



wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 10:26 AM GMT

PDB ID : 1F6N
Title : CRYSTAL STRUCTURE ANALYSIS OF THE MUTANT REACTION CENTER PRO L209-> TYR FROM THE PHOTOSYNTHETIC PURPLE BACTERIUM RHODOBACTER SPHAEROIDES
Authors : Kuglstatter, A.; Ermler, U.; Michel, H.; Baciou, L.; Fritzsche, G.
Deposited on : 2000-06-22
Resolution : 2.80 Å(reported)

This is a wwPDB 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/ValidationPDFNotes.html>

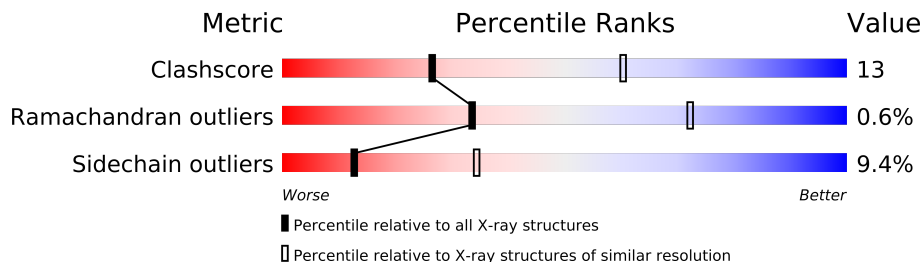
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	281	
2	M	307	
3	H	260	

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 7186 atoms, of which 0 are hydrogen and 0 are deuterium.

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 REACTION CENTER PROTEIN L CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	281	Total	C	N	O	S	0	0	0
			2237	1511	355	363	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	209	TYR	PRO	ENGINEERED	UNP P02954

- Molecule 2 is a protein called REACTION CENTER PROTEIN M CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	302	Total	C	N	O	S	0	0	0
			2404	1603	394	397	10			

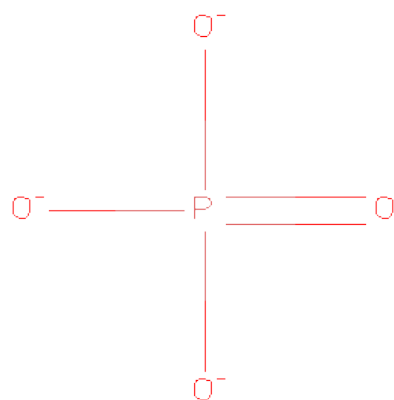
- Molecule 3 is a protein called REACTION CENTER PROTEIN H CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	240	Total	C	N	O	S	0	0	0
			1829	1169	314	337	9			

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

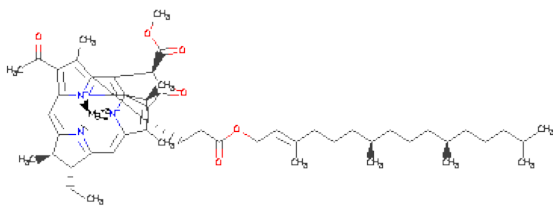
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	M	1	Total	Fe	0	0
			1	1		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

- Molecule 6 is BACTERIOCHLOROPHYLL A (three-letter code: BCL) (formula: C₅₅H₇₄MgN₄O₆).



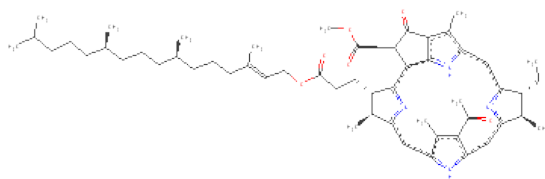
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Mg	N	O		
6	L	1	66	55	1	4	6	0	0
6	L	1	66	55	1	4	6	0	0
6	M	1	66	55	1	4	6	0	0

Continued on next page...

Continued from previous page...

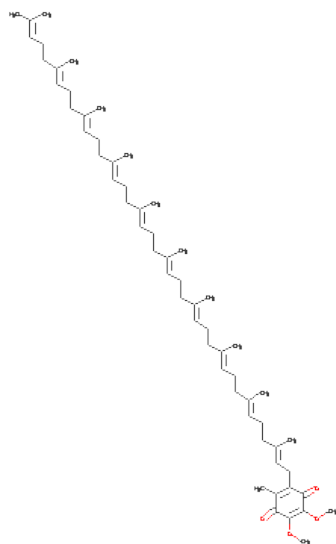
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

- Molecule 7 is BACTERIOPHEOPHYTIN A (three-letter code: BPH) (formula: $C_{55}H_{76}N_4O_6$).



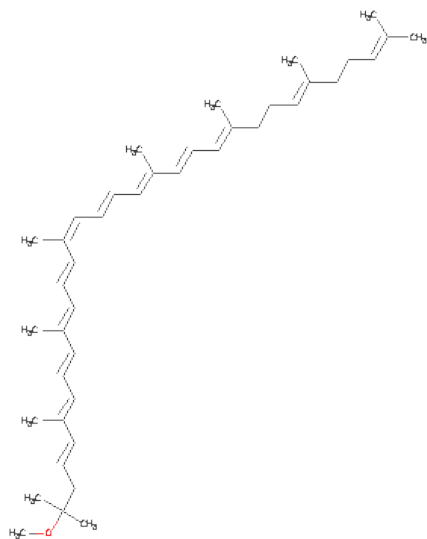
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	M	1	Total	C	N	O	0	0
			65	55	4	6		
7	L	1	Total	C	N	O	0	0
			65	55	4	6		

- Molecule 8 is UBIQUINONE-10 (three-letter code: U10) (formula: $C_{59}H_{90}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	M	1	Total	C	O	0	0
			48	44	4		
8	L	1	Total	C	O	0	0
			48	44	4		

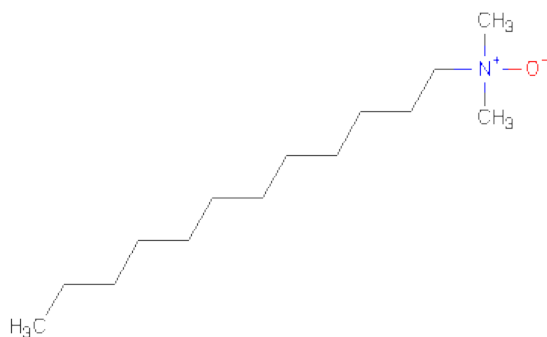
- Molecule 9 is SPHEROIDENE (three-letter code: SPO) (formula: $C_{41}H_{60}O$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	M	1	Total	C	O	0	0
			42	41	1		

- Molecule 10 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula:

C₁₄H₃₁NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	M	1	Total	C	N	O	0	0
			16	14	1	1		
10	H	1	Total	C	N	O	0	0
			16	14	1	1		
10	L	1	Total	C	N	O	0	0
			16	14	1	1		
10	M	1	Total	C	N	O	0	0
			16	14	1	1		
10	H	1	Total	C	N	O	0	0
			16	14	1	1		

- Molecule 11 is water.

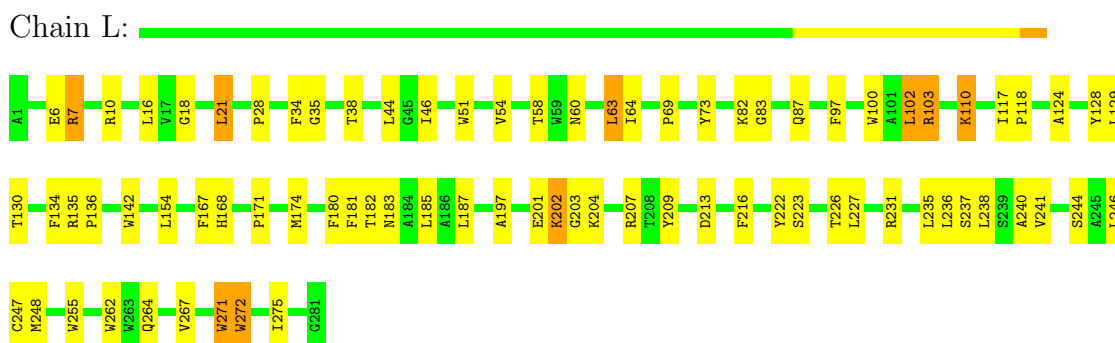
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	H	49	Total	O	0	0
			49	49		
11	L	22	Total	O	0	0
			22	22		
11	M	27	Total	O	0	0
			27	27		

3 Residue-property plots

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.

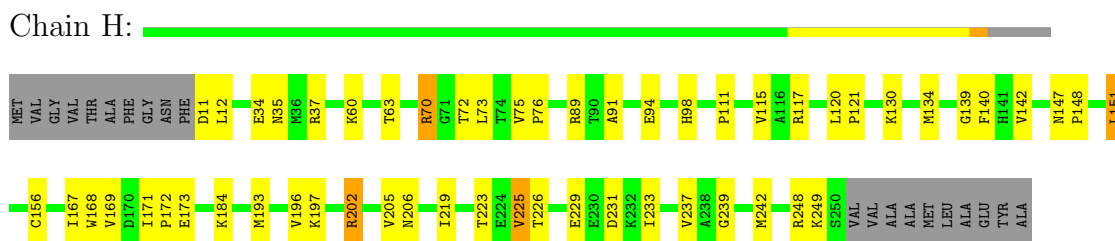
• Molecule 1: REACTION CENTER PROTEIN L CHAIN



• Molecule 2: REACTION CENTER PROTEIN M CHAIN



• Molecule 3: REACTION CENTER PROTEIN H CHAIN



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	141.53Å 141.53Å 187.91Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.80	Depositor
% Data completeness (in resolution range)	93.1 (50.00-2.80)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.3	Depositor
R, R_{free}	0.221 , 0.250	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7186	wwPDB-VP
Average B, all atoms (Å ²)	68.0	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: BCL, LDA, BPH, PO4, FE, SPO, U10

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	L	0.45	0/2325	0.51	0/3181
2	M	0.44	0/2496	0.51	1/3407 (0.0%)
3	H	0.43	0/1877	0.54	0/2553
All	All	0.44	0/6698	0.52	1/9141 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	L	0	3
2	M	0	1
All	All	0	4

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	214	LEU	CA-CB-CG	6.79	130.91	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	103	ARG	Sidechain
1	L	216	PHE	Sidechain
1	L	73	TYR	Sidechain
2	M	198	TYR	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	2237	0	2189	60	0
2	M	2404	0	2311	70	0
3	H	1829	0	1836	41	0
4	M	1	0	0	0	0
5	M	5	0	0	0	0
6	L	198	0	221	30	0
6	M	66	0	74	15	0
7	L	65	0	76	10	0
7	M	65	0	76	10	0
8	L	48	0	63	0	0
8	M	48	0	63	1	0
9	M	42	0	60	4	0
10	H	32	0	62	1	0
10	L	16	0	31	3	0
10	M	32	0	62	0	0
11	H	49	0	0	2	0
11	L	22	0	0	4	0
11	M	27	0	0	1	0
All	All	7186	0	7124	187	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

The worst 5 of 187 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
6:L:301:BCL:HHC	6:L:301:BCL:HBB2	1.40	1.02
7:L:402:BPH:HHC	7:L:402:BPH:HBB3	1.42	1.01
6:L:302:BCL:HBB3	6:L:302:BCL:HHC	1.49	0.93
2:M:153:ALA:HB2	7:M:401:BPH:HAC1	1.50	0.91
1:L:272:TRP:HA	1:L:275:ILE:HD13	1.59	0.83

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	L	279/281 (99%)	263 (94%)	15 (5%)	1 (0%)	43	80
2	M	300/307 (98%)	283 (94%)	13 (4%)	4 (1%)	18	51
3	H	238/260 (92%)	228 (96%)	10 (4%)	0	100	100
All	All	817/848 (96%)	774 (95%)	38 (5%)	5 (1%)	33	72

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	M	113	GLY
2	M	301	HIS
1	L	202	LYS
2	M	5	ASN
2	M	22	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	L	220/220 (100%)	199 (90%)	21 (10%)	12	33
2	M	235/240 (98%)	210 (89%)	25 (11%)	10	27
3	H	195/208 (94%)	180 (92%)	15 (8%)	18	45
All	All	650/668 (97%)	589 (91%)	61 (9%)	13	34

5 of 61 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	M	109	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	M	196	LEU
3	H	223	THR
2	M	133	THR
2	M	156	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
2	M	81	ASN
2	M	193	HIS
3	H	98	HIS
2	M	44	ASN
2	M	299	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 16 ligands modelled in this entry, 1 is monoatomic - leaving 15 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
10	LDA	H	702	-	15,15,15	5.05	2 (13%)	17,17,17	3.70	1 (5%)
10	LDA	H	705	-	15,15,15	4.47	2 (13%)	17,17,17	2.81	1 (5%)
6	BCL	L	301	2	74,74,74	1.56	11 (14%)	97,115,115	1.52	11 (11%)
6	BCL	L	302	1	74,74,74	1.58	12 (16%)	97,115,115	1.33	11 (11%)
6	BCL	L	304	1	74,74,74	1.93	14 (18%)	97,115,115	3.90	25 (25%)
7	BPH	L	402	-	70,70,70	1.47	9 (12%)	94,101,101	1.55	15 (15%)
8	U10	L	502	-	48,48,63	1.91	17 (35%)	59,61,79	1.18	8 (13%)
10	LDA	L	703	-	15,15,15	4.35	2 (13%)	17,17,17	4.54	1 (5%)
7	BPH	M	401	-	70,70,70	1.57	9 (12%)	94,101,101	1.60	17 (18%)
8	U10	M	501	-	48,48,63	2.20	16 (33%)	59,61,79	1.06	5 (8%)
9	SPO	M	600	-	41,41,41	3.12	23 (56%)	50,50,50	2.38	12 (24%)
10	LDA	M	701	-	15,15,15	4.52	3 (20%)	17,17,17	4.02	1 (5%)
10	LDA	M	704	-	15,15,15	4.55	4 (26%)	17,17,17	3.30	1 (5%)
5	PO4	M	800	-	4,4,4	1.07	0	6,6,6	0.31	0
6	BCL	M	801	2	74,74,74	1.47	11 (14%)	97,115,115	1.50	15 (15%)

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
10	LDA	H	702	-	-	0/13/13/13	0/0/0/0
10	LDA	H	705	-	-	0/13/13/13	0/0/0/0
6	BCL	L	301	2	1/1/21/25	0/41/137/137	0/0/9/9
6	BCL	L	302	1	-	0/41/137/137	0/0/9/9
6	BCL	L	304	1	1/1/21/25	0/41/137/137	0/0/9/9
7	BPH	L	402	-	2/2/18/22	0/49/105/105	0/0/6/6
8	U10	L	502	-	-	0/45/69/87	0/1/1/1
10	LDA	L	703	-	-	0/13/13/13	0/0/0/0
7	BPH	M	401	-	2/2/18/22	0/49/105/105	0/0/6/6
8	U10	M	501	-	-	0/45/69/87	0/1/1/1
9	SPO	M	600	-	-	0/47/47/47	0/0/0/0
10	LDA	M	701	-	-	0/13/13/13	0/0/0/0
10	LDA	M	704	-	-	0/13/13/13	0/0/0/0
5	PO4	M	800	-	-	0/0/0/0	0/0/0/0
6	BCL	M	801	2	-	0/41/137/137	0/0/9/9

The worst 5 of 135 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	H	702	LDA	O1-N1	-19.12	1.21	1.39
10	M	701	LDA	O1-N1	-17.07	1.23	1.39
10	H	705	LDA	O1-N1	-17.04	1.23	1.39
10	M	704	LDA	O1-N1	-16.65	1.23	1.39
10	L	703	LDA	O1-N1	-16.44	1.23	1.39

The worst 5 of 124 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	L	304	BCL	C1B-CHB-C4A	21.16	174.14	130.06
6	L	304	BCL	CHB-C4A-NA	-19.45	101.53	124.58
10	L	703	LDA	C2-C1-N1	18.60	145.85	113.80
10	M	701	LDA	C2-C1-N1	16.43	142.11	113.80
10	H	702	LDA	C2-C1-N1	15.11	139.85	113.80

5 of 6 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
7	M	401	BPH	C8
7	M	401	BPH	C13
6	L	304	BCL	C13
7	L	402	BPH	C8
7	L	402	BPH	C13

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

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

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

6.3 Carbohydrates ⓘ

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

6.4 Ligands ⓘ

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

6.5 Other polymers ⓘ

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