



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 03:01 AM UTC

PDB ID : 4OR0 / pdb_00004or0
Title : Crystal Structure of Bovine Serum Albumin in complex with naproxen
Authors : Zielinski, K.; Bujacz, A.; Sekula, B.; Bujacz, G.
Deposited on : 2014-02-10
Resolution : 2.58 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

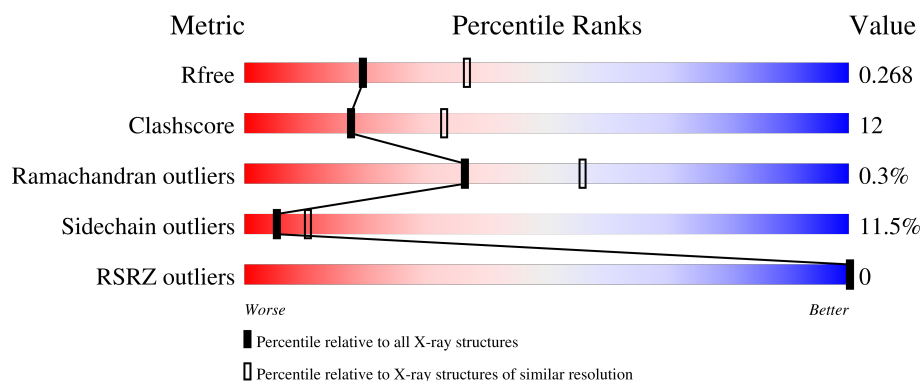
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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))
R_{free}	180053	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	583	70% 26% . .
1	B	583	64% 30% 6%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9508 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.

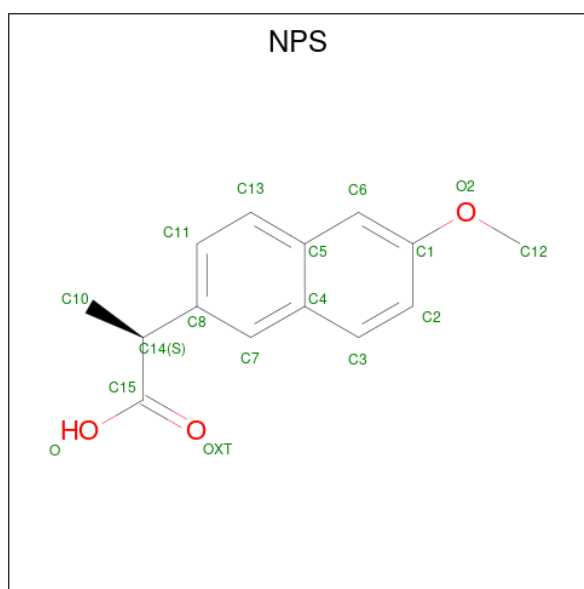
- Molecule 1 is a protein called Serum albumin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	582	Total	C	N	O	S	0	0	0
			4645	2931	780	895	39			
1	B	582	Total	C	N	O	S	0	0	0
			4645	2931	780	895	39			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	190	THR	ALA	variant	UNP P02769
B	190	THR	ALA	variant	UNP P02769

- Molecule 2 is (2S)-2-(6-methoxynaphthalen-2-yl)propanoic acid (CCD ID: NPS) (formula: C₁₄H₁₄O₃).



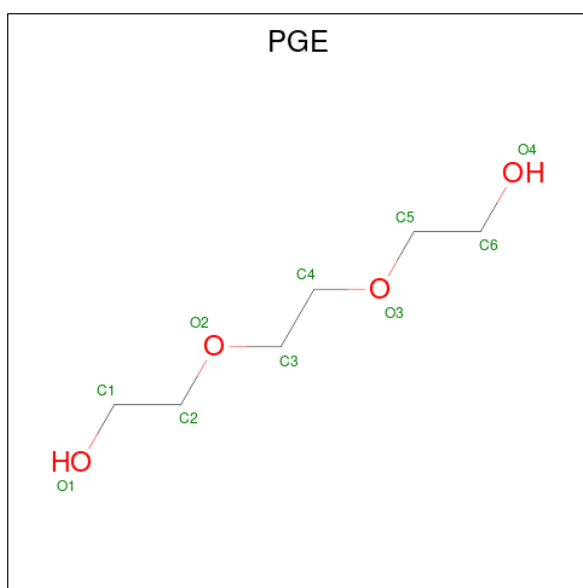
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			17	14	3		

Continued on next page...

Continued from previous page...

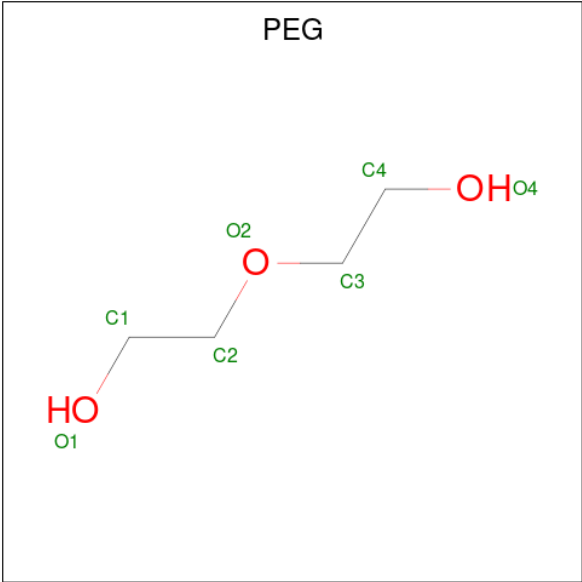
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			17	14	3		
2	A	1	Total	C	O	0	0
			17	14	3		
2	B	1	Total	C	O	0	0
			17	14	3		
2	B	1	Total	C	O	0	0
			17	14	3		
2	B	1	Total	C	O	0	0
			17	14	3		

- Molecule 3 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		

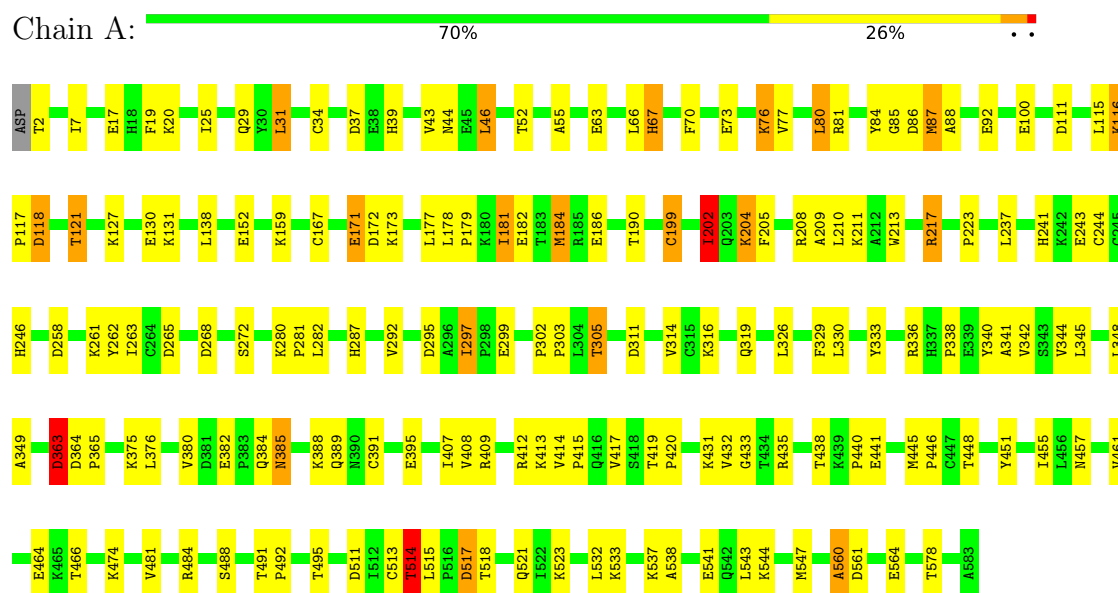
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	54	Total	O	0	0
			54	54		
5	B	45	Total	O	0	0
			45	45		

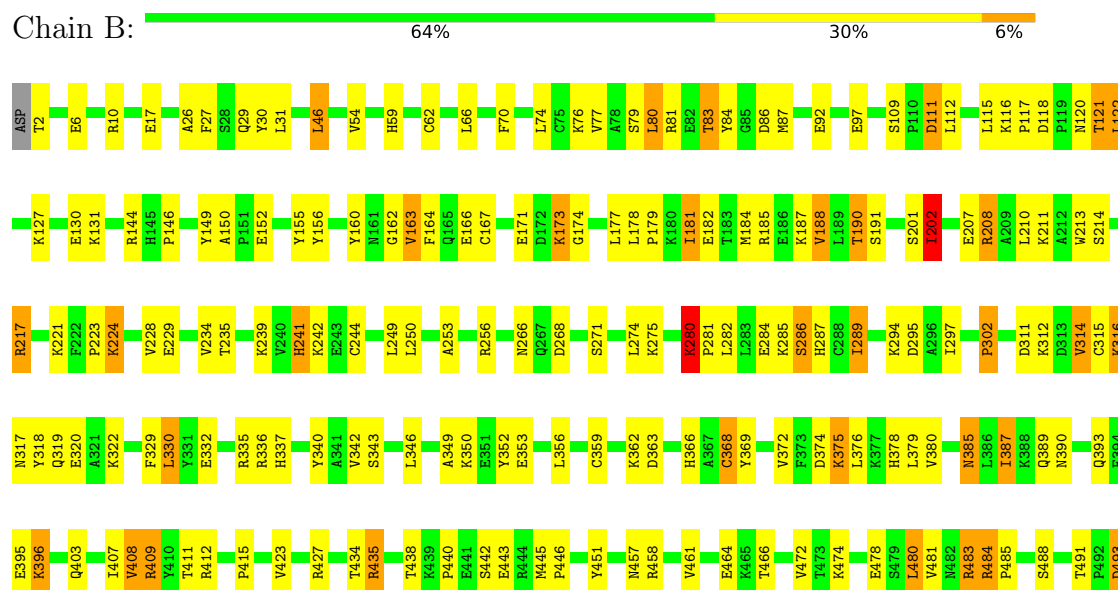
3 Residue-property plots

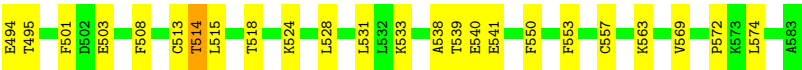
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Serum albumin



• Molecule 1: Serum albumin





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	216.04Å 44.90Å 141.70Å 90.00° 113.99° 90.00°	Depositor
Resolution (Å)	42.28 – 2.58 42.28 – 2.58	Depositor EDS
% Data completeness (in resolution range)	99.1 (42.28-2.58) 99.1 (42.28-2.58)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.56Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.183 , 0.257 (Not available) , 0.268	Depositor DCC
R_{free} test set	1244 reflections (3.11%)	wwPDB-VP
Wilson B-factor (Å ²)	58.1	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 26.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9508	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NPS, PGE, PEG

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	1.25	7/4740 (0.1%)	1.34	26/6399 (0.4%)
1	B	1.15	5/4740 (0.1%)	1.31	23/6399 (0.4%)
All	All	1.20	12/9480 (0.1%)	1.33	49/12798 (0.4%)

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	457	ASN	N-CA	6.79	1.54	1.46
1	B	461	VAL	C-O	-6.14	1.17	1.24
1	A	287	HIS	N-CA	-5.95	1.39	1.46
1	A	67	HIS	N-CA	5.82	1.53	1.46
1	A	209	ALA	C-O	5.81	1.31	1.24

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	480	LEU	N-CA-C	-9.69	101.97	113.88
1	A	202	ILE	CB-CA-C	-8.58	100.99	111.97
1	A	455	ILE	N-CA-C	7.73	117.80	110.53
1	A	391	CYS	N-CA-C	7.47	119.42	111.28
1	A	363	ASP	N-CA-C	7.13	119.13	111.36

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	4645	0	4559	90	0
1	B	4645	0	4559	137	0
2	A	51	0	39	2	0
2	B	51	0	39	7	0
3	A	10	0	13	0	0
4	A	7	0	10	2	0
5	A	54	0	0	11	0
5	B	45	0	0	13	0
All	All	9508	0	9219	230	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 12.

The worst 5 of 230 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:ILE:HG22	5:A:732:HOH:O	1.57	1.04
1:B:474:LYS:O	1:B:478:GLU:HG2	1.72	0.90
1:B:208:ARG:NH1	2:B:602:NPS:H2	1.88	0.89
1:B:202:ILE:CD1	1:B:210:LEU:HD22	2.06	0.86
1:B:202:ILE:HD11	1:B:210:LEU:HD22	1.57	0.84

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	580/583 (100%)	545 (94%)	33 (6%)	2 (0%)	36	56
1	B	580/583 (100%)	541 (93%)	38 (7%)	1 (0%)	43	63

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1160/1166 (100%)	1086 (94%)	71 (6%)	3 (0%)	36	56

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	86	ASP
1	A	265	ASP
1	B	302	PRO

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	520/521 (100%)	467 (90%)	53 (10%)	7	14
1	B	520/521 (100%)	453 (87%)	67 (13%)	4	8
All	All	1040/1042 (100%)	920 (88%)	120 (12%)	5	10

5 of 120 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	79	SER
1	B	466	THR
1	B	181	ILE
1	B	464	GLU
1	B	518	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	319	GLN
1	B	378	HIS
1	A	482	ASN
1	A	509	HIS
1	B	3	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	NPS	B	602	-	18,18,18	1.31	5 (27%)	25,25,25	1.42	6 (24%)
2	NPS	B	603	-	18,18,18	1.41	5 (27%)	25,25,25	1.89	8 (32%)
2	NPS	A	602	-	18,18,18	1.10	1 (5%)	25,25,25	1.50	4 (16%)
2	NPS	A	601	-	18,18,18	1.15	1 (5%)	25,25,25	1.47	5 (20%)
3	PGE	A	604	-	9,9,9	0.54	0	8,8,8	1.37	1 (12%)
4	PEG	A	605	-	6,6,6	0.44	0	5,5,5	1.54	1 (20%)
2	NPS	B	601	-	18,18,18	1.43	2 (11%)	25,25,25	1.43	3 (12%)
2	NPS	A	603	-	18,18,18	1.59	3 (16%)	25,25,25	2.05	10 (40%)

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	NPS	B	602	-	-	5/10/10/10	0/2/2/2
2	NPS	B	603	-	-	4/10/10/10	0/2/2/2
2	NPS	A	602	-	-	6/10/10/10	0/2/2/2
2	NPS	A	601	-	-	2/10/10/10	0/2/2/2
3	PGE	A	604	-	-	4/7/7/7	-
4	PEG	A	605	-	-	3/4/4/4	-
2	NPS	B	601	-	-	2/10/10/10	0/2/2/2
2	NPS	A	603	-	-	6/10/10/10	0/2/2/2

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	603	NPS	C8-C14	-3.13	1.47	1.52
2	A	603	NPS	C6-C1	3.00	1.42	1.37
2	B	601	NPS	C6-C1	2.94	1.42	1.37
2	A	602	NPS	C6-C1	2.58	1.41	1.37
2	B	603	NPS	C8-C14	-2.40	1.48	1.52

The worst 5 of 38 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	603	NPS	C6-C5-C4	3.86	124.32	118.96
2	B	603	NPS	C6-C5-C4	3.75	124.16	118.96
2	A	603	NPS	C8-C7-C4	-3.61	115.99	121.47
2	A	602	NPS	C7-C4-C5	3.32	123.57	118.96
2	B	603	NPS	O-C15-OXT	-3.28	116.63	124.08

There are no chirality outliers.

5 of 32 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	603	NPS	C6-C1-O2-C12
2	B	602	NPS	C6-C1-O2-C12
2	B	603	NPS	C2-C1-O2-C12
2	B	602	NPS	C2-C1-O2-C12
4	A	605	PEG	C4-C3-O2-C2

There are no ring outliers.

6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	602	NPS	3	0
2	B	603	NPS	3	0
2	A	601	NPS	1	0
4	A	605	PEG	2	0
2	B	601	NPS	2	0
2	A	603	NPS	1	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](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	582/583 (99%)	-0.68	0 100 100	33, 59, 105, 139	0
1	B	582/583 (99%)	-0.57	0 100 100	40, 68, 112, 136	0
All	All	1164/1166 (99%)	-0.63	0 100 100	33, 64, 110, 139	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	PEG	A	605	7/7	0.86	0.16	59,74,82,90	0
2	NPS	B	602	17/17	0.92	0.10	71,82,91,97	0
2	NPS	A	602	17/17	0.93	0.07	62,71,83,84	0
2	NPS	A	603	17/17	0.94	0.10	56,71,85,105	0
2	NPS	B	603	17/17	0.94	0.09	55,69,104,106	0
3	PGE	A	604	10/10	0.94	0.09	59,73,85,88	0
2	NPS	B	601	17/17	0.94	0.09	57,70,94,99	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NPS	A	601	17/17	0.95	0.08	50,57,79,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.