	110.36
1	C	21	PRO	N-CA-CB	5.07	109.04	103.26
2	B	20	LEU	CA-C-N	5.04	125.32	119.47
2	B	20	LEU	C-N-CA	5.04	125.32	119.47
2	B	156	VAL	N-CA-C	5.03	117.73	112.90
1	C	118	ILE	O-C-N	5.03	126.75	121.87
1	C	3	GLU	CB-CA-C	-5.02	100.64	109.07

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	2097	0	2086	34	0
1	C	2086	0	2071	56	0
2	B	2097	0	2084	41	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	53	0	36	1	0
4	B	53	0	36	1	0
4	C	53	0	35	2	0
5	B	5	0	0	0	0
6	B	8	0	12	0	0
7	A	240	0	0	4	0
7	B	201	0	0	1	0
7	C	157	0	0	2	0
All	All	7053	0	6360	125	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 10.

All (125) 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:139:THR:HG21	1:C:258:ALA:HA	1.46	0.97
1:C:27:LEU:HD13	1:C:32:MET:HG3	1.62	0.81
1:C:27:LEU:HD13	1:C:32:MET:CG	2.11	0.80
1:A:22:ARG:HD3	7:A:279:HOH:O	1.82	0.79
2:B:139:THR:CG2	1:C:258:ALA:HA	2.12	0.79
2:B:27:LEU:HD13	2:B:32:MET:HG3	1.64	0.78
1:C:40:ASN:O	1:C:44:GLU:HG3	1.86	0.75
2:B:27:LEU:HD13	2:B:32:MET:CG	2.18	0.72
1:C:65:TYR:CD2	1:C:68:GLU:HG3	2.24	0.72
1:C:246:MET:HE3	1:C:249:LEU:HB3	1.70	0.71
1:C:80:ASP:O	1:C:84:ILE:HD12	1.94	0.67
1:C:30:LYS:O	1:C:30:LYS:HD2	1.95	0.67
1:C:246:MET:O	1:C:250:THR:HB	1.95	0.66
2:B:65:TYR:CD2	2:B:68:GLU:HG3	2.31	0.66
2:B:32:MET:HE3	2:B:88:TRP:CB	2.26	0.65
1:A:21:PRO:HD2	7:A:279:HOH:O	1.96	0.65
1:A:171:LEU:HD21	1:A:181:LEU:HD12	1.80	0.63
1:A:171:LEU:CD2	1:A:181:LEU:HD12	2.28	0.63
2:B:32:MET:HE3	2:B:88:TRP:HB2	1.81	0.63
1:C:260:ARG:HH11	1:C:260:ARG:HG3	1.62	0.63
1:A:140:ILE:O	1:A:140:ILE:HD13	1.99	0.62
1:C:260:ARG:HG3	1:C:260:ARG:NH1	2.15	0.61
1:C:246:MET:HE3	1:C:246:MET:HA	1.84	0.60
2:B:32:MET:CE	2:B:88:TRP:HB2	2.32	0.59
1:C:21:PRO:HD2	7:C:374:HOH:O	2.03	0.59
1:C:30:LYS:HE2	1:C:34:GLU:OE2	2.02	0.59
1:C:40:ASN:OD1	1:C:95:LYS:HE2	2.03	0.59
2:B:20:LEU:N	2:B:21:PRO:HD3	2.17	0.58
1:A:242:HIS:ND1	1:A:243:PRO:HD2	2.18	0.57
2:B:139:THR:HG21	1:C:258:ALA:CA	2.29	0.57
1:A:27:LEU:HG	1:A:32:MET:HG3	1.86	0.57
1:A:32:MET:CE	1:A:88:TRP:HB2	2.35	0.56
1:A:241:THR:OG1	1:C:79:ARG:HD2	2.07	0.55
1:C:260:ARG:HH11	1:C:260:ARG:CG	2.19	0.55
1:A:32:MET:HE3	1:A:88:TRP:HB2	1.89	0.55
1:C:249:LEU:HG	1:C:253:LEU:HD22	1.89	0.54
1:C:20:LEU:N	1:C:21:PRO:HD3	2.24	0.53
1:C:57:GLU:OE1	1:C:57:GLU:HA	2.08	0.53
2:B:140:ILE:HG21	2:B:142:ILE:HD12	1.91	0.53
1:A:8:ARG:HH21	1:A:196:TRP:NE1	2.07	0.52
1:C:2:TYR:HB3	1:C:34:GLU:OE1	2.09	0.51
1:A:20:LEU:N	1:A:21:PRO:HD3	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:ASP:O	1:A:73:LYS:HG3	2.11	0.51
1:C:27:LEU:HD13	1:C:32:MET:HG2	1.91	0.51
2:B:140:ILE:CG2	2:B:142:ILE:HD12	2.41	0.51
2:B:29:VAL:HA	2:B:32:MET:HE2	1.92	0.51
2:B:172:TYR:HB3	2:B:173:PRO:HD2	1.93	0.50
2:B:72:ASP:OD1	2:B:73:LYS:HG2	2.12	0.50
1:C:102:PRO:HG3	1:C:210:GLN:OE1	2.11	0.49
1:C:56:ALA:O	1:C:57:GLU:HB2	2.11	0.49
1:A:95:LYS:NZ	7:A:488:HOH:O	2.45	0.49
1:C:2:TYR:CD1	1:C:34:GLU:HB3	2.47	0.49
2:B:88:TRP:HE3	7:B:1082:HOH:O	1.96	0.49
2:B:1:MET:HG2	2:B:2:TYR:N	2.27	0.48
1:C:181:LEU:HD23	1:C:181:LEU:O	2.14	0.48
2:B:20:LEU:N	2:B:21:PRO:CD	2.77	0.48
1:C:229:THR:O	1:C:233:ILE:HG13	2.13	0.47
1:A:72:ASP:C	1:A:73:LYS:HG3	2.39	0.47
1:C:32:MET:CE	1:C:88:TRP:HB2	2.44	0.47
1:C:73:LYS:HB2	1:C:77:GLY:HA3	1.95	0.47
1:C:30:LYS:HD2	1:C:30:LYS:C	2.40	0.47
2:B:65:TYR:CD2	2:B:68:GLU:CG	2.97	0.47
1:C:211:VAL:O	1:C:215:GLU:HG3	2.15	0.47
1:A:162:MET:HB3	1:A:162:MET:HE2	1.54	0.46
1:A:242:HIS:CG	1:A:243:PRO:HD2	2.50	0.46
1:A:64:PHE:HB2	4:A:272:BCA:C2B	2.46	0.45
1:C:249:LEU:O	1:C:253:LEU:HB2	2.17	0.45
2:B:20:LEU:HD12	2:B:20:LEU:HA	1.68	0.45
2:B:24:ARG:O	2:B:25:ASN:HB2	2.16	0.45
1:C:154:ARG:HG2	1:C:221:TRP:CE2	2.52	0.45
1:A:216:ARG:HB3	1:C:149:SER:CB	2.47	0.45
2:B:32:MET:HE2	2:B:32:MET:HB2	1.82	0.44
2:B:66:LEU:HD22	2:B:140:ILE:HD12	2.00	0.44
1:A:190:GLU:O	1:A:194:VAL:HG23	2.18	0.44
2:B:124:MET:HB3	2:B:124:MET:HE3	1.54	0.44
1:A:163:GLU:OE1	7:A:453:HOH:O	2.21	0.44
1:A:197:LYS:O	1:A:201:GLU:HG3	2.18	0.44
2:B:136:ALA:N	2:B:168:ASP:OD1	2.48	0.44
2:B:242:HIS:ND1	2:B:243:PRO:HD2	2.32	0.44
2:B:64:PHE:HB2	4:B:272:BCA:C2B	2.48	0.44
1:A:27:LEU:CG	1:A:32:MET:HG3	2.47	0.44
2:B:107:ILE:HD11	2:B:119[B]:SER:HB3	2.00	0.44
1:A:27:LEU:CD2	1:A:32:MET:HG3	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:ARG:NH2	1:A:196:TRP:CE2	2.86	0.43
1:C:260:ARG:NH1	1:C:260:ARG:CG	2.80	0.43
1:A:137:TRP:CD1	1:A:144:ASN:HA	2.52	0.43
1:A:171:LEU:CD2	1:A:181:LEU:CD1	2.94	0.43
2:B:58:ASP:CG	2:B:188:LYS:HZ1	2.25	0.43
1:C:205:ALA:HB3	1:C:210:GLN:NE2	2.33	0.43
1:C:123:ASP:OD1	1:C:214:LYS:HE3	2.19	0.43
2:B:30:LYS:HE3	2:B:34:GLU:OE2	2.19	0.43
1:C:197:LYS:O	1:C:201:GLU:HG3	2.18	0.43
1:A:206:PRO:O	1:A:210:GLN:HG3	2.19	0.43
1:A:32:MET:HE2	1:A:32:MET:HB2	1.63	0.42
2:B:252:PHE:CD1	2:B:252:PHE:C	2.97	0.42
1:C:129:ASP:OD2	1:C:187:PRO:HA	2.19	0.42
2:B:27:LEU:HD13	2:B:32:MET:HG2	1.98	0.42
2:B:58:ASP:OD1	2:B:188:LYS:NZ	2.52	0.42
2:B:140:ILE:HG23	2:B:140:ILE:O	2.18	0.42
1:C:253:LEU:HD12	1:C:253:LEU:HA	1.65	0.42
1:C:124:MET:HB3	1:C:124:MET:HE2	1.68	0.42
1:A:258:ALA:N	4:C:272:BCA:O9P	2.46	0.42
2:B:140:ILE:CG2	2:B:142:ILE:CD1	2.97	0.42
1:C:249:LEU:O	1:C:250:THR:C	2.62	0.42
2:B:1:MET:CG	2:B:2:TYR:N	2.82	0.41
2:B:159:ARG:NH2	1:C:122:SER:O	2.51	0.41
1:C:116:LEU:O	1:C:119[A]:SER:OG	2.25	0.41
2:B:72:ASP:C	2:B:73:LYS:HG2	2.45	0.41
1:C:90:HIS:HE1	1:C:145:ASP:OD2	2.03	0.41
1:C:197:LYS:HD3	7:C:370:HOH:O	2.19	0.41
1:C:265:LEU:HD12	1:C:265:LEU:HA	1.81	0.41
2:B:246:MET:N	2:B:247:PRO:CD	2.83	0.41
1:C:32:MET:HE2	1:C:32:MET:HB2	1.70	0.41
2:B:18:ILE:HD13	2:B:61:CYS:HB2	2.03	0.41
1:C:23:HIS:HA	4:C:272:BCA:O4D	2.21	0.41
1:C:52:MET:HE1	1:C:196:TRP:CD1	2.56	0.41
1:C:246:MET:HA	1:C:246:MET:CE	2.50	0.41
1:A:32:MET:HE3	1:A:88:TRP:CB	2.51	0.41
2:B:37:ASP:O	2:B:41:ARG:HG2	2.21	0.41
1:C:20:LEU:N	1:C:21:PRO:CD	2.81	0.41
1:A:116:LEU:O	1:A:120:LEU:HG	2.21	0.41
1:C:95:LYS:HD2	1:C:95:LYS:HA	1.93	0.40
1:A:20:LEU:HD12	1:A:20:LEU:HA	1.91	0.40
1:C:246:MET:CE	1:C:246:MET:CA	2.99	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:GLU:HG3	1:A:52:MET:HE2	2.04	0.40

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	A	268/269 (100%)	256 (96%)	12 (4%)	0	100	100
1	C	265/269 (98%)	250 (94%)	15 (6%)	0	100	100
2	B	268/269 (100%)	257 (96%)	11 (4%)	0	100	100
All	All	801/807 (99%)	763 (95%)	38 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	215/214 (100%)	208 (97%)	7 (3%)	33	21
1	C	214/214 (100%)	204 (95%)	10 (5%)	23	11
2	B	215/214 (100%)	203 (94%)	12 (6%)	19	8
All	All	644/642 (100%)	615 (96%)	29 (4%)	24	12

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

Mol	Chain	Res	Type
1	A	1	MET
1	A	3	GLU
1	A	22	ARG
1	A	32	MET
1	A	140	ILE
1	A	162	MET
1	A	163	GLU
2	B	1	MET
2	B	32	MET
2	B	71	LEU
2	B	72	ASP
2	B	73	LYS
2	B	75	VAL
2	B	79	ARG
2	B	139	THR
2	B	140	ILE
2	B	181	LEU
2	B	260	ARG
2	B	264	GLU
1	C	20	LEU
1	C	47	SER
1	C	57	GLU
1	C	241	THR
1	C	246	MET
1	C	250	THR
1	C	253	LEU
1	C	260	ARG
1	C	262	GLN
1	C	264	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
2	B	33	GLN
2	B	235	ASN
1	C	33	GLN
1	C	235	ASN
1	C	262	GLN

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 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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	$\# Z > 2$
6	EDO	B	900	-	3,3,3	0.36	0	2,2,2	0.76	0
4	BCA	C	272	-	54,56,60	1.83	7 (12%)	76,82,89	1.75	12 (15%)
6	EDO	B	901	-	3,3,3	0.40	0	2,2,2	0.23	0
4	BCA	A	272	-	54,56,60	2.09	11 (20%)	76,82,89	1.38	5 (6%)
4	BCA	B	272	-	54,56,60	1.81	6 (11%)	76,82,89	1.83	12 (15%)
5	PO4	B	902	-	4,4,4	1.16	0	6,6,6	0.79	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
6	EDO	B	900	-	-	0/1/1/1	-
4	BCA	C	272	-	-	1/50/66/71	0/4/4/4
6	EDO	B	901	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	BCA	A	272	-	-	1/50/66/71	0/4/4/4
4	BCA	B	272	-	-	6/50/66/71	0/4/4/4

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	272	BCA	P2A-O3A	7.07	1.67	1.59
4	B	272	BCA	C9P-N8P	6.72	1.49	1.33
4	A	272	BCA	C5P-N4P	6.49	1.48	1.33
4	C	272	BCA	C9P-N8P	6.40	1.48	1.33
4	A	272	BCA	P1A-O3A	-6.25	1.52	1.59
4	A	272	BCA	C9P-N8P	6.04	1.47	1.33
4	C	272	BCA	C5P-N4P	5.87	1.47	1.33
4	B	272	BCA	C5P-N4P	5.38	1.46	1.33
4	C	272	BCA	P1A-O3A	-5.11	1.54	1.59
4	C	272	BCA	P2A-O3A	4.53	1.64	1.59
4	A	272	BCA	C7B-C2B	4.06	1.45	1.39
4	A	272	BCA	C6B-C5B	3.37	1.45	1.39
4	A	272	BCA	C4B-C3B	3.31	1.44	1.38
4	A	272	BCA	C7B-C6B	3.30	1.44	1.38
4	A	272	BCA	O2B-C5B	3.00	1.43	1.37
4	B	272	BCA	C4B-C5B	2.82	1.44	1.39
4	A	272	BCA	C2B-C1B	-2.80	1.44	1.49
4	C	272	BCA	C6B-C5B	2.75	1.44	1.39
4	A	272	BCA	C4B-C5B	2.58	1.43	1.39
4	B	272	BCA	C1B-S1P	2.41	1.81	1.76
4	C	272	BCA	C3B-C2B	2.25	1.42	1.39
4	C	272	BCA	C2A-N1A	2.16	1.37	1.33
4	B	272	BCA	C2A-N1A	2.07	1.37	1.33
4	A	272	BCA	C2A-N1A	2.02	1.37	1.33

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	272	BCA	C3P-N4P-C5P	-7.54	108.78	122.82
4	B	272	BCA	C2P-S1P-C1B	6.90	107.98	99.85
4	C	272	BCA	C3P-N4P-C5P	-6.70	110.35	122.82
4	C	272	BCA	O6A-P2A-O4A	-6.26	84.11	108.94
4	A	272	BCA	CEP-CBP-CAP	6.12	119.20	108.77
4	B	272	BCA	C7P-N8P-C9P	-5.64	112.41	122.55
4	A	272	BCA	C7P-N8P-C9P	-5.21	113.18	122.55
4	C	272	BCA	C7P-N8P-C9P	-4.94	113.67	122.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	272	BCA	C3P-N4P-C5P	-4.75	113.99	122.82
4	B	272	BCA	C6P-C5P-N4P	-4.47	108.19	116.34
4	C	272	BCA	C2P-S1P-C1B	3.49	103.97	99.85
4	B	272	BCA	O5P-C5P-N4P	-3.20	116.75	123.03
4	C	272	BCA	CDP-CBP-CAP	3.19	114.21	108.77
4	B	272	BCA	O6A-CCP-CBP	-3.06	105.63	110.55
4	C	272	BCA	C6P-C5P-N4P	-3.04	110.81	116.34
4	C	272	BCA	C2D-C1D-N9A	-2.87	106.16	113.30
4	C	272	BCA	CAP-C9P-N8P	-2.87	111.03	116.48
4	C	272	BCA	C7B-C2B-C3B	2.84	122.18	118.57
4	B	272	BCA	O5A-P2A-O3A	2.59	114.28	107.27
4	A	272	BCA	CDP-CBP-CAP	-2.46	104.58	108.77
4	B	272	BCA	O5D-C5D-C4D	-2.37	100.92	108.99
4	B	272	BCA	CDP-CBP-CAP	2.36	112.79	108.77
4	B	272	BCA	O4D-C1D-N9A	2.28	112.46	108.09
4	C	272	BCA	P2A-O6A-CCP	-2.23	108.96	121.12
4	B	272	BCA	C2B-C1B-S1P	2.19	119.00	114.97
4	A	272	BCA	CDP-CBP-CCP	2.14	111.75	108.22
4	B	272	BCA	C7B-C2B-C3B	2.13	121.28	118.57
4	C	272	BCA	C2B-C1B-S1P	2.12	118.87	114.97
4	C	272	BCA	OAP-CAP-CBP	-2.03	105.47	110.18

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	272	BCA	CDP-CBP-CCP-O6A
4	B	272	BCA	CEP-CBP-CCP-O6A
4	B	272	BCA	CAP-CBP-CCP-O6A
4	B	272	BCA	O5P-C5P-N4P-C3P
4	C	272	BCA	O1B-C1B-S1P-C2P
4	B	272	BCA	P1A-O3A-P2A-O6A
4	B	272	BCA	S1P-C2P-C3P-N4P
6	B	901	EDO	O1-C1-C2-O2
4	A	272	BCA	O1B-C1B-S1P-C2P

There are no ring outliers.

3 monomers are involved in 4 short contacts:

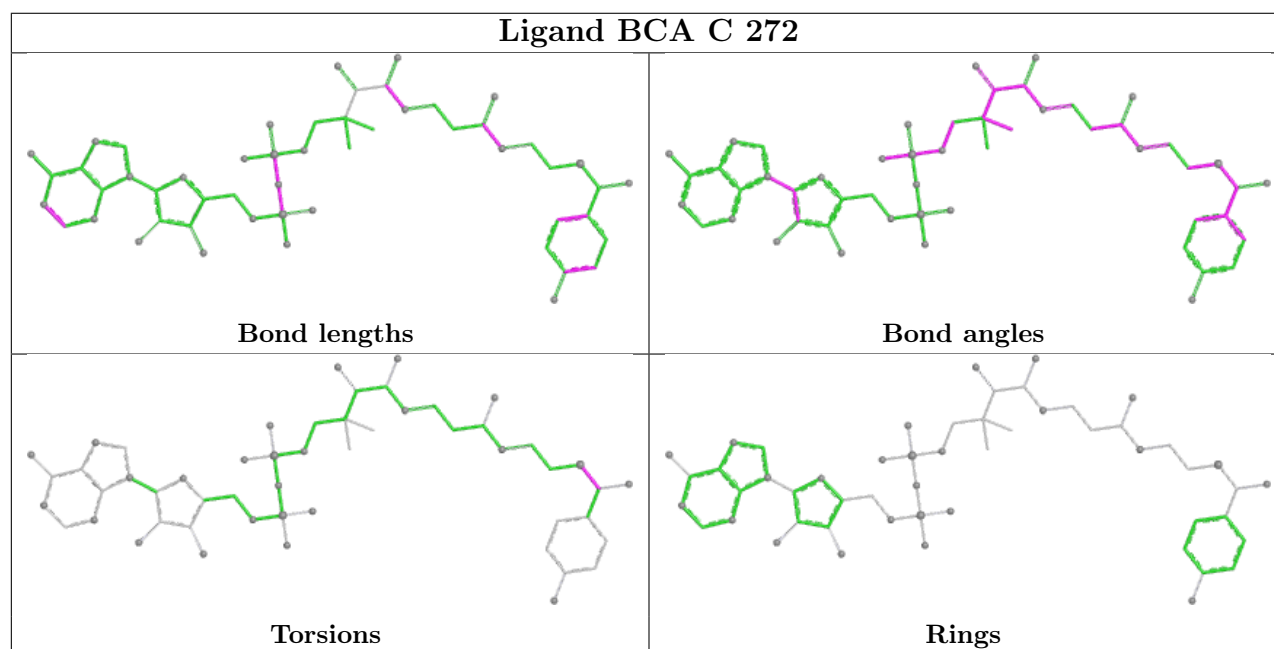
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	272	BCA	2	0

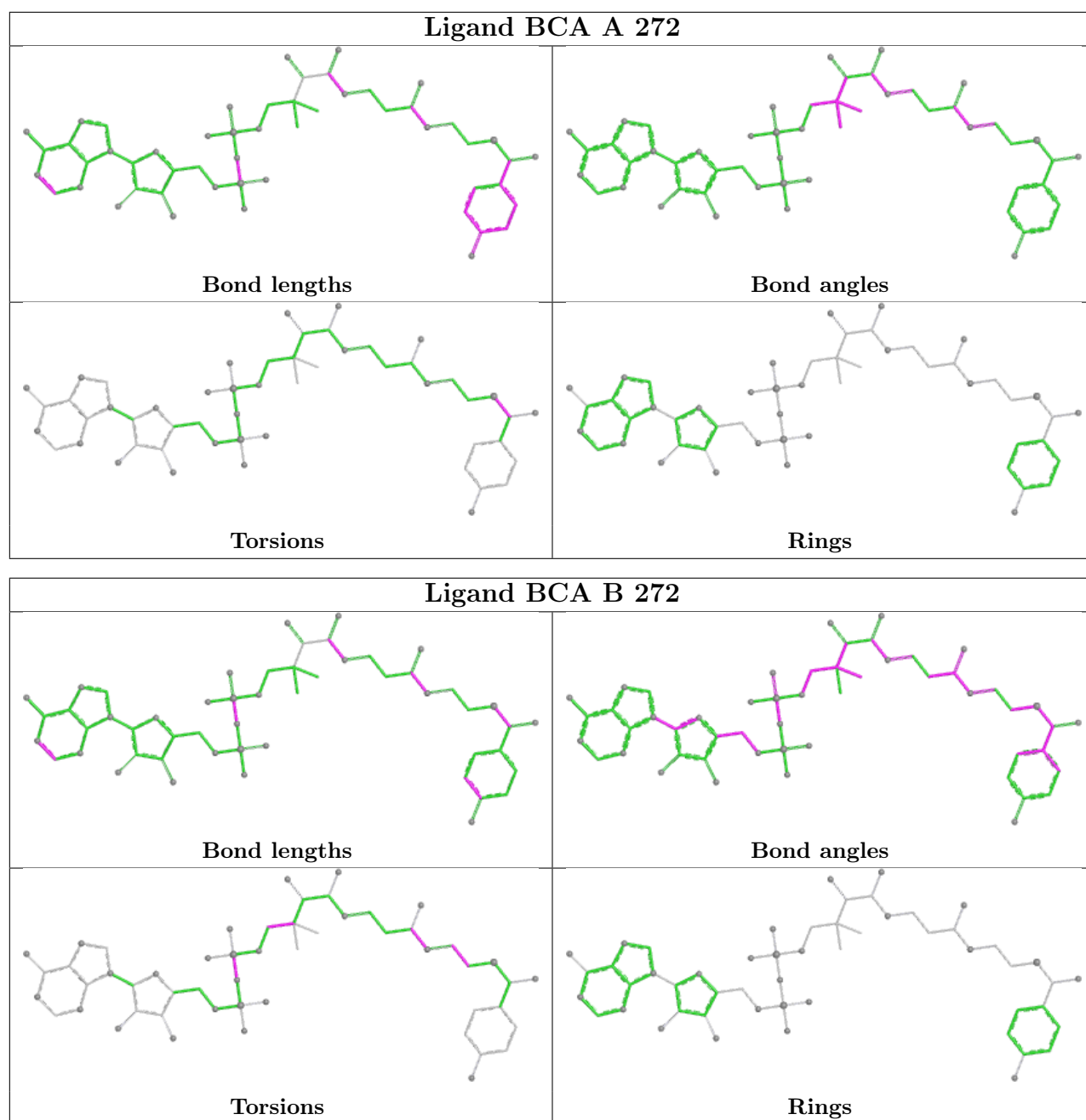
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	272	BCA	1	0
4	B	272	BCA	1	0

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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.