



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 28, 2023 – 10:59 PM EDT

PDB ID : 3PRY  
Title : Crystal structure of the middle domain of human HSP90-beta refined at 2.3 Å resolution  
Authors : Chaikuad, A.; Pilka, E.; Sharpe, T.D.; Cooper, C.D.O.; Phillips, C.; Berridge, G.; Ayinampudi, V.; Fedorov, O.; Keates, T.; Thangaratnarajah, C.; Zimmermann, T.; Vollmar, M.; Yue, W.W.; Che, K.H.; Krojer, T.; Muniz, J.R.C.; von Delft, F.; Bountra, C.; Arrowsmith, C.H.; Weigelt, J.; Edwards, A.; Bullock, A.; Structural Genomics Consortium (SGC)  
Deposited on : 2010-11-30  
Resolution : 2.28 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](mailto: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>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)

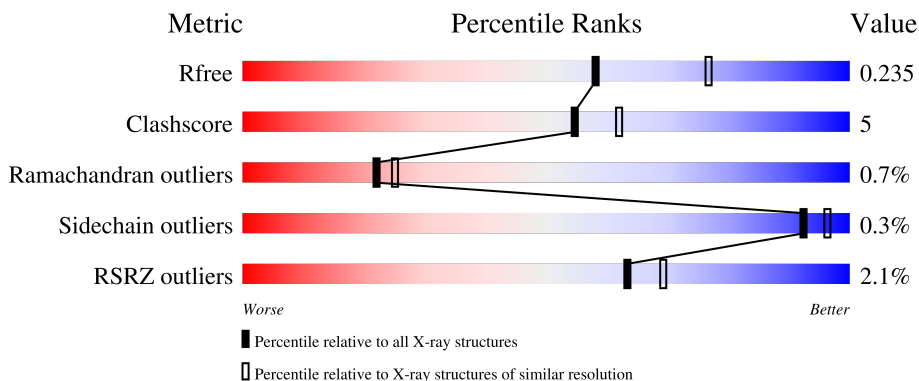
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.28 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))
$R_{free}$	130704	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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  $\geq 3$ , 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	268	 2% 81% 13% 5%
1	B	268	 % 81% 13% 6%
1	C	268	 3% 81% 11% 7%

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35

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
2	SO4	B	3	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6396 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 Heat shock protein HSP 90-beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	255	2074	1319	349	398	8	0	1	0
1	B	251	2028	1292	345	383	8	0	2	0
1	C	250	2038	1295	345	390	8	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

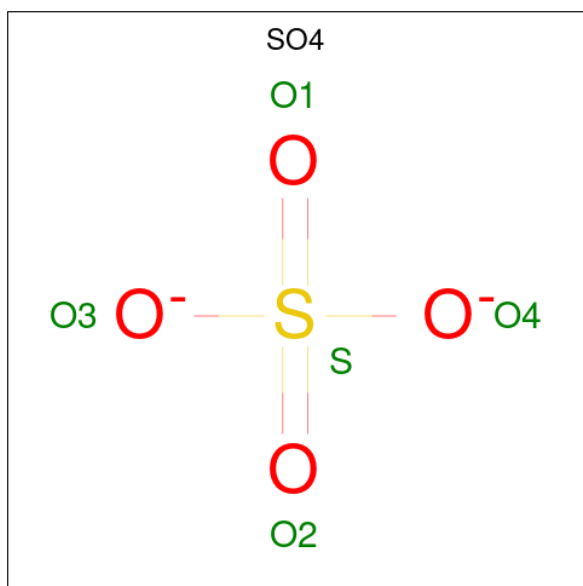
Chain	Residue	Modelled	Actual	Comment	Reference
A	283	MET	-	initiating methionine	UNP P08238
A	544	ALA	-	expression tag	UNP P08238
A	545	GLU	-	expression tag	UNP P08238
A	546	ASN	-	expression tag	UNP P08238
A	547	LEU	-	expression tag	UNP P08238
A	548	TYR	-	expression tag	UNP P08238
A	549	PHE	-	expression tag	UNP P08238
A	550	GLN	-	expression tag	UNP P08238
B	283	MET	-	initiating methionine	UNP P08238
B	544	ALA	-	expression tag	UNP P08238
B	545	GLU	-	expression tag	UNP P08238
B	546	ASN	-	expression tag	UNP P08238
B	547	LEU	-	expression tag	UNP P08238
B	548	TYR	-	expression tag	UNP P08238
B	549	PHE	-	expression tag	UNP P08238
B	550	GLN	-	expression tag	UNP P08238
C	283	MET	-	initiating methionine	UNP P08238
C	544	ALA	-	expression tag	UNP P08238
C	545	GLU	-	expression tag	UNP P08238
C	546	ASN	-	expression tag	UNP P08238
C	547	LEU	-	expression tag	UNP P08238
C	548	TYR	-	expression tag	UNP P08238
C	549	PHE	-	expression tag	UNP P08238

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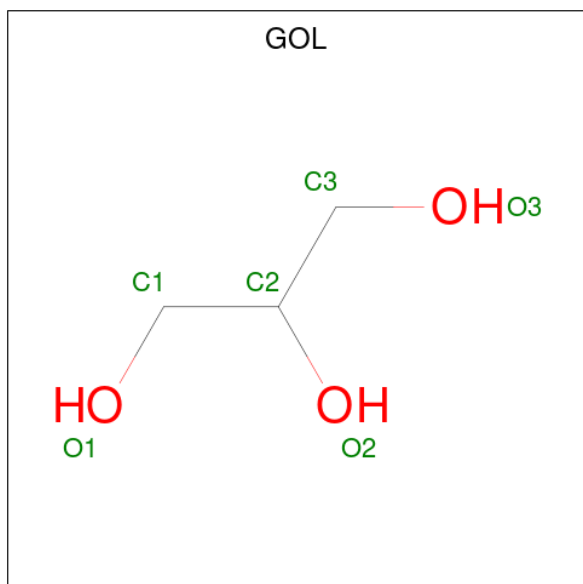
Chain	Residue	Modelled	Actual	Comment	Reference
C	550	GLN	-	expression tag	UNP P08238

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



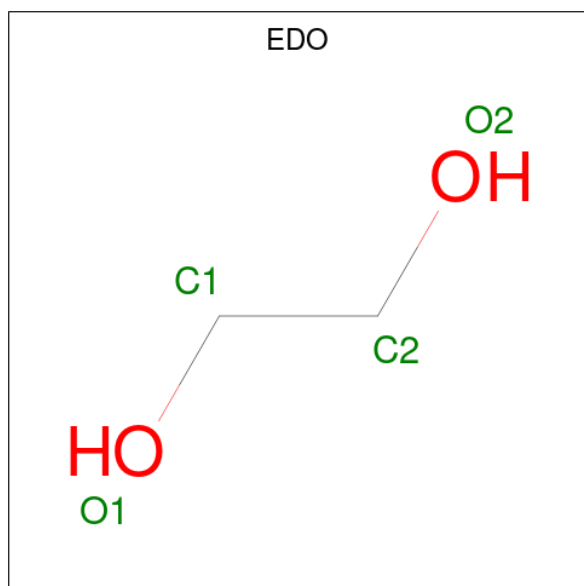
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



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

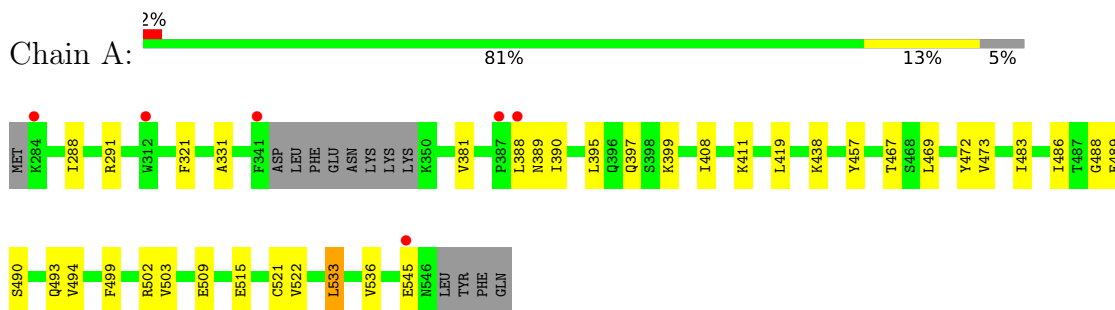
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	88	Total	O	0	0
			88	88		
5	B	67	Total	O	0	0
			67	67		
5	C	65	Total	O	0	0
			65	65		

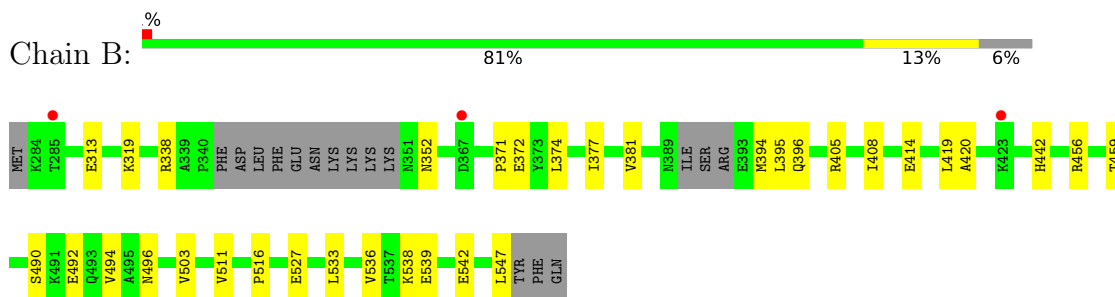
### 3 Residue-property plots [i](#)

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 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.

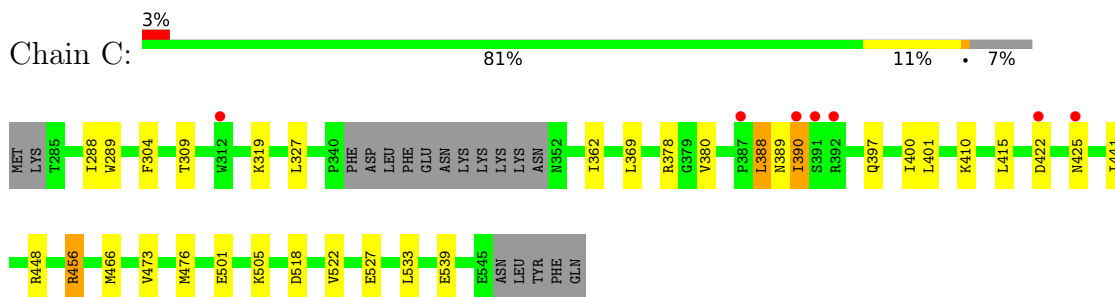
- Molecule 1: Heat shock protein HSP 90-beta



- Molecule 1: Heat shock protein HSP 90-beta



- Molecule 1: Heat shock protein HSP 90-beta



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.71Å 73.71Å 149.05Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.19 – 2.28 35.78 – 2.28	Depositor EDS
% Data completeness (in resolution range)	99.6 (39.19-2.28) 100.0 (35.78-2.28)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.64 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.172 , 0.217 0.191 , 0.235	Depositor DCC
$R_{free}$ test set	2027 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.6	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 59.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.005 for -h,-k,l 0.049 for h,-h-k,-l 0.004 for -k,-h,-l	Xtriage
Reported twinning fraction	0.675 for H, K, L 0.325 for -H, H+K, -L	Depositor
Outliers	0 of 41275 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6396	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, SO4, EDO

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	0.71	0/2115	0.73	0/2850
1	B	0.65	0/2069	0.67	0/2784
1	C	0.61	0/2075	0.70	1/2794 (0.0%)
All	All	0.66	0/6259	0.70	1/8428 (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	A	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	456	ARG	NE-CZ-NH1	5.17	122.88	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	545	GLU	Peptide

### 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	2074	0	2012	22	0
1	B	2028	0	1984	23	0
1	C	2038	0	1996	20	0
2	A	5	0	0	0	0
2	B	5	0	0	2	0
3	B	6	0	8	2	0
4	B	16	0	24	2	0
4	C	4	0	6	0	0
5	A	88	0	0	1	0
5	B	67	0	0	2	0
5	C	65	0	0	1	0
All	All	6396	0	6030	64	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 5.

All (64) 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:388:LEU:N	1:A:389:ASN:HA	2.04	0.72
1:C:441:ILE:HG23	1:C:448:ARG:HG3	1.77	0.66
1:C:362:ILE:HD12	1:C:388:LEU:HD13	1.81	0.61
1:B:319:LYS:HZ2	3:B:4:GOL:H32	1.67	0.59
1:C:397:GLN:O	1:C:401:LEU:HG	2.02	0.59
1:B:319:LYS:NZ	3:B:4:GOL:H32	2.17	0.59
1:A:288:ILE:HA	1:A:291:ARG:HD3	1.86	0.58
1:B:372:GLU:HG2	4:B:6:EDO:H12	1.84	0.58
1:B:527:GLU:HA	1:B:533:LEU:HD23	1.86	0.57
1:C:518:ASP:O	1:C:522:VAL:HG12	2.05	0.56
1:A:503:VAL:HG21	1:A:536:VAL:HG12	1.88	0.55
1:A:489:GLU:OE1	1:A:493:GLN:HG3	2.08	0.54
1:C:289:TRP:CE3	1:C:380:VAL:HG21	2.44	0.53
1:B:405:ARG:NH1	4:B:6:EDO:O2	2.42	0.53
1:C:309:THR:HG22	1:C:378:ARG:NH1	2.24	0.53
1:B:313:GLU:OE1	1:B:338:ARG:NH1	2.41	0.52
1:B:496:ASN:HB2	5:B:125:HOH:O	2.07	0.52
1:C:319:LYS:HB2	1:C:415:LEU:HD22	1.91	0.52
1:A:397:GLN:HE22	1:B:516:PRO:HB3	1.74	0.52
1:B:503:VAL:HG21	1:B:536:VAL:HG12	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:492:GLU:N	1:B:492:GLU:OE1	2.42	0.51
1:A:490:SER:O	1:A:494:VAL:HG12	2.10	0.51
1:C:473:VAL:O	1:C:476:MET:HG2	2.11	0.51
1:B:538:LYS:NZ	2:B:3:SO4:O3	2.38	0.51
1:B:459:THR:HG22	1:B:511:VAL:HA	1.93	0.51
1:C:539:GLU:H	1:C:539:GLU:CD	2.13	0.51
1:C:288:ILE:HG21	1:C:304:PHE:CD1	2.46	0.50
1:A:381:VAL:HG21	1:A:408:ILE:HD11	1.92	0.50
1:B:442:HIS:O	1:B:442:HIS:ND1	2.45	0.50
1:B:539:GLU:HB3	2:B:3:SO4:O4	2.12	0.49
1:A:483:ILE:HB	1:A:533:LEU:HD13	1.95	0.48
1:B:381:VAL:HG21	1:B:408:ILE:HD11	1.96	0.48
1:A:457:TYR:O	1:A:467:THR:HG22	2.14	0.48
1:C:501:GLU:HG3	1:C:505:LYS:HE3	1.95	0.48
1:A:321:PHE:CD2	1:A:411:LYS:HG2	2.48	0.47
1:B:371:PRO:HD2	1:B:374:LEU:HD12	1.96	0.47
1:A:395:LEU:HD11	1:A:399:LYS:HE3	1.97	0.47
1:A:321:PHE:CE2	1:A:331:ALA:HB3	2.49	0.47
1:B:456:ARG:O	5:B:80:HOH:O	2.21	0.47
1:C:327:LEU:HD21	1:C:400:ILE:HG23	1.98	0.45
1:A:469:LEU:O	1:A:473:VAL:HG23	2.17	0.45
1:B:542:GLU:OE2	1:B:547:LEU:HD11	2.17	0.45
1:A:472:TYR:CE1	1:A:509:GLU:HB3	2.52	0.44
1:A:515:GLU:OE2	5:A:97:HOH:O	2.21	0.44
1:C:456:ARG:HB3	1:C:466:MET:HB3	2.00	0.43
1:C:369:LEU:HD23	1:C:369:LEU:HA	1.83	0.43
1:C:422:ASP:OD1	1:C:425:ASN:HB2	2.18	0.43
1:B:419:LEU:HD12	1:B:419:LEU:HA	1.77	0.43
1:A:438:LYS:HG2	1:A:521:CYS:SG	2.59	0.43
1:C:527:GLU:HA	1:C:533:LEU:HD23	2.01	0.42
1:A:389:ASN:OD1	1:A:389:ASN:C	2.58	0.42
1:A:522:VAL:HG23	1:A:533:LEU:HG	2.01	0.42
1:A:419:LEU:HD12	1:A:419:LEU:HA	1.93	0.42
1:B:394:MET:C	1:B:396:GLN:H	2.22	0.42
1:C:539:GLU:N	1:C:539:GLU:OE1	2.53	0.41
1:B:352:ASN:HB2	1:B:377:ILE:O	2.20	0.41
1:A:488:GLY:C	1:A:489:GLU:HG2	2.40	0.41
1:B:490:SER:O	1:B:494:VAL:HG22	2.21	0.41
1:A:486:ILE:HD11	1:A:489:GLU:CD	2.41	0.41
1:B:319:LYS:NZ	1:B:414:GLU:OE2	2.52	0.41
1:A:499:PHE:O	1:A:502:ARG:HG2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:388:LEU:O	1:C:390:ILE:N	2.51	0.40
1:C:410:LYS:HE3	5:C:168:HOH:O	2.22	0.40
1:C:533:LEU:HD22	1:C:533:LEU:N	2.36	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	252/268 (94%)	245 (97%)	6 (2%)	1 (0%)	34	40
1	B	247/268 (92%)	236 (96%)	9 (4%)	2 (1%)	19	22
1	C	246/268 (92%)	230 (94%)	14 (6%)	2 (1%)	19	22
All	All	745/804 (93%)	711 (95%)	29 (4%)	5 (1%)	22	25

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	420	ALA
1	C	390	ILE
1	C	389	ASN
1	B	395	LEU
1	A	390	ILE

### 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	225/251 (90%)	224 (100%)	1 (0%)	91	95
1	B	218/251 (87%)	218 (100%)	0	100	100
1	C	223/251 (89%)	222 (100%)	1 (0%)	91	95
All	All	666/753 (88%)	664 (100%)	2 (0%)	92	96

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	533	LEU
1	C	388	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides 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
4	EDO	B	6	-	3,3,3	0.47	0	2,2,2	0.67	0
2	SO4	B	3	-	4,4,4	0.16	0	6,6,6	0.25	0
4	EDO	B	5	-	3,3,3	0.48	0	2,2,2	0.43	0
4	EDO	C	9	-	3,3,3	0.75	0	2,2,2	0.24	0
4	EDO	B	7	-	3,3,3	0.33	0	2,2,2	1.24	0
2	SO4	A	1	-	4,4,4	0.21	0	6,6,6	0.18	0
3	GOL	B	4	-	5,5,5	0.36	0	5,5,5	0.68	0
4	EDO	B	8	-	3,3,3	0.58	0	2,2,2	0.43	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
4	EDO	B	6	-	-	1/1/1/1	-
4	EDO	B	5	-	-	1/1/1/1	-
4	EDO	C	9	-	-	1/1/1/1	-
4	EDO	B	7	-	-	1/1/1/1	-
3	GOL	B	4	-	-	2/4/4/4	-
4	EDO	B	8	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	4	GOL	C1-C2-C3-O3
3	B	4	GOL	O2-C2-C3-O3
4	B	5	EDO	O1-C1-C2-O2
4	B	8	EDO	O1-C1-C2-O2
4	C	9	EDO	O1-C1-C2-O2
4	B	7	EDO	O1-C1-C2-O2
4	B	6	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	6	EDO	2	0
2	B	3	SO4	2	0
3	B	4	GOL	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	255/268 (95%)	-0.04	6 (2%) 59 65	23, 41, 67, 87	0
1	B	251/268 (93%)	-0.03	3 (1%) 79 82	28, 44, 67, 86	0
1	C	250/268 (93%)	0.13	7 (2%) 53 59	28, 46, 74, 111	0
All	All	756/804 (94%)	0.02	16 (2%) 63 69	23, 44, 70, 111	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	392	ARG	7.9
1	C	391	SER	4.8
1	C	390	ILE	4.8
1	A	388	LEU	4.2
1	C	312	TRP	3.7
1	B	285	THR	3.5
1	C	425	ASN	3.5
1	B	367	ASP	2.8
1	A	341	PHE	2.6
1	A	545	GLU	2.5
1	A	312	TRP	2.4
1	B	423	LYS	2.4
1	A	387	PRO	2.3
1	A	284	LYS	2.2
1	C	387	PRO	2.1
1	C	422	ASP	2.0

### 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 monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	GOL	B	4	6/6	0.86	0.14	37,40,61,61	0
4	EDO	B	6	4/4	0.86	0.20	29,33,34,35	0
4	EDO	C	9	4/4	0.92	0.21	29,31,39,43	0
4	EDO	B	8	4/4	0.93	0.25	36,40,41,44	0
4	EDO	B	7	4/4	0.94	0.15	27,29,34,34	0
2	SO4	A	1	5/5	0.97	0.07	37,47,54,59	0
4	EDO	B	5	4/4	0.97	0.16	40,40,44,46	0
2	SO4	B	3	5/5	0.98	0.07	35,48,53,55	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.