



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 12:51 AM GMT

PDB ID : 2BT4
Title : TYPE II DEHYDROQUINASE INHIBITOR COMPLEX
Authors : Toscano, M.D.; Stewart, K.A.; Coggins, J.R.; Lapthorn, A.J.; Abell, C.
Deposited on : 2005-05-26
Resolution : 1.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

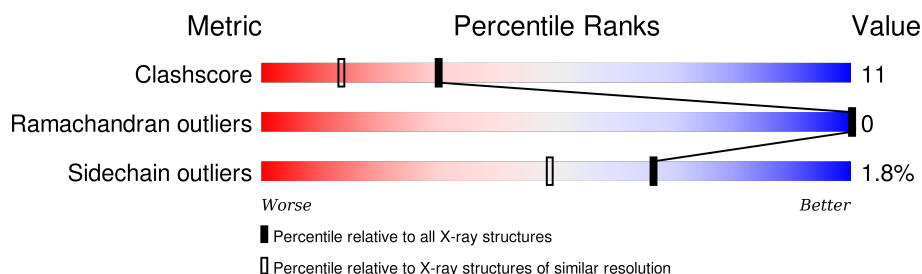
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)



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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	157	
1	B	157	
1	C	157	
1	D	157	
1	E	157	
1	F	157	
1	G	157	

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Mol	Chain	Length	Quality of chain
1	H	157	 84% 10% • 5%
1	I	157	 80% 15% 5%
1	J	157	 82% 11% • 5%
1	K	157	 83% 12% 5%
1	L	157	 82% 12% • 5%

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	PO4	A	2401	-	-	X	-
5	GOL	B	2414	-	-	X	-
5	GOL	C	2415	-	-	X	-
5	GOL	D	2416	-	-	X	-
5	GOL	E	2419	-	-	X	-
5	GOL	E	2422	-	-	X	-
5	GOL	F	2418	-	-	X	-
5	GOL	G	2417	-	-	X	-
5	GOL	H	2411	-	-	X	-
5	GOL	I	2420	-	-	X	-
5	GOL	J	2412	-	-	X	-
5	GOL	K	2421	-	-	X	-

2 Entry composition

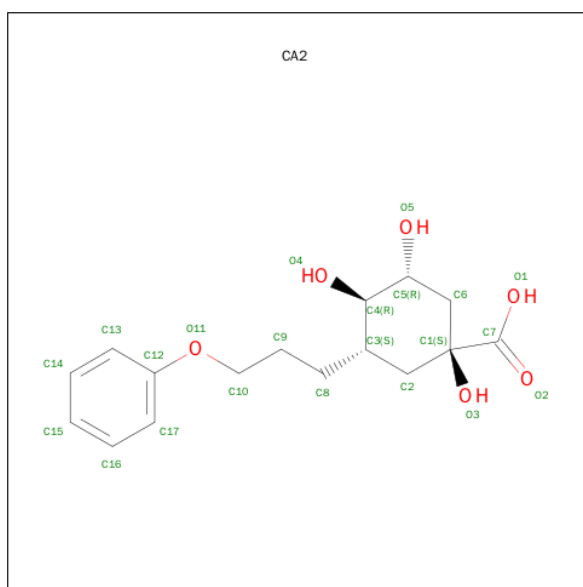
There are 6 unique types of molecules in this entry. The entry contains 15641 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 3-DEHYDROQUINATE DEHYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	2	0
			1127	701	210	211	5			
1	B	149	Total	C	N	O	S	0	1	0
			1126	701	210	210	5			
1	C	149	Total	C	N	O	S	0	1	0
			1126	701	210	210	5			
1	D	149	Total	C	N	O	S	0	2	0
			1127	701	210	211	5			
1	E	149	Total	C	N	O	S	0	1	0
			1126	701	210	210	5			
1	F	149	Total	C	N	O	S	0	1	0
			1126	701	210	210	5			
1	G	149	Total	C	N	O	S	0	2	0
			1127	701	210	211	5			
1	H	149	Total	C	N	O	S	0	1	0
			1126	701	210	210	5			
1	I	149	Total	C	N	O	S	0	1	0
			1126	701	210	210	5			
1	J	149	Total	C	N	O	S	0	2	0
			1132	704	213	210	5			
1	K	149	Total	C	N	O	S	0	1	0
			1126	701	210	210	5			
1	L	149	Total	C	N	O	S	0	2	0
			1127	701	210	211	5			

- Molecule 2 is (1S,3R,4R,5S)-1,3,4-TRIHIDROXY-5-(3-PHENOXYPROPYL)CYCLOHEXANECARBOXYLIC ACID (three-letter code: CA2) (formula: C₁₆H₂₂O₆).



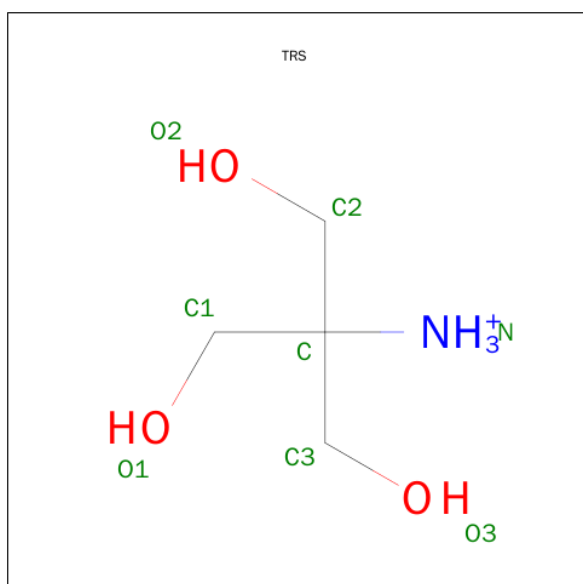
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			22	16	6		
2	B	1	Total	C	O	0	0
			22	16	6		
2	C	1	Total	C	O	0	0
			22	16	6		
2	D	1	Total	C	O	0	0
			22	16	6		
2	E	1	Total	C	O	0	0
			22	16	6		
2	F	1	Total	C	O	0	0
			22	16	6		
2	G	1	Total	C	O	0	0
			22	16	6		
2	H	1	Total	C	O	0	0
			22	16	6		
2	I	1	Total	C	O	0	0
			22	16	6		
2	J	1	Total	C	O	0	0
			22	16	6		
2	K	1	Total	C	O	0	0
			22	16	6		
2	L	1	Total	C	O	0	0
			22	16	6		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



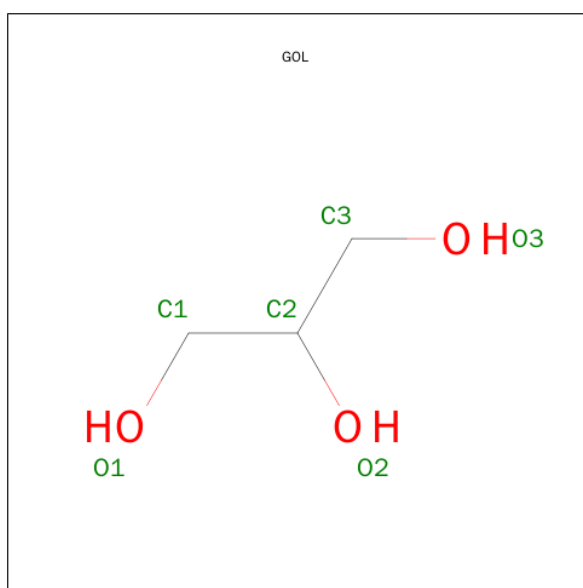
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		
3	I	1	Total	O	P	0	0
			5	4	1		
3	L	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			8	4	1	3		
4	F	1	Total	C	N	O	0	0
			8	4	1	3		
4	H	1	Total	C	N	O	0	0
			8	4	1	3		
4	L	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		
5	F	1	Total	C	O	0	0
			6	3	3		
5	G	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	H	1	Total	C	O	0	0
			6	3	3		
5	I	1	Total	C	O	0	0
			6	3	3		
5	J	1	Total	C	O	0	0
			6	3	3		
5	K	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	135	Total	O	0	0
			135	135		
6	B	152	Total	O	0	0
			152	152		
6	C	146	Total	O	0	0
			146	146		
6	D	146	Total	O	0	0
			146	146		
6	E	186	Total	O	0	0
			186	186		
6	F	146	Total	O	0	0
			146	146		
6	G	148	Total	O	0	0
			148	148		
6	H	124	Total	O	0	0
			124	124		
6	I	156	Total	O	0	0
			156	156		
6	J	120	Total	O	0	0
			120	120		
6	K	123	Total	O	0	0
			123	123		
6	L	149	Total	O	0	0
			149	149		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.


Note EDS was not executed.

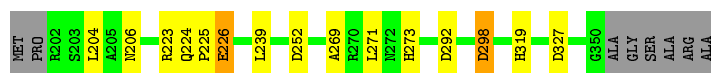
• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain A: 




• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain B: 




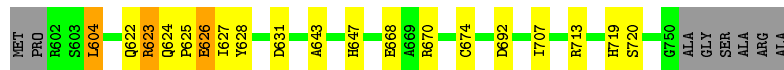
• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain C: 




• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain D: 




• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain E: 



• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain F:  81% 13% 5%




• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain G:  86% 8% 5%



• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain H:  84% 10% 5%




• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain I:  80% 15% 5%




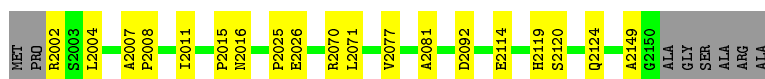
• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain J:  82% 11% 5%




• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain K:  83% 12% 5%



• Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain L:  82% 12% 5%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	196.62Å 196.49Å 240.63Å 65.91° 65.91° 90.01°	Depositor
Resolution (Å)	27.00 – 1.70	Depositor
% Data completeness (in resolution range)	75.8 (27.00-1.70)	Depositor
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.197 , 0.248	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	15641	wwPDB-VP
Average B, all atoms (Å ²)	15.0	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: GOL, PO4, TRS, CA2

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	1.02	0/1161	1.04	4/1582 (0.3%)
1	B	1.05	0/1155	0.99	4/1574 (0.3%)
1	C	1.01	0/1155	0.95	3/1574 (0.2%)
1	D	1.05	0/1161	0.95	2/1582 (0.1%)
1	E	1.01	2/1155 (0.2%)	0.94	3/1574 (0.2%)
1	F	1.06	1/1155 (0.1%)	1.02	6/1574 (0.4%)
1	G	1.03	1/1161 (0.1%)	1.01	3/1582 (0.2%)
1	H	1.04	0/1155	1.02	2/1574 (0.1%)
1	I	1.00	0/1155	0.96	2/1574 (0.1%)
1	J	1.02	1/1166 (0.1%)	0.98	3/1588 (0.2%)
1	K	1.00	0/1155	0.93	0/1574
1	L	1.02	0/1161	0.98	3/1582 (0.2%)
All	All	1.03	5/13895 (0.0%)	0.98	35/18934 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	1083	TYR	CD1-CE1	6.32	1.48	1.39
1	G	1305	VAL	CB-CG2	-5.66	1.41	1.52
1	E	828	TYR	CD2-CE2	-5.37	1.31	1.39
1	E	939	VAL	CB-CG1	-5.24	1.41	1.52
1	J	1940	PHE	CD2-CE2	5.09	1.49	1.39

The worst 5 of 35 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	1527	ASP	CB-CG-OD2	13.19	130.17	118.30
1	G	1327	ASP	CB-CG-OD2	8.20	125.68	118.30
1	C	527	ASP	CB-CG-OD2	7.63	125.17	118.30
1	B	252	ASP	CB-CG-OD1	7.24	124.82	118.30
1	D	631	ASP	CB-CG-OD2	7.06	124.65	118.30

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	1127	0	1093	20	0
1	B	1126	0	1092	14	0
1	C	1126	0	1092	17	0
1	D	1127	0	1093	18	0
1	E	1126	0	1092	38	0
1	F	1126	0	1092	36	0
1	G	1127	0	1093	16	0
1	H	1126	0	1092	27	0
1	I	1126	0	1092	41	0
1	J	1132	0	1102	19	0
1	K	1126	0	1092	20	0
1	L	1127	0	1093	22	0
2	A	22	0	21	1	0
2	B	22	0	21	2	0
2	C	22	0	21	0	0
2	D	22	0	21	0	0
2	E	22	0	21	2	0
2	F	22	0	21	2	0
2	G	22	0	21	1	0
2	H	22	0	21	1	0
2	I	22	0	21	0	0
2	J	22	0	21	1	0
2	K	22	0	21	0	0
2	L	22	0	21	0	0
3	A	5	0	0	2	0
3	D	5	0	0	1	0
3	I	5	0	0	1	0
3	L	5	0	0	0	0
4	A	8	0	12	0	0
4	F	8	0	12	0	0
4	H	8	0	12	0	0
4	L	8	0	12	0	0
5	A	6	0	8	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	6	0	8	6	0
5	C	6	0	8	9	0
5	D	6	0	8	9	0
5	E	12	0	16	15	0
5	F	6	0	8	13	0
5	G	6	0	8	7	0
5	H	6	0	8	7	0
5	I	6	0	8	21	0
5	J	6	0	8	4	0
5	K	6	0	8	12	0
6	A	135	0	0	10	0
6	B	152	0	0	8	0
6	C	146	0	0	10	0
6	D	146	0	0	5	0
6	E	186	0	0	17	0
6	F	146	0	0	10	0
6	G	148	0	0	0	0
6	H	124	0	0	18	0
6	I	156	0	0	22	0
6	J	120	0	0	13	0
6	K	123	0	0	2	0
6	L	149	0	0	13	0
All	All	15641	0	13514	305	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 11.

The worst 5 of 305 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:I:1719[A]:HIS:CD2	5:I:2420:GOL:H11	1.62	1.33
1:F:1098:ASP:CB	6:F:2109:HOH:O	1.68	1.31
1:H:1519[A]:HIS:CD2	5:H:2411:GOL:H12	1.72	1.25
5:B:2414:GOL:C1	6:B:2150:HOH:O	1.88	1.22
1:I:1602:ARG:N	6:I:2022:HOH:O	1.76	1.18

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	149/157 (95%)	146 (98%)	3 (2%)	0	100	100
1	B	148/157 (94%)	145 (98%)	3 (2%)	0	100	100
1	C	148/157 (94%)	145 (98%)	3 (2%)	0	100	100
1	D	149/157 (95%)	146 (98%)	3 (2%)	0	100	100
1	E	148/157 (94%)	144 (97%)	4 (3%)	0	100	100
1	F	148/157 (94%)	145 (98%)	3 (2%)	0	100	100
1	G	149/157 (95%)	145 (97%)	4 (3%)	0	100	100
1	H	148/157 (94%)	145 (98%)	3 (2%)	0	100	100
1	I	148/157 (94%)	145 (98%)	3 (2%)	0	100	100
1	J	148/157 (94%)	144 (97%)	4 (3%)	0	100	100
1	K	148/157 (94%)	145 (98%)	3 (2%)	0	100	100
1	L	149/157 (95%)	145 (97%)	4 (3%)	0	100	100
All	All	1780/1884 (94%)	1740 (98%)	40 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	119/121 (98%)	117 (98%)	2 (2%)	68	51
1	B	118/121 (98%)	114 (97%)	4 (3%)	44	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	118/121 (98%)	116 (98%)	2 (2%)	68	51
1	D	119/121 (98%)	116 (98%)	3 (2%)	55	34
1	E	118/121 (98%)	116 (98%)	2 (2%)	68	51
1	F	118/121 (98%)	117 (99%)	1 (1%)	86	79
1	G	119/121 (98%)	117 (98%)	2 (2%)	68	51
1	H	118/121 (98%)	116 (98%)	2 (2%)	68	51
1	I	118/121 (98%)	116 (98%)	2 (2%)	68	51
1	J	119/121 (98%)	117 (98%)	2 (2%)	68	51
1	K	118/121 (98%)	118 (100%)	0	100	100
1	L	119/121 (98%)	116 (98%)	3 (2%)	55	34
All	All	1421/1452 (98%)	1396 (98%)	25 (2%)	66	49

5 of 25 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	823	ARG
1	F	1026	GLU
1	L	2223	ARG
1	E	826	GLU
1	G	1211	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 26 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	1222	GLN
1	H	1422	GLN
1	L	2222	GLN
1	G	1247	HIS
1	G	1324	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

32 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$
2	CA2	A	160	-	20,23,23	1.42	3 (15%)	23,32,32	2.20	5 (21%)
3	PO4	A	2401	-	4,4,4	1.14	0	6,6,6	0.36	0
4	TRS	A	2405	-	7,7,7	0.99	1 (14%)	9,9,9	2.28	3 (33%)
5	GOL	A	2413	-	5,5,5	0.81	0	5,5,5	0.72	0
5	GOL	B	2414	-	5,5,5	0.54	0	5,5,5	0.47	0
2	CA2	B	360	-	20,23,23	1.03	1 (5%)	23,32,32	2.36	9 (39%)
5	GOL	C	2415	-	5,5,5	0.83	0	5,5,5	0.72	0
2	CA2	C	560	-	20,23,23	0.98	1 (5%)	23,32,32	1.53	4 (17%)
3	PO4	D	2402	-	4,4,4	1.01	0	6,6,6	0.28	0
5	GOL	D	2416	-	5,5,5	0.57	0	5,5,5	0.52	0
2	CA2	D	760	-	20,23,23	0.94	1 (5%)	23,32,32	2.42	7 (30%)
5	GOL	E	2419	-	5,5,5	0.75	0	5,5,5	0.83	0
5	GOL	E	2422	-	5,5,5	0.95	0	5,5,5	0.42	0
2	CA2	E	960	-	20,23,23	1.16	1 (5%)	23,32,32	2.11	7 (30%)
2	CA2	F	1160	-	20,23,23	1.04	1 (5%)	23,32,32	1.56	5 (21%)
4	TRS	F	2406	-	7,7,7	0.82	0	9,9,9	1.10	1 (11%)
5	GOL	F	2418	-	5,5,5	0.96	0	5,5,5	1.95	2 (40%)
2	CA2	G	1360	-	20,23,23	1.67	3 (15%)	23,32,32	2.43	6 (26%)
5	GOL	G	2417	-	5,5,5	0.76	0	5,5,5	1.41	1 (20%)
2	CA2	H	1560	-	20,23,23	1.21	2 (10%)	23,32,32	1.93	6 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	TRS	H	2407	-	7,7,7	1.29	0	9,9,9	1.24	1 (11%)
5	GOL	H	2411	-	5,5,5	0.38	0	5,5,5	0.54	0
2	CA2	I	1760	-	20,23,23	1.18	1 (5%)	23,32,32	1.59	5 (21%)
3	PO4	I	2403	-	4,4,4	0.97	0	6,6,6	0.34	0
5	GOL	I	2420	-	5,5,5	0.56	0	5,5,5	0.74	0
2	CA2	J	1960	-	20,23,23	1.35	3 (15%)	23,32,32	1.48	4 (17%)
5	GOL	J	2412	-	5,5,5	0.52	0	5,5,5	0.99	0
2	CA2	K	2160	-	20,23,23	0.64	0	23,32,32	2.04	5 (21%)
5	GOL	K	2421	-	5,5,5	0.67	0	5,5,5	0.87	0
2	CA2	L	2360	-	20,23,23	1.04	1 (5%)	23,32,32	1.66	6 (26%)
3	PO4	L	2404	-	4,4,4	0.67	0	6,6,6	0.41	0
4	TRS	L	2408	-	7,7,7	0.70	0	9,9,9	1.49	2 (22%)

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	CA2	A	160	-	-	0/7/31/31	0/2/2/2
3	PO4	A	2401	-	-	0/0/0/0	0/0/0/0
4	TRS	A	2405	-	-	0/9/9/9	0/0/0/0
5	GOL	A	2413	-	-	0/4/4/4	0/0/0/0
5	GOL	B	2414	-	-	0/4/4/4	0/0/0/0
2	CA2	B	360	-	-	0/7/31/31	0/2/2/2
5	GOL	C	2415	-	-	0/4/4/4	0/0/0/0
2	CA2	C	560	-	-	0/7/31/31	0/2/2/2
3	PO4	D	2402	-	-	0/0/0/0	0/0/0/0
5	GOL	D	2416	-	-	0/4/4/4	0/0/0/0
2	CA2	D	760	-	-	0/7/31/31	0/2/2/2
5	GOL	E	2419	-	-	0/4/4/4	0/0/0/0
5	GOL	E	2422	-	-	0/4/4/4	0/0/0/0
2	CA2	E	960	-	-	0/7/31/31	0/2/2/2
2	CA2	F	1160	-	-	0/7/31/31	0/2/2/2
4	TRS	F	2406	-	-	0/9/9/9	0/0/0/0
5	GOL	F	2418	-	-	0/4/4/4	0/0/0/0
2	CA2	G	1360	-	-	0/7/31/31	0/2/2/2
5	GOL	G	2417	-	-	0/4/4/4	0/0/0/0
2	CA2	H	1560	-	-	0/7/31/31	0/2/2/2
4	TRS	H	2407	-	-	0/9/9/9	0/0/0/0
5	GOL	H	2411	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CA2	I	1760	-	-	0/7/31/31	0/2/2/2
3	PO4	I	2403	-	-	0/0/0/0	0/0/0/0
5	GOL	I	2420	-	-	0/4/4/4	0/0/0/0
2	CA2	J	1960	-	-	0/7/31/31	0/2/2/2
5	GOL	J	2412	-	-	0/4/4/4	0/0/0/0
2	CA2	K	2160	-	-	0/7/31/31	0/2/2/2
5	GOL	K	2421	-	-	0/4/4/4	0/0/0/0
2	CA2	L	2360	-	-	0/7/31/31	0/2/2/2
3	PO4	L	2404	-	-	0/0/0/0	0/0/0/0
4	TRS	L	2408	-	-	0/9/9/9	0/0/0/0

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	1360	CA2	C2-C1	-5.42	1.48	1.53
2	I	1760	CA2	C6-C1	-3.99	1.50	1.53
2	B	360	CA2	C2-C1	-3.49	1.50	1.53
2	L	2360	CA2	C2-C1	-3.30	1.50	1.53
2	H	1560	CA2	C2-C1	-3.20	1.50	1.53

The worst 5 of 79 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	1360	CA2	C2-C3-C8	-7.50	102.23	112.11
2	K	2160	CA2	C6-C5-C4	-5.86	105.45	110.84
2	D	760	CA2	C2-C3-C8	-5.71	104.59	112.11
2	D	760	CA2	C6-C5-C4	-5.67	105.62	110.84
2	G	1360	CA2	C6-C5-C4	-4.51	106.70	110.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

22 monomers are involved in 118 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	160	CA2	1	0
3	A	2401	PO4	2	0
5	A	2413	GOL	1	0
5	B	2414	GOL	6	0
2	B	360	CA2	2	0
5	C	2415	GOL	9	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	2402	PO4	1	0
5	D	2416	GOL	9	0
5	E	2419	GOL	10	0
5	E	2422	GOL	5	0
2	E	960	CA2	2	0
2	F	1160	CA2	2	0
5	F	2418	GOL	13	0
2	G	1360	CA2	1	0
5	G	2417	GOL	7	0
2	H	1560	CA2	1	0
5	H	2411	GOL	7	0
3	I	2403	PO4	1	0
5	I	2420	GOL	21	0
2	J	1960	CA2	1	0
5	J	2412	GOL	4	0
5	K	2421	GOL	12	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.