



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:27 PM UTC

PDB ID : 6EZT / pdb_00006ezt
Title : Crystal structure of GH20 Exo beta-N-Acetylglucosaminidase D437A inactive mutant from *Vibrio harveyi*
Authors : Porfetye, A.T.; Meekrathok, P.; Burger, M.; Vetter, I.R.; Suginta, W.
Deposited on : 2017-11-16
Resolution : 2.60 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)	:	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

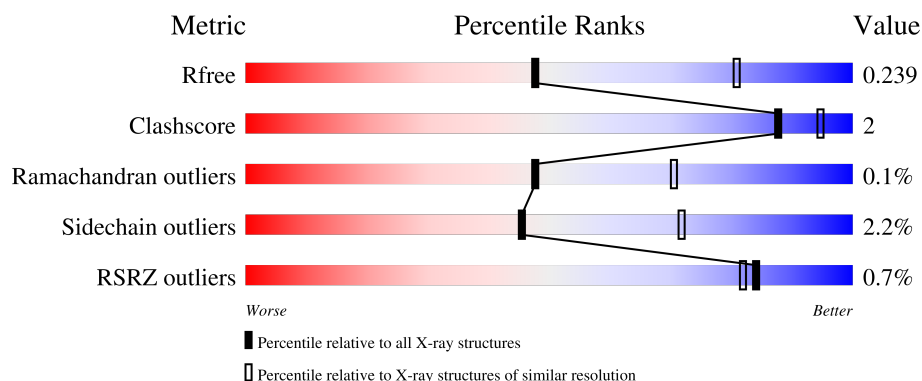
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	649	% 91% 7% .
1	B	649	% 91% 7% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10764 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 Beta-N-acetylglucosaminidase Nag2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	639	Total	C	N	O	S	0	1	0
			5145	3275	879	969	22			
1	B	639	Total	C	N	O	S	0	2	0
			5155	3281	881	971	22			

There are 22 discrepancies between the modelled and reference sequences:

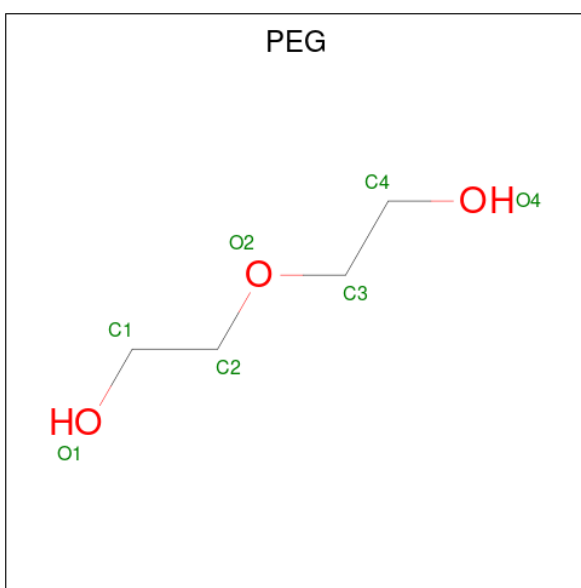
Chain	Residue	Modelled	Actual	Comment	Reference
A	437	ALA	ASP	engineered mutation	UNP D9ISE0
A	643	ARG	-	expression tag	UNP D9ISE0
A	644	SER	-	expression tag	UNP D9ISE0
A	645	ARG	-	expression tag	UNP D9ISE0
A	646	SER	-	expression tag	UNP D9ISE0
A	647	HIS	-	expression tag	UNP D9ISE0
A	648	HIS	-	expression tag	UNP D9ISE0
A	649	HIS	-	expression tag	UNP D9ISE0
A	650	HIS	-	expression tag	UNP D9ISE0
A	651	HIS	-	expression tag	UNP D9ISE0
A	652	HIS	-	expression tag	UNP D9ISE0
B	437	ALA	ASP	engineered mutation	UNP D9ISE0
B	643	ARG	-	expression tag	UNP D9ISE0
B	644	SER	-	expression tag	UNP D9ISE0
B	645	ARG	-	expression tag	UNP D9ISE0
B	646	SER	-	expression tag	UNP D9ISE0
B	647	HIS	-	expression tag	UNP D9ISE0
B	648	HIS	-	expression tag	UNP D9ISE0
B	649	HIS	-	expression tag	UNP D9ISE0
B	650	HIS	-	expression tag	UNP D9ISE0
B	651	HIS	-	expression tag	UNP D9ISE0
B	652	HIS	-	expression tag	UNP D9ISE0

- Molecule 2 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	8	5		
2	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

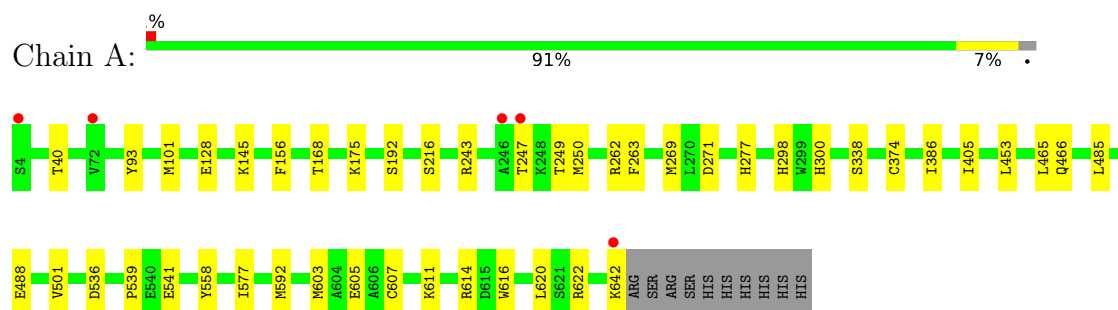
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	215	Total 215	O 215	0	0
4	B	209	Total 209	O 209	0	0

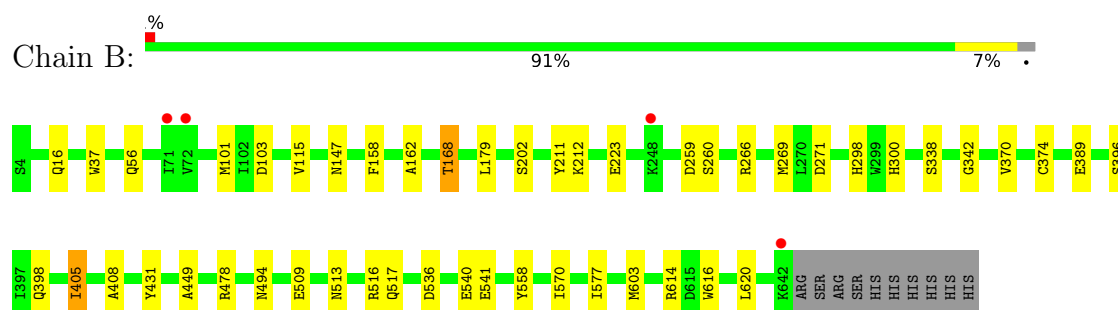
3 Residue-property plots [i](#)

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: Beta-N-acetylglucosaminidase Nag2



• Molecule 1: Beta-N-acetylglucosaminidase Nag2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.42Å 129.24Å 98.36Å 90.00° 112.20° 90.00°	Depositor
Resolution (Å)	91.09 – 2.60 91.07 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (91.09-2.60) 99.9 (91.07-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.195 , 0.236 0.200 , 0.239	Depositor DCC
R_{free} test set	3180 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.661	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10764	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.06% 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: PEG, PG4

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	0.52	0/5285	0.78	0/7187
1	B	0.53	0/5295	0.79	1/7200 (0.0%)
All	All	0.53	0/10580	0.78	1/14387 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	342	GLY	N-CA-C	5.58	118.12	110.58

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	5145	0	4985	21	0
1	B	5155	0	4994	19	0
2	A	13	0	18	0	0
2	B	13	0	18	0	0
3	A	7	0	10	0	0
3	B	7	0	10	1	0
4	A	215	0	0	1	0
4	B	209	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	10764	0	10035	40	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 2.

All (40) 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:101:MET:HE1	1:A:338:SER:HA	1.81	0.61
1:B:509:GLU:OE2	1:B:516:ARG:NH2	2.36	0.59
1:B:558:TYR:O	1:B:614:ARG:NH2	2.40	0.54
1:A:466:GLN:NE2	4:A:802:HOH:O	2.41	0.52
1:B:271:ASP:HA	1:B:300:HIS:HB3	1.91	0.51
1:B:266:ARG:HH12	3:B:701:PEG:H22	1.75	0.50
1:A:269:MET:HG3	1:A:298:HIS:CG	2.46	0.50
1:A:536:ASP:HB3	1:A:541:GLU:HG3	1.94	0.50
1:B:211:TYR:CE1	1:B:259:ASP:HB3	2.46	0.50
1:B:101:MET:HE1	1:B:338:SER:HA	1.94	0.49
1:B:269:MET:HG3	1:B:298:HIS:CG	2.47	0.49
1:A:616:TRP:CZ2	1:A:620:LEU:HD11	2.47	0.49
1:A:93:TYR:CE2	1:A:539:PRO:HB2	2.48	0.48
1:B:212:LYS:HB2	1:B:223:GLU:HB2	1.96	0.48
1:B:536:ASP:HB3	1:B:541:GLU:HG3	1.95	0.48
1:A:271:ASP:HA	1:A:300:HIS:HB3	1.96	0.47
1:B:370:VAL:HG12	1:B:405:ILE:HD12	1.97	0.46
1:B:603:MET:HA	1:B:603:MET:HE2	1.98	0.46
1:A:262:ARG:HG2	1:A:263:PHE:CD2	2.52	0.45
1:A:247:THR:HG22	1:A:249:THR:HG23	1.98	0.44
1:A:269:MET:HG3	1:A:298:HIS:CD2	2.52	0.44
1:A:156:PHE:HB2	1:A:216:SER:O	2.17	0.44
1:A:277:HIS:CG	1:A:592:MET:HE1	2.53	0.43
1:B:37:TRP:HA	1:B:103:ASP:O	2.19	0.43
1:A:485:LEU:HD22	1:A:501:VAL:HB	2.01	0.43
1:B:513:ASN:O	1:B:517:GLN:HG2	2.19	0.43
1:A:243:ARG:HD2	1:A:250:MET:HE2	2.01	0.42
1:A:145:LYS:HB2	1:A:622:ARG:CZ	2.49	0.42
1:A:453:LEU:HD23	1:A:465:LEU:HD22	2.02	0.42
1:A:605:GLU:HG2	1:A:614:ARG:HD3	2.02	0.42
1:B:616:TRP:CZ2	1:B:620:LEU:HD11	2.54	0.42
1:B:147:ASN:ND2	1:B:260:SER:O	2.53	0.41
1:A:603:MET:HA	1:A:603:MET:HE2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:PHE:HB2	1:B:162:ALA:HB2	2.02	0.41
1:B:168:THR:CG2	1:B:202:SER:HB2	2.50	0.41
1:A:558:TYR:O	1:A:614:ARG:NH2	2.53	0.41
1:A:175:LYS:NZ	1:A:192:SER:OG	2.54	0.41
1:A:611:LYS:HA	1:A:614:ARG:HG3	2.03	0.41
1:B:101:MET:HA	1:B:115:VAL:HG23	2.03	0.41
1:B:408:ALA:HB1	1:B:449:ALA:HB1	2.04	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	638/649 (98%)	622 (98%)	15 (2%)	1 (0%)	43	66
1	B	639/649 (98%)	623 (98%)	16 (2%)	0	100	100
All	All	1277/1298 (98%)	1245 (98%)	31 (2%)	1 (0%)	48	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	488	GLU

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	556/565 (98%)	547 (98%)	9 (2%)	55	79
1	B	557/565 (99%)	541 (97%)	16 (3%)	37	65
All	All	1113/1130 (98%)	1088 (98%)	25 (2%)	45	72

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	THR
1	A	128	GLU
1	A	168	THR
1	A	374	CYS
1	A	386	ILE
1	A	405	ILE
1	A	577	ILE
1	A	607	CYS
1	A	642	LYS
1	B	16	GLN
1	B	56[A]	GLN
1	B	56[B]	GLN
1	B	168	THR
1	B	179	LEU
1	B	374	CYS
1	B	389	GLU
1	B	396	SER
1	B	398	GLN
1	B	405	ILE
1	B	431	TYR
1	B	478	ARG
1	B	494	ASN
1	B	540	GLU
1	B	570	ILE
1	B	577	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	339	GLN
1	A	399	HIS
1	A	468	HIS
1	A	588	ASN
1	A	613	HIS

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Mol	Chain	Res	Type
1	B	19	ASN
1	B	111	GLN
1	B	359	GLN
1	B	441	ASN
1	B	527	GLN
1	B	535	GLN
1	B	578	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	PG4	A	701	-	12,12,12	0.52	0	11,11,11	0.29	0
3	PEG	B	701	-	6,6,6	0.60	0	5,5,5	0.45	0
3	PEG	A	702	-	6,6,6	0.52	0	5,5,5	0.28	0
2	PG4	B	702	-	12,12,12	0.50	0	11,11,11	0.27	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
2	PG4	A	701	-	-	5/10/10/10	-
3	PEG	B	701	-	-	2/4/4/4	-
3	PEG	A	702	-	-	3/4/4/4	-
2	PG4	B	702	-	-	5/10/10/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	702	PEG	O1-C1-C2-O2
2	A	701	PG4	O2-C3-C4-O3
3	B	701	PEG	O2-C3-C4-O4
2	B	702	PG4	O1-C1-C2-O2
3	B	701	PEG	O1-C1-C2-O2
2	A	701	PG4	O4-C7-C8-O5
2	A	701	PG4	O1-C1-C2-O2
2	A	701	PG4	C8-C7-O4-C6
3	A	702	PEG	C1-C2-O2-C3
2	B	702	PG4	C5-C6-O4-C7
2	B	702	PG4	O2-C3-C4-O3
2	A	701	PG4	C3-C4-O3-C5
3	A	702	PEG	C4-C3-O2-C2
2	B	702	PG4	C6-C5-O3-C4
2	B	702	PG4	O3-C5-C6-O4

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	701	PEG	1	0

5.7 Other polymers [i](#)

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 [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	639/649 (98%)	0.02	5 (0%)	82 80	44, 62, 86, 127	1 (0%)
1	B	639/649 (98%)	0.01	4 (0%)	85 83	35, 63, 88, 122	2 (0%)
All	All	1278/1298 (98%)	0.01	9 (0%)	84 82	35, 63, 87, 127	3 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	71	ILE	2.7
1	A	72	VAL	2.3
1	B	72	VAL	2.3
1	A	246	ALA	2.3
1	A	4	SER	2.2
1	B	642	LYS	2.2
1	A	247	THR	2.2
1	A	642	LYS	2.1
1	B	248	LYS	2.1

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
3	PEG	A	702	7/7	0.77	0.22	95,97,98,98	0
3	PEG	B	701	7/7	0.79	0.18	72,73,74,74	0
2	PG4	B	702	13/13	0.84	0.21	95,98,105,105	0
2	PG4	A	701	13/13	0.85	0.20	92,97,101,101	0

6.5 Other polymers [i](#)

There are no such residues in this entry.