



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:13 PM UTC

PDB ID : 1KRV / pdb_00001krv
Title : Galactoside Acetyltransferase in Complex with CoA and PNP-beta-Gal
Authors : Wang, X.-G.; Olsen, L.R.; Roderick, S.L.
Deposited on : 2002-01-10
Resolution : 2.80 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

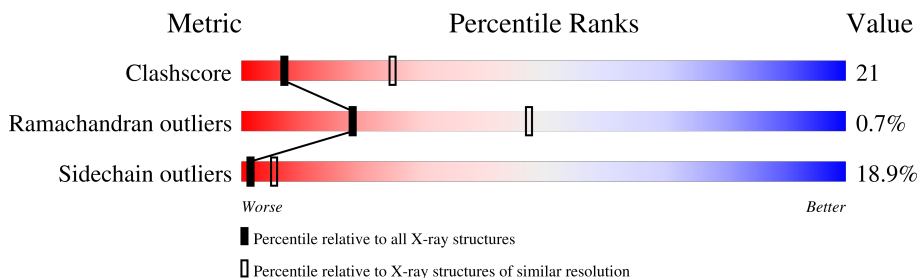
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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	203	
1	B	203	
1	C	203	

2 Entry composition [i](#)

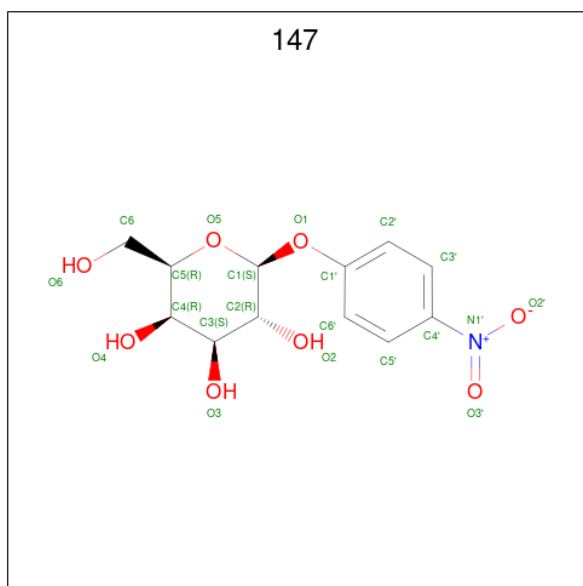
There are 4 unique types of molecules in this entry. The entry contains 5029 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GALACTOSIDE O-ACETYLTRANSFERASE.

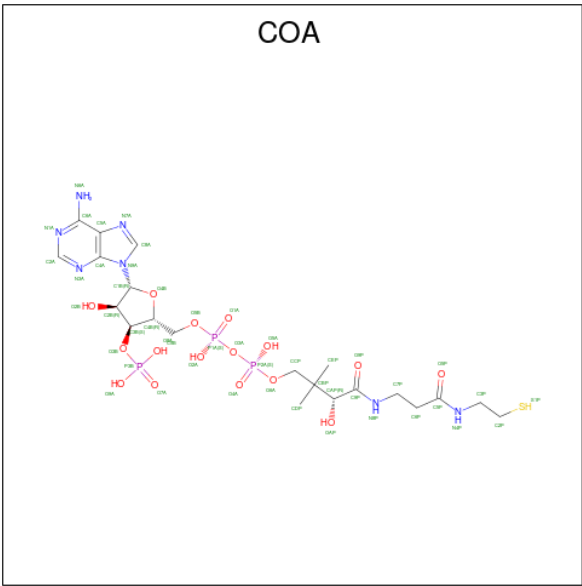
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	201	Total	C	N	O	S	0	0	0
			1584	1007	274	295	8			
1	B	201	Total	C	N	O	S	0	0	0
			1584	1007	274	295	8			
1	C	201	Total	C	N	O	S	0	0	0
			1584	1007	274	295	8			

- Molecule 2 is 4-nitrophenyl beta-D-galactopyranoside (CCD ID: 147) (formula: $C_{12}H_{15}NO_8$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			21	12	1	8		
2	B	1	Total	C	N	O	0	0
			21	12	1	8		
2	C	1	Total	C	N	O	0	0
			21	12	1	8		

- Molecule 3 is COENZYME A (CCD ID: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
3	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
3	C	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 4 is water.

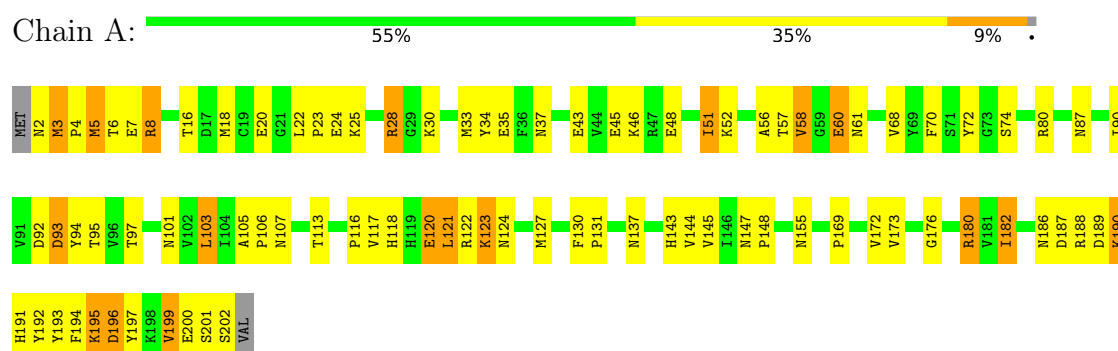
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total	O	0	0
			25	25		
4	B	24	Total	O	0	0
			24	24		
4	C	21	Total	O	0	0
			21	21		

3 Residue-property plots [i](#)

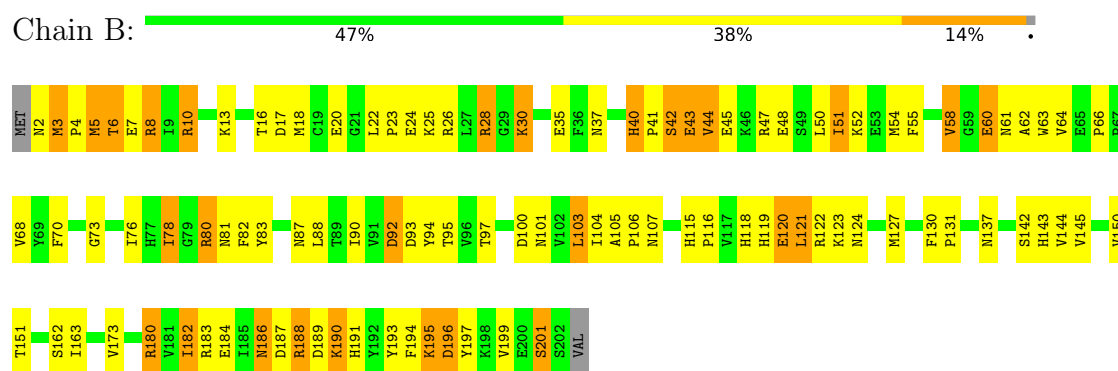
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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.

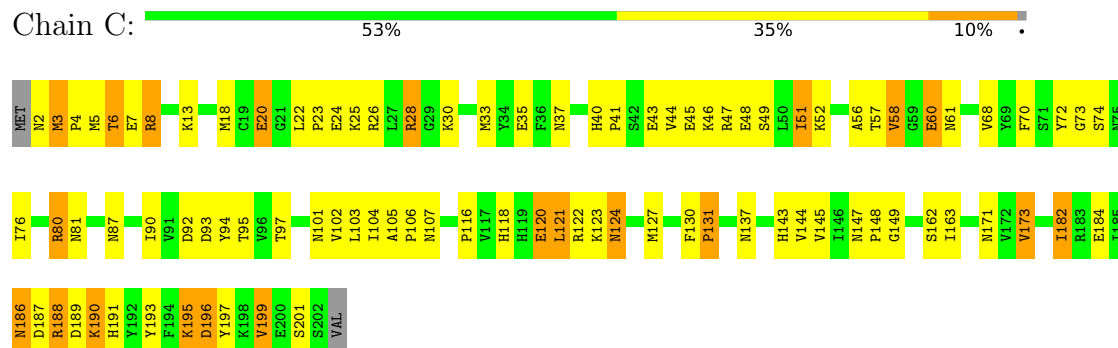
• Molecule 1: GALACTOSIDE O-ACETYLTRANSFERASE



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• Molecule 1: GALACTOSIDE O-ACETYLTRANSFERASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	66.30Å 183.80Å 121.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.00 – 2.80	Depositor
% Data completeness (in resolution range)	94.1 (31.00-2.80)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.174 , 0.253	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5029	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: COA, 147

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	0.58	0/1624	1.04	7/2209 (0.3%)
1	B	0.58	0/1624	1.08	12/2209 (0.5%)
1	C	0.57	0/1624	1.00	7/2209 (0.3%)
All	All	0.58	0/4872	1.04	26/6627 (0.4%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	201	SER	N-CA-C	10.08	132.26	110.80
1	B	95	THR	N-CA-C	7.83	121.75	110.24
1	B	87	ASN	N-CA-C	7.51	121.75	111.39
1	C	95	THR	N-CA-C	7.02	120.42	110.23
1	A	93	ASP	N-CA-C	-6.29	105.57	113.18
1	B	130	PHE	CA-C-N	6.29	126.31	119.90
1	B	130	PHE	C-N-CA	6.29	126.31	119.90
1	B	93	ASP	N-CA-C	-6.23	105.78	113.19
1	B	92	ASP	N-CA-C	6.17	117.87	109.11
1	B	88	LEU	N-CA-C	-5.76	101.75	110.28
1	A	130	PHE	CA-C-N	5.59	126.82	119.84
1	A	130	PHE	C-N-CA	5.59	126.82	119.84
1	B	17	ASP	N-CA-C	5.54	117.97	110.88
1	A	87	ASN	N-CA-C	5.52	118.86	111.24
1	C	87	ASN	N-CA-C	5.44	118.89	111.39
1	C	93	ASP	N-CA-C	-5.41	105.94	112.54
1	B	44	VAL	N-CA-C	5.39	116.16	110.72
1	A	176	GLY	N-CA-C	5.31	117.46	112.04
1	B	40	HIS	CA-C-N	5.15	125.19	119.32
1	B	40	HIS	C-N-CA	5.15	125.19	119.32
1	A	95	THR	N-CA-C	5.13	117.47	110.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	173	VAL	N-CA-C	-5.12	99.79	107.37
1	C	92	ASP	N-CA-C	5.12	117.10	110.65
1	C	44	VAL	N-CA-C	5.07	115.79	110.62
1	A	92	ASP	N-CA-C	5.06	116.29	109.11
1	B	78	ILE	N-CA-C	5.04	115.84	108.53

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	A	1584	0	1548	58	0
1	B	1584	0	1548	80	0
1	C	1584	0	1548	68	0
2	A	21	0	14	1	0
2	B	21	0	14	1	0
2	C	21	0	14	2	0
3	A	48	0	32	1	0
3	B	48	0	32	2	0
3	C	48	0	32	2	0
4	A	25	0	0	2	0
4	B	24	0	0	3	0
4	C	21	0	0	0	0
All	All	5029	0	4782	204	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 21.

All (204) 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:24:GLU:HB3	1:B:28:ARG:HH12	1.16	1.09
1:C:24:GLU:HB3	1:C:28:ARG:HH12	1.25	1.02
1:B:4:PRO:HG2	1:B:7:GLU:HG3	1.49	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:GLU:HB3	1:B:28:ARG:NH1	1.85	0.90
1:A:24:GLU:HB3	1:A:28:ARG:HH12	1.38	0.89
1:C:4:PRO:HG2	1:C:7:GLU:HG3	1.56	0.88
1:C:2:ASN:CG	1:C:3:MET:H	1.83	0.85
1:C:51:ILE:HD11	1:C:58:VAL:HG11	1.60	0.82
1:C:56:ALA:HB2	1:C:74:SER:O	1.79	0.82
1:C:24:GLU:HB3	1:C:28:ARG:NH1	1.95	0.80
1:C:118:HIS:ND1	1:C:120:GLU:HB2	1.98	0.79
1:A:6:THR:HG23	1:A:20:GLU:OE1	1.84	0.78
1:A:51:ILE:HD11	1:A:58:VAL:HG11	1.66	0.78
1:C:191:HIS:O	1:C:199:VAL:HG12	1.84	0.77
1:C:173:VAL:O	1:C:182:ILE:HG12	1.84	0.76
1:B:51:ILE:HD11	1:B:58:VAL:HG11	1.67	0.76
1:A:191:HIS:O	1:A:199:VAL:HG12	1.88	0.74
1:C:60:GLU:CD	1:C:80:ARG:HH21	1.98	0.70
1:A:120:GLU:HA	1:A:123:LYS:HE2	1.73	0.70
1:B:62:ALA:HB2	4:B:2066:HOH:O	1.90	0.70
1:C:22:LEU:HB2	1:C:23:PRO:HD3	1.74	0.70
1:C:37:ASN:HD21	1:C:68:VAL:H	1.38	0.69
1:B:6:THR:HG23	1:B:20:GLU:OE1	1.93	0.69
1:B:60:GLU:CD	1:B:80:ARG:HH21	2.00	0.69
1:C:2:ASN:CG	1:C:3:MET:N	2.50	0.69
1:B:45:GLU:H	1:B:45:GLU:CD	2.01	0.68
1:B:127:MET:HE1	2:B:204:147:O5	1.93	0.68
1:C:173:VAL:HG12	1:C:182:ILE:HD11	1.76	0.68
1:A:180:ARG:HG3	4:A:2041:HOH:O	1.94	0.67
1:A:137:ASN:HB2	1:A:155:ASN:HD22	1.60	0.67
1:B:191:HIS:O	1:B:199:VAL:HG12	1.95	0.66
1:A:4:PRO:HG2	1:A:7:GLU:HG3	1.78	0.65
1:A:116:PRO:HG2	1:A:122:ARG:HB3	1.77	0.65
1:C:45:GLU:H	1:C:45:GLU:CD	2.05	0.64
1:B:2:ASN:CG	1:B:3:MET:H	2.05	0.64
1:C:173:VAL:CG1	1:C:182:ILE:HD11	2.29	0.63
1:A:56:ALA:HB2	1:A:74:SER:O	2.00	0.62
1:B:4:PRO:HG2	1:B:7:GLU:CG	2.26	0.62
1:A:107:ASN:O	1:A:143:HIS:HA	1.98	0.61
1:A:127:MET:HE1	2:A:206:147:O5	2.00	0.60
1:B:42:SER:O	1:B:44:VAL:HG23	2.00	0.60
1:C:127:MET:HE1	2:C:205:147:O5	2.01	0.60
1:C:131:PRO:O	1:C:149:GLY:HA2	2.02	0.60
1:B:116:PRO:HG2	1:B:122:ARG:HB3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:TYR:CE2	1:C:40:HIS:HD2	2.21	0.58
1:A:194:PHE:CZ	1:A:195:LYS:HD3	2.39	0.58
1:C:187:ASP:O	1:C:190:LYS:HB2	2.03	0.58
1:C:73:GLY:HA2	1:C:76:ILE:HD12	1.86	0.58
1:A:22:LEU:HB2	1:A:23:PRO:HD3	1.85	0.57
1:C:4:PRO:HG2	1:C:7:GLU:CG	2.30	0.57
1:C:186:ASN:HD22	1:C:188:ARG:H	1.53	0.56
1:C:105:ALA:HB1	1:C:106:PRO:HD2	1.87	0.56
1:A:24:GLU:HB3	1:A:28:ARG:NH1	2.14	0.56
1:B:24:GLU:CG	1:B:28:ARG:HH22	2.18	0.56
1:B:50:LEU:O	1:B:54:MET:HG2	2.05	0.56
1:C:24:GLU:O	1:C:28:ARG:HG2	2.04	0.56
1:B:186:ASN:C	1:B:186:ASN:HD22	2.13	0.56
1:C:171:ASN:O	1:C:184:GLU:HG3	2.06	0.55
1:A:118:HIS:ND1	1:A:120:GLU:HB2	2.21	0.55
1:A:193:TYR:CZ	1:A:197:TYR:HB3	2.41	0.55
1:A:105:ALA:HB1	1:A:106:PRO:HD2	1.88	0.55
1:A:105:ALA:HB1	1:A:106:PRO:CD	2.37	0.55
1:A:101:ASN:HB2	1:A:137:ASN:HD22	1.72	0.55
1:C:116:PRO:HG2	1:C:122:ARG:HB3	1.87	0.55
1:A:121:LEU:HD11	1:C:189:ASP:O	2.07	0.55
1:B:61:ASN:HB2	1:B:81:ASN:HD22	1.70	0.55
1:C:101:ASN:HB2	1:C:137:ASN:HD22	1.70	0.55
1:B:60:GLU:HG2	1:B:61:ASN:CG	2.32	0.54
1:C:147:ASN:HB3	1:C:148:PRO:HD2	1.89	0.54
1:A:5:MET:HG3	1:A:16:THR:O	2.06	0.54
1:A:45:GLU:H	1:A:45:GLU:CD	2.14	0.54
1:B:2:ASN:CG	1:B:3:MET:N	2.64	0.54
1:C:118:HIS:CE1	1:C:120:GLU:HB2	2.42	0.54
1:C:101:ASN:HB2	1:C:137:ASN:ND2	2.22	0.54
1:A:94:TYR:HD2	1:A:131:PRO:HD3	1.72	0.54
1:A:103:LEU:HD22	1:B:115:HIS:NE2	2.23	0.54
1:B:42:SER:O	1:B:44:VAL:N	2.35	0.53
1:B:106:PRO:HD2	1:B:142:SER:OG	2.09	0.53
1:A:3:MET:HG2	4:A:2034:HOH:O	2.09	0.52
1:A:2:ASN:CG	1:A:3:MET:H	2.18	0.52
1:B:22:LEU:N	1:B:23:PRO:CD	2.73	0.52
1:B:80:ARG:HG3	1:B:100:ASP:OD1	2.09	0.52
1:B:187:ASP:O	1:B:190:LYS:HB2	2.09	0.52
1:C:186:ASN:ND2	1:C:188:ARG:H	2.08	0.51
1:C:124:ASN:N	1:C:124:ASN:ND2	2.57	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ASP:O	1:B:118:HIS:HB2	2.11	0.51
1:A:4:PRO:HG2	1:A:7:GLU:CG	2.41	0.51
1:B:5:MET:HE3	1:B:16:THR:OG1	2.10	0.51
1:C:147:ASN:HB3	1:C:148:PRO:CD	2.40	0.51
1:A:80:ARG:CZ	1:A:80:ARG:HB3	2.38	0.50
1:A:94:TYR:CD2	1:A:131:PRO:HD3	2.46	0.50
1:B:22:LEU:O	1:B:26:ARG:HG3	2.11	0.50
1:B:24:GLU:HG3	1:B:28:ARG:HH22	1.75	0.50
1:B:182:ILE:O	1:B:183:ARG:HB3	2.11	0.50
1:A:189:ASP:O	1:B:121:LEU:HD11	2.11	0.50
1:B:94:TYR:CD2	1:B:131:PRO:HD3	2.46	0.50
1:A:3:MET:O	1:A:8:ARG:HD2	2.12	0.50
1:B:37:ASN:HD21	1:B:68:VAL:H	1.59	0.50
1:B:105:ALA:HB1	1:B:106:PRO:CD	2.42	0.50
1:C:60:GLU:HG2	1:C:61:ASN:CG	2.37	0.50
1:B:40:HIS:ND1	1:B:41:PRO:HD2	2.27	0.50
1:B:118:HIS:ND1	1:B:120:GLU:HB2	2.27	0.49
1:A:144:VAL:CG1	1:A:145:VAL:N	2.76	0.49
1:C:94:TYR:CD2	1:C:131:PRO:HD3	2.47	0.49
1:A:18:MET:HE3	1:A:18:MET:HA	1.94	0.49
1:A:118:HIS:CE1	1:A:120:GLU:HB2	2.48	0.49
1:C:120:GLU:HA	1:C:123:LYS:HE2	1.94	0.49
1:C:186:ASN:HD22	1:C:188:ARG:N	2.10	0.49
1:B:26:ARG:O	1:B:30:LYS:HB2	2.13	0.48
1:B:193:TYR:CZ	1:B:197:TYR:HB3	2.48	0.48
1:B:105:ALA:HB1	1:B:106:PRO:HD2	1.96	0.48
1:A:60:GLU:HG2	1:A:61:ASN:N	2.27	0.48
1:C:61:ASN:HB2	1:C:81:ASN:HD22	1.79	0.48
1:A:2:ASN:CG	1:A:3:MET:N	2.72	0.48
1:C:107:ASN:O	1:C:143:HIS:HA	2.14	0.48
1:B:173:VAL:HG12	1:B:182:ILE:HD11	1.96	0.47
1:C:105:ALA:HB1	1:C:106:PRO:CD	2.44	0.47
1:C:6:THR:HG23	1:C:20:GLU:OE1	2.14	0.47
1:A:37:ASN:HD21	1:A:68:VAL:H	1.63	0.47
1:C:18:MET:HE3	1:C:18:MET:HA	1.97	0.46
1:C:144:VAL:HG12	1:C:145:VAL:N	2.30	0.46
1:A:4:PRO:HG2	1:A:7:GLU:CD	2.40	0.46
1:C:37:ASN:ND2	1:C:68:VAL:H	2.09	0.46
1:C:80:ARG:HB3	1:C:80:ARG:CZ	2.45	0.46
1:C:47:ARG:O	1:C:51:ILE:HG22	2.16	0.46
1:A:144:VAL:HG12	1:A:145:VAL:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASN:HB2	1:A:155:ASN:ND2	2.27	0.46
1:B:64:VAL:O	1:B:66:PRO:HD3	2.16	0.46
1:A:187:ASP:O	1:A:190:LYS:HB2	2.14	0.46
1:B:145:VAL:HB	1:B:163:ILE:HD13	1.98	0.46
1:C:193:TYR:CZ	1:C:197:TYR:HB3	2.51	0.45
1:A:70:PHE:HA	1:A:90:ILE:O	2.17	0.45
1:B:5:MET:HG3	1:B:16:THR:O	2.15	0.45
1:B:10:ARG:HB2	1:B:10:ARG:HH11	1.80	0.45
1:B:107:ASN:O	1:B:143:HIS:HA	2.17	0.45
1:B:144:VAL:HG12	1:B:145:VAL:N	2.32	0.45
1:B:173:VAL:O	1:B:182:ILE:HG12	2.17	0.45
1:C:8:ARG:H	1:C:8:ARG:HG2	1.45	0.45
1:B:61:ASN:O	1:B:81:ASN:HA	2.17	0.45
1:B:101:ASN:HB2	1:B:137:ASN:HD22	1.82	0.45
1:B:119:HIS:HD2	4:B:2054:HOH:O	1.99	0.45
1:B:22:LEU:HB2	1:B:23:PRO:HD3	1.99	0.44
1:B:189:ASP:O	1:C:121:LEU:HD11	2.16	0.44
1:A:33:MET:HB2	1:A:70:PHE:CZ	2.52	0.44
1:B:180:ARG:HG3	4:B:2011:HOH:O	2.17	0.44
1:C:47:ARG:O	1:C:51:ILE:CG2	2.66	0.44
1:B:144:VAL:CG1	1:B:145:VAL:N	2.81	0.44
1:B:145:VAL:HB	1:B:163:ILE:CD1	2.48	0.44
1:A:113:THR:HG21	1:A:127:MET:HE2	2.00	0.44
1:A:192:TYR:HE2	1:A:196:ASP:HA	1.83	0.44
1:B:195:LYS:HB3	1:B:196:ASP:H	1.51	0.44
1:A:169:PRO:O	1:A:172:VAL:HG22	2.16	0.44
1:B:2:ASN:O	1:B:3:MET:SD	2.76	0.44
1:B:63:TRP:HD1	1:B:83:TYR:CE1	2.34	0.44
1:B:61:ASN:HB2	1:B:81:ASN:ND2	2.33	0.44
1:C:33:MET:HB2	1:C:70:PHE:CZ	2.53	0.44
1:A:147:ASN:HB3	1:A:148:PRO:HD2	1.99	0.43
1:C:70:PHE:HA	1:C:90:ILE:O	2.18	0.43
1:A:60:GLU:HG2	1:A:61:ASN:CG	2.43	0.43
1:B:92:ASP:C	1:B:94:TYR:H	2.25	0.43
1:C:26:ARG:HG2	1:C:72:TYR:CE2	2.53	0.43
1:C:40:HIS:CG	1:C:41:PRO:HD2	2.53	0.43
1:C:116:PRO:HB2	1:C:121:LEU:HB2	2.00	0.43
1:B:70:PHE:HA	1:B:90:ILE:O	2.18	0.43
1:A:72:TYR:CE1	1:A:93:ASP:HA	2.54	0.43
1:B:45:GLU:CD	1:B:45:GLU:N	2.74	0.43
1:B:55:PHE:HA	1:B:73:GLY:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:MET:HE3	1:B:18:MET:HA	2.01	0.43
1:B:51:ILE:HA	1:B:54:MET:HG2	2.01	0.43
1:B:78:ILE:HG21	1:B:82:PHE:CG	2.54	0.42
1:B:186:ASN:ND2	1:B:188:ARG:H	2.17	0.42
1:C:61:ASN:OD1	1:C:81:ASN:ND2	2.52	0.42
1:B:92:ASP:C	1:B:94:TYR:N	2.77	0.42
1:B:118:HIS:HB3	1:B:121:LEU:HD12	2.00	0.42
1:C:116:PRO:CB	1:C:121:LEU:HB2	2.49	0.42
1:C:186:ASN:HD22	1:C:186:ASN:C	2.27	0.42
1:A:173:VAL:O	1:A:182:ILE:HG12	2.20	0.42
1:C:81:ASN:HB2	1:C:101:ASN:ND2	2.34	0.42
1:B:73:GLY:HA2	1:B:76:ILE:HD12	2.02	0.42
1:C:144:VAL:CG1	1:C:145:VAL:N	2.83	0.41
1:B:47:ARG:O	1:B:51:ILE:CG2	2.68	0.41
3:B:208:COA:OAP	3:B:208:COA:N7A	2.53	0.41
1:B:47:ARG:O	1:B:51:ILE:HG23	2.19	0.41
1:B:150:VAL:HG12	1:B:151:THR:N	2.35	0.41
1:A:60:GLU:CD	1:A:80:ARG:HH21	2.28	0.41
3:A:207:COA:OAP	3:A:207:COA:N7A	2.53	0.41
1:C:145:VAL:HB	1:C:163:ILE:HD12	2.01	0.41
1:B:81:ASN:HB2	1:B:101:ASN:ND2	2.36	0.41
1:C:4:PRO:HG2	1:C:7:GLU:CD	2.46	0.41
1:A:195:LYS:HB3	1:A:196:ASP:H	1.64	0.41
3:C:209:COA:N7A	3:C:209:COA:OAP	2.53	0.41
1:A:118:HIS:HB3	1:A:121:LEU:HD12	2.02	0.41
1:B:8:ARG:H	1:B:8:ARG:HG2	1.38	0.41
1:B:50:LEU:HA	1:B:50:LEU:HD23	1.90	0.41
3:B:208:COA:HN8	3:B:208:COA:H141	1.85	0.41
3:C:209:COA:H141	3:C:209:COA:HN8	1.86	0.41
1:A:201:SER:O	1:A:202:SER:C	2.63	0.41
1:B:194:PHE:CD2	1:C:130:PHE:CE2	3.09	0.41
1:A:116:PRO:HB2	1:A:121:LEU:HB2	2.03	0.40
1:B:40:HIS:CG	1:B:41:PRO:HD2	2.57	0.40
1:C:173:VAL:HG12	1:C:182:ILE:CD1	2.49	0.40
1:B:103:LEU:HD21	2:C:205:147:H3'	2.04	0.40
1:C:80:ARG:CZ	1:C:80:ARG:CB	2.99	0.40
1:C:195:LYS:HB3	1:C:196:ASP:H	1.50	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	199/203 (98%)	188 (94%)	10 (5%)	1 (0%)	24	55
1	B	199/203 (98%)	186 (94%)	11 (6%)	2 (1%)	12	38
1	C	199/203 (98%)	188 (94%)	10 (5%)	1 (0%)	24	55
All	All	597/609 (98%)	562 (94%)	31 (5%)	4 (1%)	18	47

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	201	SER
1	B	43	GLU
1	A	200	GLU
1	C	131	PRO

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	176/179 (98%)	146 (83%)	30 (17%)	2	7
1	B	176/179 (98%)	141 (80%)	35 (20%)	1	5
1	C	176/179 (98%)	141 (80%)	35 (20%)	1	5
All	All	528/537 (98%)	428 (81%)	100 (19%)	1	5

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	MET
1	A	5	MET
1	A	8	ARG
1	A	25	LYS
1	A	28	ARG
1	A	30	LYS
1	A	35	GLU
1	A	43	GLU
1	A	46	LYS
1	A	48	GLU
1	A	51	ILE
1	A	52	LYS
1	A	57	THR
1	A	58	VAL
1	A	60	GLU
1	A	97	THR
1	A	103	LEU
1	A	117	VAL
1	A	120	GLU
1	A	121	LEU
1	A	123	LYS
1	A	124	ASN
1	A	180	ARG
1	A	182	ILE
1	A	186	ASN
1	A	188	ARG
1	A	190	LYS
1	A	195	LYS
1	A	196	ASP
1	A	199	VAL
1	B	3	MET
1	B	5	MET
1	B	6	THR
1	B	8	ARG
1	B	10	ARG
1	B	13	LYS
1	B	25	LYS
1	B	28	ARG
1	B	30	LYS
1	B	35	GLU
1	B	42	SER
1	B	43	GLU
1	B	48	GLU

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Mol	Chain	Res	Type
1	B	51	ILE
1	B	52	LYS
1	B	58	VAL
1	B	60	GLU
1	B	80	ARG
1	B	97	THR
1	B	103	LEU
1	B	104	ILE
1	B	120	GLU
1	B	121	LEU
1	B	123	LYS
1	B	124	ASN
1	B	162	SER
1	B	180	ARG
1	B	182	ILE
1	B	184	GLU
1	B	186	ASN
1	B	188	ARG
1	B	190	LYS
1	B	195	LYS
1	B	196	ASP
1	B	201	SER
1	C	3	MET
1	C	5	MET
1	C	6	THR
1	C	8	ARG
1	C	13	LYS
1	C	20	GLU
1	C	25	LYS
1	C	28	ARG
1	C	30	LYS
1	C	35	GLU
1	C	43	GLU
1	C	46	LYS
1	C	48	GLU
1	C	49	SER
1	C	51	ILE
1	C	52	LYS
1	C	57	THR
1	C	58	VAL
1	C	60	GLU
1	C	80	ARG

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Mol	Chain	Res	Type
1	C	97	THR
1	C	102	VAL
1	C	103	LEU
1	C	104	ILE
1	C	120	GLU
1	C	121	LEU
1	C	124	ASN
1	C	162	SER
1	C	182	ILE
1	C	186	ASN
1	C	188	ARG
1	C	190	LYS
1	C	195	LYS
1	C	196	ASP
1	C	199	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	38	HIS
1	A	77	HIS
1	A	81	ASN
1	A	85	ASN
1	A	137	ASN
1	A	147	ASN
1	A	155	ASN
1	A	171	ASN
1	A	186	ASN
1	B	37	ASN
1	B	77	HIS
1	B	81	ASN
1	B	101	ASN
1	B	119	HIS
1	B	137	ASN
1	B	147	ASN
1	B	155	ASN
1	B	171	ASN
1	B	186	ASN
1	C	37	ASN
1	C	61	ASN
1	C	77	HIS

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Mol	Chain	Res	Type
1	C	81	ASN
1	C	101	ASN
1	C	137	ASN
1	C	147	ASN
1	C	155	ASN
1	C	186	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 ⓘ

6 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	COA	A	207	-	47,50,50	1.64	4 (8%)	69,75,75	1.08	4 (5%)
2	147	A	206	-	22,22,22	3.21	13 (59%)	29,31,31	2.06	5 (17%)
2	147	C	205	-	22,22,22	3.21	12 (54%)	29,31,31	2.01	4 (13%)
2	147	B	204	-	22,22,22	3.25	12 (54%)	29,31,31	2.00	4 (13%)
3	COA	C	209	-	47,50,50	1.58	3 (6%)	69,75,75	1.09	3 (4%)
3	COA	B	208	-	47,50,50	1.67	3 (6%)	69,75,75	1.07	3 (4%)

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	COA	A	207	-	-	11/48/64/64	0/3/3/3
2	147	A	206	-	-	4/8/30/30	0/2/2/2
2	147	C	205	-	-	4/8/30/30	0/2/2/2
2	147	B	204	-	-	4/8/30/30	0/2/2/2
3	COA	C	209	-	-	12/48/64/64	0/3/3/3
3	COA	B	208	-	-	12/48/64/64	0/3/3/3

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	208	COA	P2A-O3A	8.23	1.68	1.59
3	A	207	COA	P2A-O3A	7.96	1.68	1.59
3	C	209	COA	P2A-O3A	7.69	1.67	1.59
2	C	205	147	C6'-C1'	7.17	1.52	1.38
2	B	204	147	C6'-C1'	7.09	1.52	1.38
2	A	206	147	C6'-C1'	7.00	1.51	1.38
2	A	206	147	O1-C1'	6.97	1.51	1.38
2	B	204	147	O1-C1'	6.84	1.51	1.38
2	C	205	147	O1-C1'	6.70	1.51	1.38
2	B	204	147	C3'-C4'	6.01	1.49	1.38
2	C	205	147	C3'-C4'	5.96	1.49	1.38
2	A	206	147	C3'-C4'	5.82	1.49	1.38
3	B	208	COA	CCP-CBP	4.18	1.59	1.52
2	C	205	147	O5-C5	3.76	1.53	1.44
2	B	204	147	O5-C5	3.62	1.53	1.44
2	A	206	147	O5-C5	3.58	1.53	1.44
3	A	207	COA	CCP-CBP	3.42	1.58	1.52
2	C	205	147	C4'-N1'	3.33	1.53	1.45
3	C	209	COA	CCP-CBP	3.32	1.58	1.52
2	B	204	147	C4-C3	3.27	1.60	1.52
2	B	204	147	C4'-N1'	3.25	1.52	1.45
2	B	204	147	O5-C1	3.23	1.50	1.41
2	A	206	147	C4-C3	3.20	1.60	1.52
2	A	206	147	O5-C1	3.19	1.50	1.41
2	A	206	147	C2'-C1'	3.13	1.44	1.38
2	C	205	147	C4-C5	3.09	1.59	1.53
2	B	204	147	C4-C5	3.07	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	206	147	C4'-N1'	2.97	1.52	1.45
2	B	204	147	C2'-C1'	2.92	1.44	1.38
2	C	205	147	C2'-C1'	2.89	1.44	1.38
2	C	205	147	C4-C3	2.89	1.59	1.52
2	A	206	147	C4-C5	2.78	1.59	1.53
2	C	205	147	O5-C1	2.73	1.48	1.41
2	B	204	147	C3-C2	2.67	1.59	1.52
3	A	207	COA	O6A-CCP	2.57	1.52	1.43
2	A	206	147	C3-C2	2.51	1.58	1.52
3	C	209	COA	O6A-CCP	2.49	1.51	1.43
2	C	205	147	C3-C2	2.44	1.58	1.52
2	A	206	147	C6'-C5'	2.39	1.42	1.38
3	B	208	COA	O6A-CCP	2.37	1.51	1.43
2	B	204	147	C6'-C5'	2.21	1.42	1.38
2	C	205	147	C6'-C5'	2.17	1.42	1.38
2	B	204	147	O2'-N1'	-2.14	1.21	1.35
2	C	205	147	O2'-N1'	-2.10	1.21	1.35
2	A	206	147	O2'-N1'	-2.07	1.21	1.35
2	A	206	147	C3'-C2'	2.04	1.42	1.38
3	A	207	COA	C4A-N3A	2.02	1.38	1.34

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	206	147	C3'-C4'-N1'	-6.72	113.47	119.34
2	B	204	147	C3'-C4'-N1'	-6.42	113.73	119.34
2	C	205	147	C3'-C4'-N1'	-6.41	113.74	119.34
2	C	205	147	C1'-O1-C1	5.19	125.13	117.82
2	A	206	147	C1'-O1-C1	4.93	124.77	117.82
2	B	204	147	C1'-O1-C1	4.84	124.64	117.82
2	A	206	147	C5'-C4'-N1'	4.67	123.41	119.34
2	B	204	147	C5'-C4'-N1'	4.59	123.34	119.34
2	C	205	147	C5'-C4'-N1'	4.57	123.33	119.34
3	A	207	COA	O6A-CCP-CBP	4.28	117.44	110.55
3	C	209	COA	O6A-CCP-CBP	4.25	117.38	110.55
3	B	208	COA	O6A-CCP-CBP	4.15	117.22	110.55
2	B	204	147	O5-C5-C6	2.58	112.83	106.44
2	A	206	147	O5-C5-C6	2.55	112.77	106.44
3	C	209	COA	O4B-C1B-N9A	2.49	112.87	108.09
2	C	205	147	O5-C5-C6	2.49	112.60	106.44
3	B	208	COA	O4B-C1B-N9A	2.47	112.83	108.09
3	C	209	COA	C3B-C2B-C1B	2.33	105.02	99.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	207	COA	C3B-C2B-C1B	2.30	104.94	99.89
3	A	207	COA	O4B-C1B-N9A	2.26	112.43	108.09
3	B	208	COA	C3B-C2B-C1B	2.25	104.85	99.89
3	A	207	COA	C6P-C7P-N8P	-2.14	107.45	112.00
2	A	206	147	C5'-C4'-C3'	2.01	123.11	119.88

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	206	147	C3'-C4'-N1'-O3'
2	A	206	147	C5'-C4'-N1'-O3'
2	B	204	147	C3'-C4'-N1'-O3'
2	B	204	147	C5'-C4'-N1'-O3'
2	C	205	147	C3'-C4'-N1'-O3'
2	C	205	147	C5'-C4'-N1'-O3'
3	A	207	COA	C5B-O5B-P1A-O1A
3	A	207	COA	C5B-O5B-P1A-O3A
3	A	207	COA	CDP-CBP-CCP-O6A
3	A	207	COA	CEP-CBP-CCP-O6A
3	A	207	COA	CAP-CBP-CCP-O6A
3	A	207	COA	O9P-C9P-CAP-CBP
3	A	207	COA	N8P-C9P-CAP-CBP
3	A	207	COA	N8P-C9P-CAP-OAP
3	B	208	COA	C5B-O5B-P1A-O1A
3	B	208	COA	C5B-O5B-P1A-O3A
3	B	208	COA	CDP-CBP-CCP-O6A
3	B	208	COA	CEP-CBP-CCP-O6A
3	B	208	COA	CAP-CBP-CCP-O6A
3	B	208	COA	O9P-C9P-CAP-CBP
3	B	208	COA	N8P-C9P-CAP-CBP
3	B	208	COA	N8P-C9P-CAP-OAP
3	C	209	COA	C5B-O5B-P1A-O1A
3	C	209	COA	C5B-O5B-P1A-O3A
3	C	209	COA	CDP-CBP-CCP-O6A
3	C	209	COA	CEP-CBP-CCP-O6A
3	C	209	COA	CAP-CBP-CCP-O6A
3	C	209	COA	O9P-C9P-CAP-CBP
3	C	209	COA	N8P-C9P-CAP-CBP
3	C	209	COA	N8P-C9P-CAP-OAP
3	A	207	COA	C2B-C3B-O3B-P3B
3	A	207	COA	C4B-C3B-O3B-P3B

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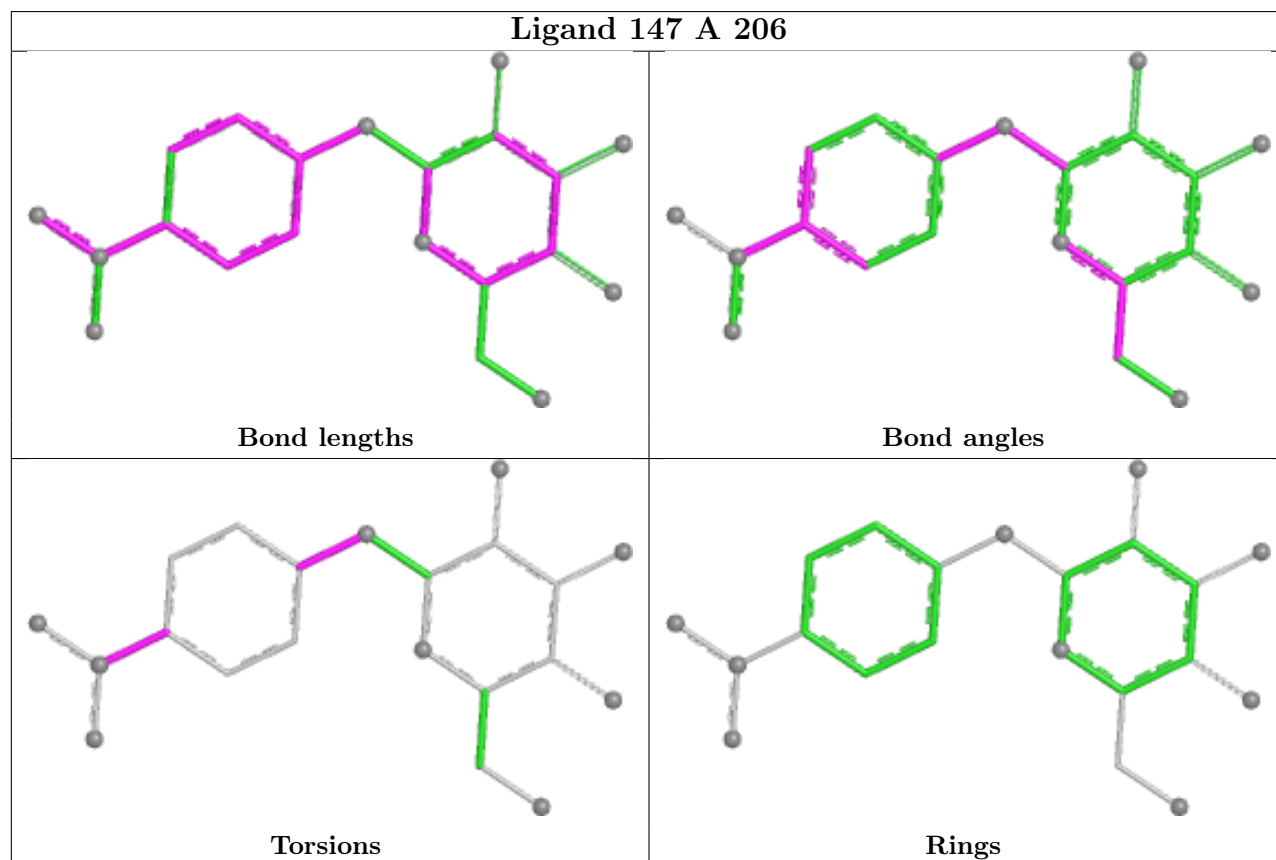
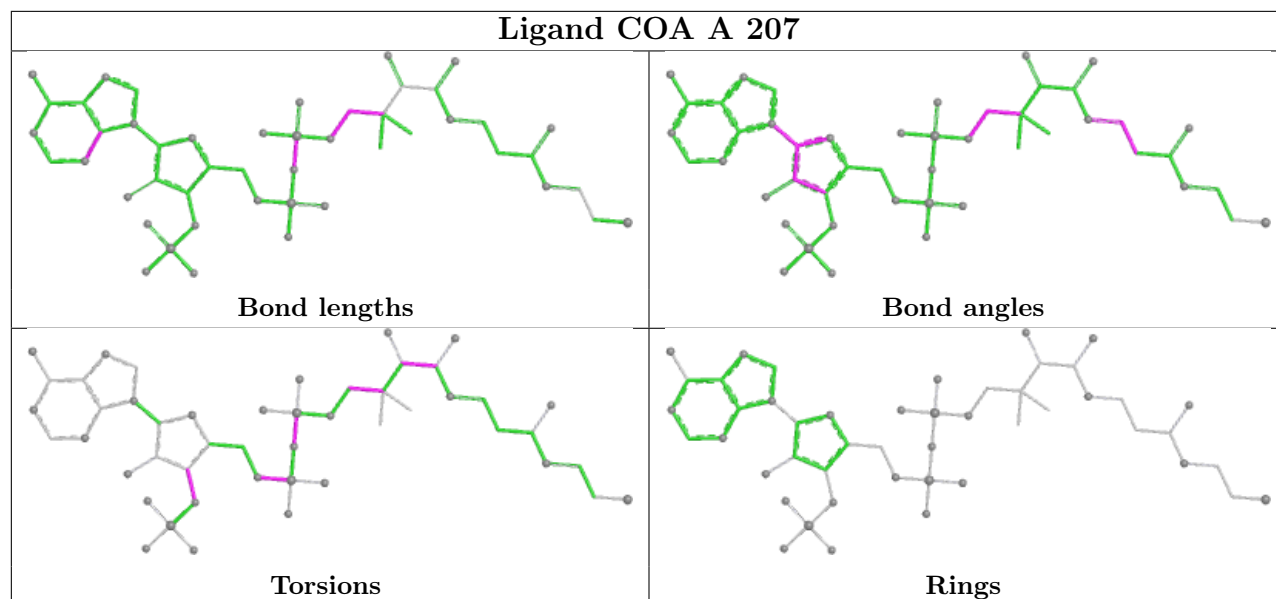
Mol	Chain	Res	Type	Atoms
3	B	208	COA	C2B-C3B-O3B-P3B
3	B	208	COA	C4B-C3B-O3B-P3B
3	C	209	COA	C2B-C3B-O3B-P3B
3	C	209	COA	C4B-C3B-O3B-P3B
2	C	205	147	C6'-C1'-O1-C1
3	A	207	COA	P1A-O3A-P2A-O5A
3	B	208	COA	P1A-O3A-P2A-O5A
3	C	209	COA	P1A-O3A-P2A-O5A
2	A	206	147	C6'-C1'-O1-C1
2	B	204	147	C6'-C1'-O1-C1
2	A	206	147	C2'-C1'-O1-C1
2	B	204	147	C2'-C1'-O1-C1
2	C	205	147	C2'-C1'-O1-C1
3	B	208	COA	C2B-C1B-N9A-C8A
3	C	209	COA	C2B-C1B-N9A-C8A

There are no ring outliers.

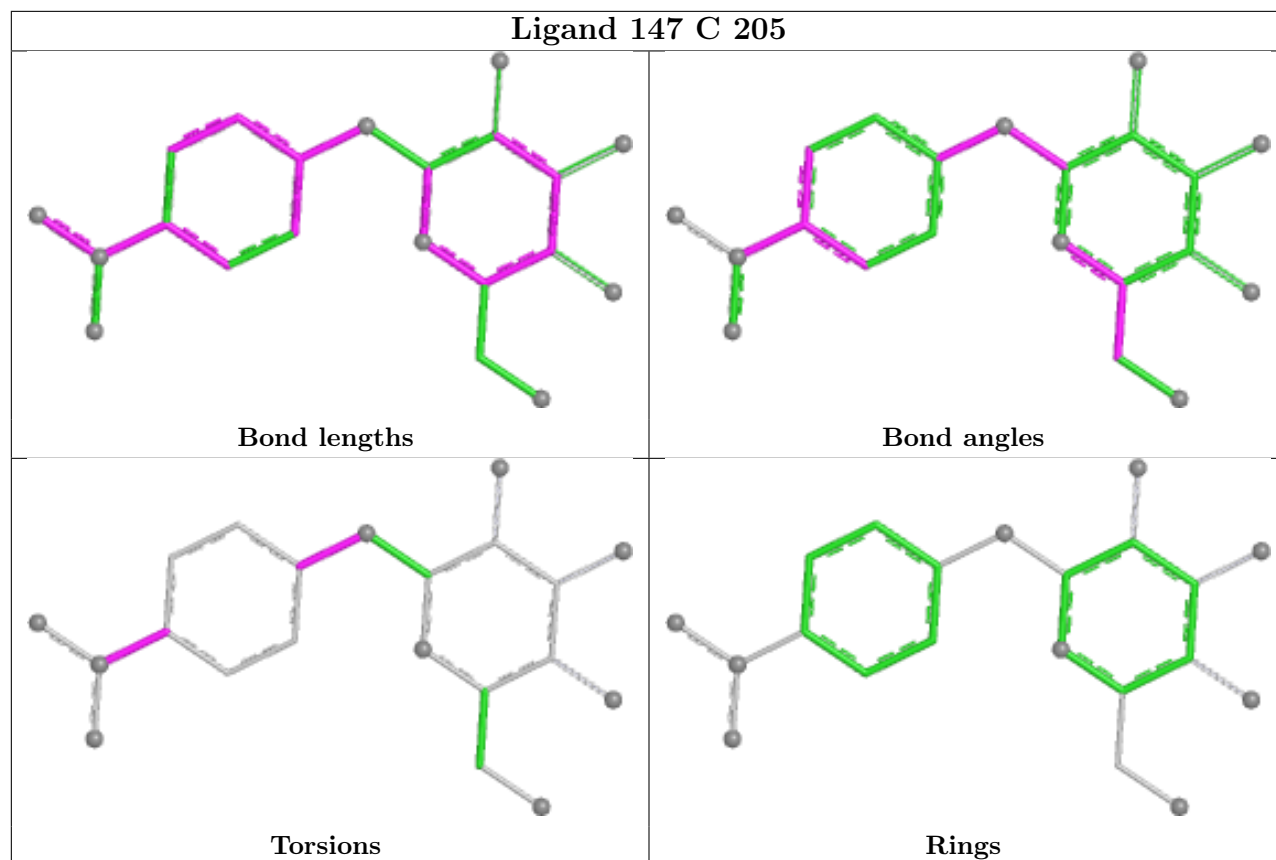
6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	207	COA	1	0
2	A	206	147	1	0
2	C	205	147	2	0
2	B	204	147	1	0
3	C	209	COA	2	0
3	B	208	COA	2	0

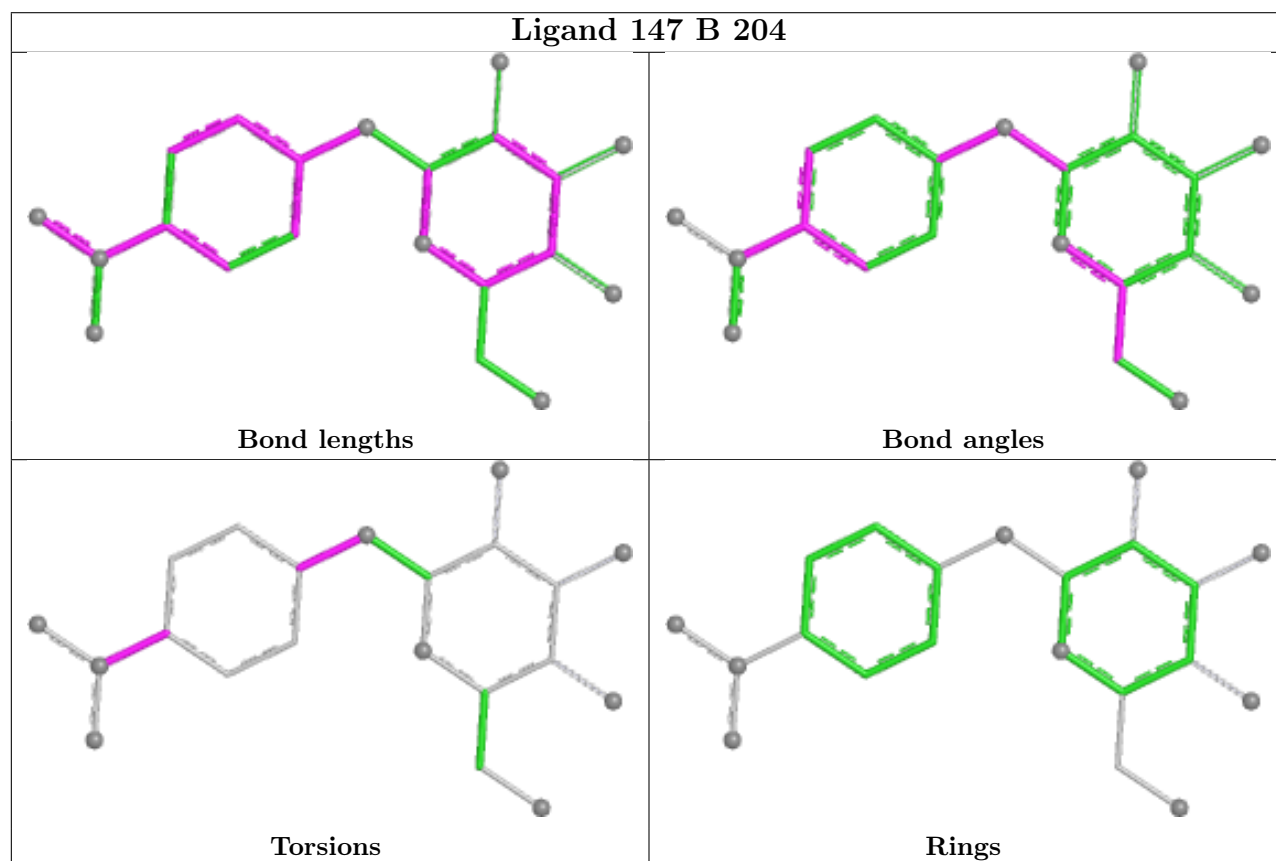
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

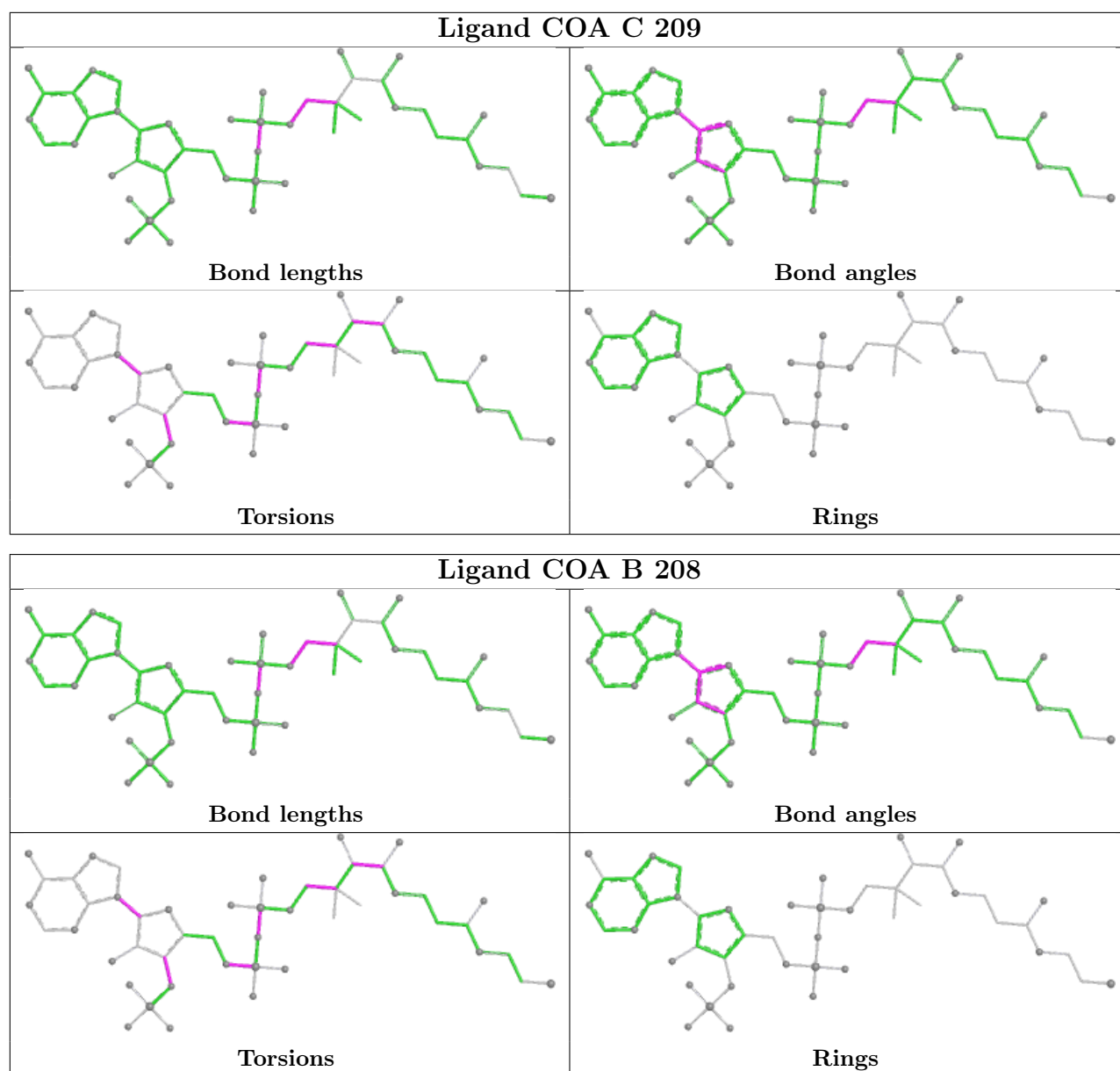


Ligand 147 C 205



Ligand 147 B 204





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.