



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:02 AM UTC

PDB ID : 6QYH / pdb_00006qyh
Title : Structure of Apo HPAB from E.coli
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Deposited on : 2019-03-09
Resolution : 1.94 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

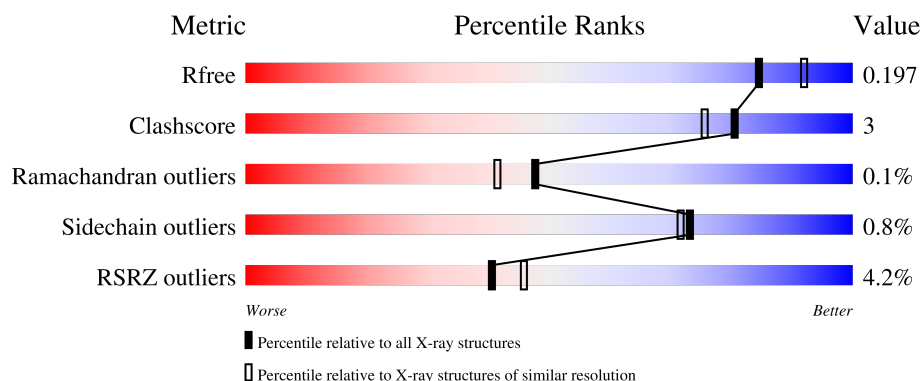
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	520	4% 89% 9%
1	B	520	4% 89% 9%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8811 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.

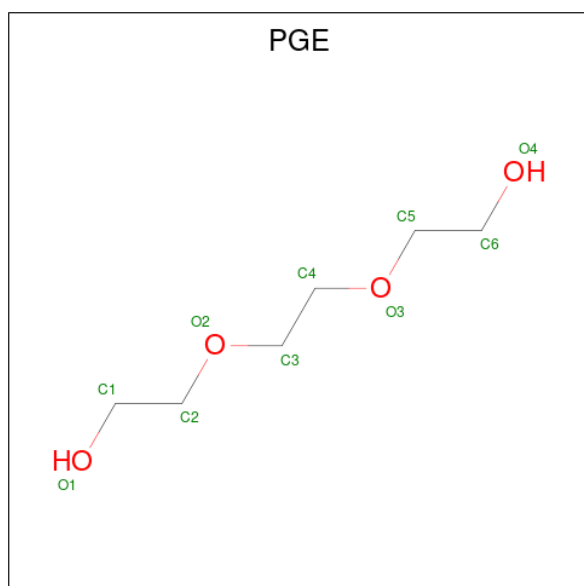
- Molecule 1 is a protein called 4-hydroxyphenylacetate 3-monooxygenase oxygenase component.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	511	Total	C	N	O	S	0	8	0
			4095	2591	709	766	29			
1	B	512	Total	C	N	O	S	0	8	0
			4107	2599	714	765	29			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

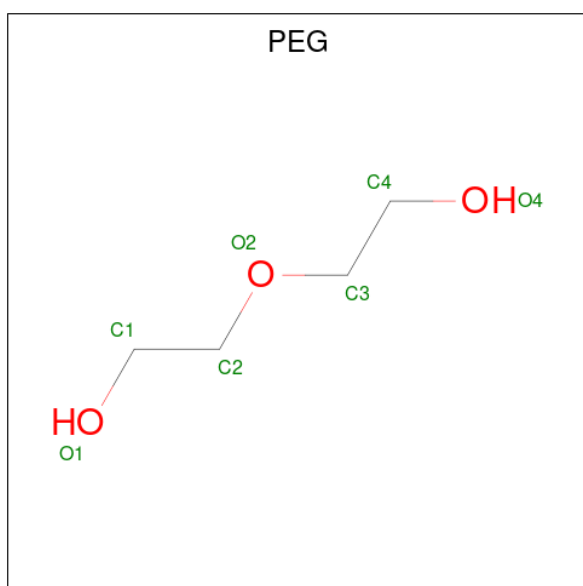
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	2	Total	Ca	0	0
			2	2		

- Molecule 3 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



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

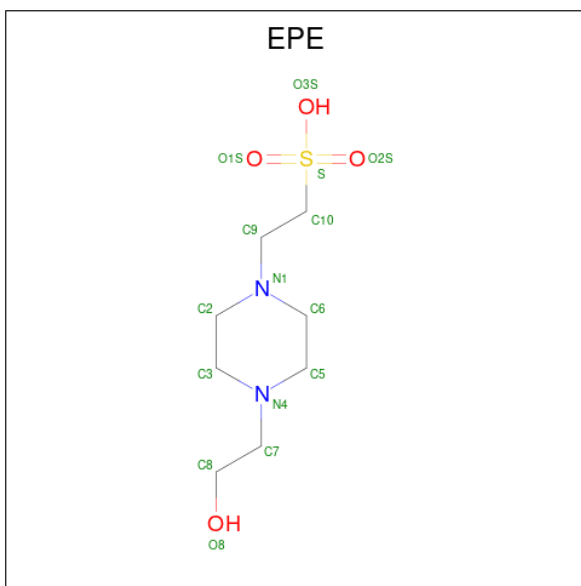
- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD

ID: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

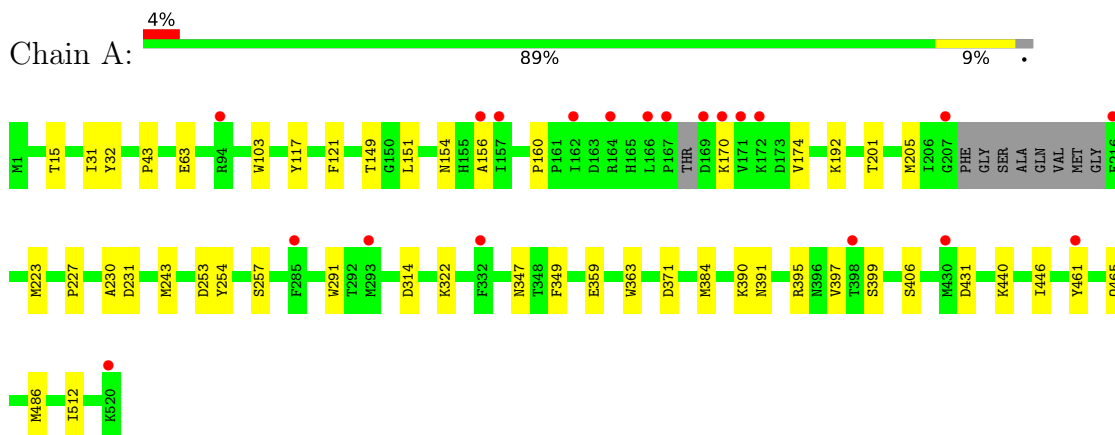
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	271	Total	O	0	0
			271	271		
6	B	231	Total	O	0	0
			231	231		

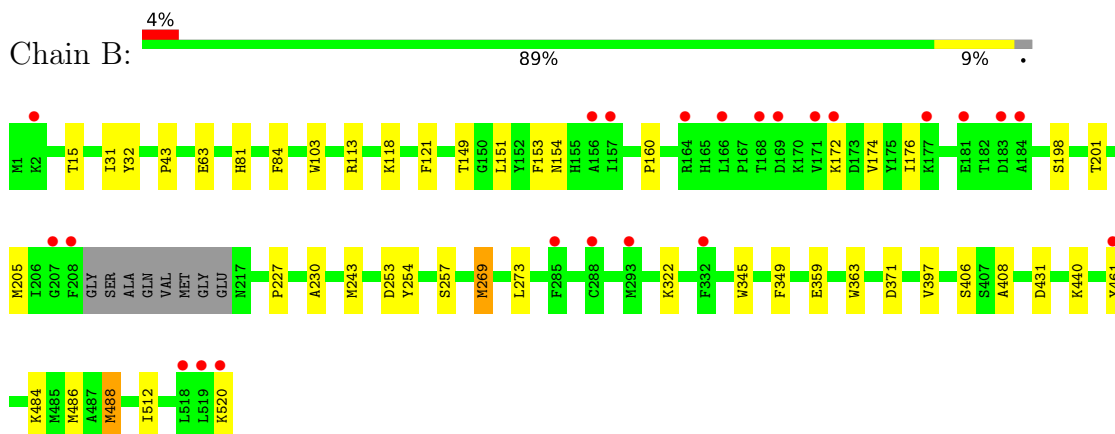
3 Residue-property plots

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: 4-hydroxyphenylacetate 3-monooxygenase oxygenase component



- Molecule 1: 4-hydroxyphenylacetate 3-monooxygenase oxygenase component



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	189.28Å 189.28Å 163.53Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.45 – 1.94 19.45 – 1.94	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.45-1.94) 99.9 (19.45-1.94)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 1.94Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.166 , 0.184 0.174 , 0.197	Depositor DCC
R_{free} test set	2220 reflections (1.74%)	wwPDB-VP
Wilson B-factor (Å ²)	39.6	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 60.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8811	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% 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

5.1 Standard geometry

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

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	0.87	1/4196 (0.0%)	1.14	12/5689 (0.2%)
1	B	0.86	3/4204 (0.1%)	1.15	7/5702 (0.1%)
All	All	0.87	4/8400 (0.0%)	1.15	19/11391 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	269	MET	SD-CE	-5.61	1.65	1.79
1	A	363	TRP	CA-C	5.61	1.55	1.52
1	B	363	TRP	CA-C	5.48	1.55	1.52
1	B	488	MET	SD-CE	-5.42	1.66	1.79

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	15	THR	N-CA-C	-6.47	101.21	110.59
1	B	15	THR	N-CA-C	-6.34	101.39	110.59
1	B	349	PHE	CA-CB-CG	-6.30	107.50	113.80
1	A	349	PHE	CA-CB-CG	-6.01	107.79	113.80
1	B	431	ASP	CA-CB-CG	5.96	118.56	112.60
1	A	431	ASP	CA-CB-CG	5.79	118.39	112.60
1	A	397	VAL	CA-C-N	5.78	128.61	120.28
1	A	397	VAL	C-N-CA	5.78	128.61	120.28
1	A	170	LYS	CA-C-N	5.49	127.47	120.56
1	A	170	LYS	C-N-CA	5.49	127.47	120.56
1	B	397	VAL	CA-C-N	5.29	127.90	120.28
1	B	397	VAL	C-N-CA	5.29	127.90	120.28
1	A	231	ASP	CA-CB-CG	5.29	117.89	112.60
1	B	406	SER	N-CA-C	5.22	117.65	111.33
1	A	399	SER	CA-C-N	5.15	126.61	120.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	399	SER	C-N-CA	5.15	126.61	120.13
1	A	347	ASN	CA-CB-CG	5.13	117.73	112.60
1	B	408	ALA	N-CA-C	-5.08	106.18	112.38
1	A	406	SER	N-CA-C	5.01	117.39	111.33

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	4095	0	3920	25	0
1	B	4107	0	3929	24	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
3	A	30	0	42	2	0
3	B	10	0	14	0	0
4	A	42	0	60	0	0
4	B	7	0	10	0	0
5	B	15	0	17	0	0
6	A	271	0	0	1	0
6	B	231	0	0	0	0
All	All	8811	0	7992	43	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 3.

All (43) 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:B:176:ILE:HD12	1:B:269:MET:HE3	1.61	0.81
1:A:154:ASN:HB2	1:A:201:THR:HG21	1.78	0.65
1:B:154:ASN:HD22	1:B:201:THR:HG21	1.68	0.59
1:A:149:THR:OG1	1:A:151:LEU:HG	2.04	0.57
1:A:391[B]:ASN:ND2	1:A:395:ARG:HH11	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:THR:OG1	1:B:151:LEU:HG	2.05	0.56
1:A:390:LYS:HE2	1:A:446:ILE:O	2.06	0.55
1:A:486:MET:SD	1:B:322:LYS:HG3	2.49	0.52
1:B:154:ASN:HB2	1:B:201:THR:HG21	1.91	0.52
1:B:160:PRO:HG2	1:B:174:VAL:HG12	1.92	0.51
1:B:113[B]:ARG:HH12	1:B:461[B]:TYR:HE1	1.57	0.51
1:A:227:PRO:HG2	1:A:230:ALA:HB2	1.93	0.51
1:A:160:PRO:HG2	1:A:174:VAL:HG12	1.93	0.51
3:A:602:PGE:H52	1:B:484:LYS:HB3	1.93	0.51
1:A:322:LYS:HG3	1:B:486:MET:SD	2.50	0.50
1:B:227:PRO:HG2	1:B:230:ALA:HB2	1.93	0.49
1:B:176:ILE:CD1	1:B:269:MET:HE3	2.36	0.49
1:A:223:MET:HE1	1:A:291:TRP:CD2	2.48	0.49
1:B:253:ASP:HA	1:B:440:LYS:HD3	1.95	0.48
1:A:359:GLU:HB2	1:A:371:ASP:HB2	1.95	0.47
1:B:81:HIS:HB3	1:B:84:PHE:CD2	2.49	0.47
1:B:118[B]:LYS:HG2	1:B:153:PHE:HB3	1.97	0.47
1:B:359:GLU:HB2	1:B:371:ASP:HB2	1.97	0.46
1:A:253:ASP:HA	1:A:440:LYS:HD3	1.97	0.46
1:A:465:GLN:HG3	3:A:603:PGE:H32	1.97	0.46
1:A:223:MET:HE1	1:A:291:TRP:CE2	2.51	0.46
1:A:117:TYR:HE1	1:A:461:TYR:HH	1.63	0.45
1:A:384[B]:MET:HE1	1:B:345:TRP:HZ2	1.82	0.45
1:A:512:ILE:HG21	1:B:43:PRO:HB3	1.99	0.45
1:B:254:TYR:HB3	1:B:257:SER:HB2	1.99	0.45
1:A:43:PRO:HB2	1:A:243:MET:HG3	1.99	0.45
1:A:117:TYR:HE1	1:A:461:TYR:OH	1.99	0.45
1:A:254:TYR:HB3	1:A:257:SER:HB2	1.98	0.44
1:B:43:PRO:HB2	1:B:243:MET:HG3	2.00	0.43
1:B:118[A]:LYS:HG3	1:B:154:ASN:HA	2.00	0.43
1:A:384[A]:MET:HG3	6:A:734:HOH:O	2.19	0.43
1:A:43:PRO:HB3	1:B:512:ILE:HG21	2.00	0.42
1:A:31:ILE:HG12	1:A:32:TYR:CD2	2.55	0.42
1:A:121:PHE:CZ	1:A:205:MET:HG3	2.54	0.41
1:A:314:ASP:HB2	1:B:488:MET:HE1	2.02	0.41
1:B:121:PHE:CZ	1:B:205:MET:HG3	2.55	0.41
1:A:156:ALA:HB1	1:A:192:LYS:HB2	2.03	0.41
1:B:31:ILE:HG12	1:B:32:TYR:CD2	2.56	0.41

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	513/520 (99%)	500 (98%)	13 (2%)	0	100	100
1	B	516/520 (99%)	503 (98%)	12 (2%)	1 (0%)	43	37
All	All	1029/1040 (99%)	1003 (98%)	25 (2%)	1 (0%)	48	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	172	LYS

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	423/436 (97%)	421 (100%)	2 (0%)	81	82
1	B	423/436 (97%)	418 (99%)	5 (1%)	63	57
All	All	846/872 (97%)	839 (99%)	7 (1%)	73	71

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	GLU
1	A	103	TRP
1	B	63	GLU
1	B	103	TRP
1	B	198	SER

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Mol	Chain	Res	Type
1	B	273	LEU
1	B	520	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	128	ASN
1	A	376	GLN
1	A	477	GLN
1	A	498	GLN
1	B	71	ASN
1	B	204	ASN
1	B	376	GLN
1	B	477	GLN
1	B	498	GLN

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](#)

Of 15 ligands modelled in this entry, 3 are monoatomic - leaving 12 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
4	PEG	A	607	-	6,6,6	0.10	0	5,5,5	0.04	0
3	PGE	A	603	-	9,9,9	0.12	0	8,8,8	0.19	0
3	PGE	A	604	-	9,9,9	0.18	0	8,8,8	0.18	0
3	PGE	A	602	-	9,9,9	0.19	0	8,8,8	0.25	0
4	PEG	A	608	-	6,6,6	0.11	0	5,5,5	0.07	0
4	PEG	A	610	-	6,6,6	0.12	0	5,5,5	0.07	0
4	PEG	A	609	-	6,6,6	0.12	0	5,5,5	0.05	0
4	PEG	B	604	-	6,6,6	0.22	0	5,5,5	0.21	0
4	PEG	A	605	-	6,6,6	0.06	0	5,5,5	0.06	0
4	PEG	A	606	-	6,6,6	0.18	0	5,5,5	0.15	0
3	PGE	B	603	-	9,9,9	0.18	0	8,8,8	0.18	0
5	EPE	B	605	-	15,15,15	1.29	1 (6%)	19,20,20	0.49	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	PEG	A	607	-	-	1/4/4/4	-
3	PGE	A	603	-	-	5/7/7/7	-
3	PGE	A	604	-	-	4/7/7/7	-
3	PGE	A	602	-	-	3/7/7/7	-
4	PEG	A	608	-	-	4/4/4/4	-
4	PEG	A	610	-	-	0/4/4/4	-
4	PEG	A	609	-	-	3/4/4/4	-
4	PEG	B	604	-	-	3/4/4/4	-
4	PEG	A	605	-	-	2/4/4/4	-
4	PEG	A	606	-	-	0/4/4/4	-
3	PGE	B	603	-	-	1/7/7/7	-
5	EPE	B	605	-	-	4/9/19/19	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	605	EPE	C10-S	-4.93	1.70	1.77

There are no bond angle outliers.

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	605	EPE	C9-C10-S-O2S
5	B	605	EPE	C9-C10-S-O3S
3	A	603	PGE	O1-C1-C2-O2
3	A	602	PGE	O2-C3-C4-O3
5	B	605	EPE	S-C10-C9-N1
4	B	604	PEG	O2-C3-C4-O4
3	A	602	PGE	O3-C5-C6-O4
5	B	605	EPE	C9-C10-S-O1S
4	A	608	PEG	C1-C2-O2-C3
4	A	609	PEG	C1-C2-O2-C3
4	B	604	PEG	C4-C3-O2-C2
3	B	603	PGE	C1-C2-O2-C3
4	A	609	PEG	C4-C3-O2-C2
4	A	607	PEG	C4-C3-O2-C2
4	A	608	PEG	C4-C3-O2-C2
4	A	605	PEG	C4-C3-O2-C2
3	A	603	PGE	C6-C5-O3-C4
3	A	604	PGE	O2-C3-C4-O3
4	A	605	PEG	O1-C1-C2-O2
3	A	604	PGE	C1-C2-O2-C3
4	A	609	PEG	O2-C3-C4-O4
3	A	603	PGE	C1-C2-O2-C3
3	A	604	PGE	C6-C5-O3-C4
3	A	602	PGE	C3-C4-O3-C5
4	A	608	PEG	O2-C3-C4-O4
3	A	603	PGE	C4-C3-O2-C2
4	B	604	PEG	C1-C2-O2-C3
3	A	603	PGE	O2-C3-C4-O3
3	A	604	PGE	O1-C1-C2-O2
4	A	608	PEG	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	603	PGE	1	0
3	A	602	PGE	1	0

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

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	511/520 (98%)	-0.00	20 (3%)	43 49	18, 42, 67, 99	8 (1%)
1	B	512/520 (98%)	0.07	23 (4%)	38 43	17, 45, 78, 111	8 (1%)
All	All	1023/1040 (98%)	0.04	43 (4%)	40 46	17, 44, 74, 111	16 (1%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	208	PHE	6.6
1	B	171	VAL	5.6
1	B	168	THR	4.9
1	A	167	PRO	4.3
1	A	207	GLY	4.2
1	B	172	LYS	4.2
1	B	520	LYS	4.1
1	A	157	ILE	4.0
1	A	169	ASP	3.8
1	B	285	PHE	3.5
1	B	164	ARG	3.4
1	A	461	TYR	3.3
1	B	169	ASP	3.3
1	B	157	ILE	3.1
1	A	162	ILE	3.0
1	B	184	ALA	2.9
1	A	171	VAL	2.8
1	A	398	THR	2.8
1	A	170	LYS	2.8
1	A	520	LYS	2.8
1	A	285	PHE	2.7
1	A	216	GLU	2.7
1	B	166	LEU	2.6
1	A	172	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	177	LYS	2.5
1	A	293	MET	2.4
1	A	332	PHE	2.4
1	B	332	PHE	2.4
1	B	293	MET	2.4
1	B	461[A]	TYR	2.3
1	A	430	MET	2.3
1	B	519	LEU	2.3
1	B	156	ALA	2.3
1	A	94[A]	ARG	2.2
1	B	288	CYS	2.2
1	B	207	GLY	2.2
1	A	156	ALA	2.2
1	B	181	GLU	2.1
1	B	518	LEU	2.1
1	B	2	LYS	2.1
1	A	166	LEU	2.0
1	B	183	ASP	2.0
1	A	164	ARG	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 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
5	EPE	B	605	15/15	0.79	0.25	76,77,81,82	15
4	PEG	B	604	7/7	0.83	0.15	74,75,77,77	0
4	PEG	A	608	7/7	0.84	0.20	84,85,87,88	0
3	PGE	A	602	10/10	0.85	0.19	70,78,80,80	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	PEG	A	609	7/7	0.85	0.15	69,70,73,75	0
3	PGE	A	603	10/10	0.86	0.18	63,71,82,83	0
4	PEG	A	606	7/7	0.86	0.16	67,68,70,72	0
3	PGE	B	603	10/10	0.87	0.15	62,64,64,65	0
4	PEG	A	607	7/7	0.87	0.14	76,76,80,80	0
4	PEG	A	605	7/7	0.88	0.14	74,74,75,77	0
3	PGE	A	604	10/10	0.89	0.17	71,76,83,85	0
4	PEG	A	610	7/7	0.89	0.14	69,70,75,75	0
2	CA	B	601	1/1	0.94	0.06	45,45,45,45	1
2	CA	B	602	1/1	0.96	0.04	45,45,45,45	1
2	CA	A	601	1/1	0.98	0.12	44,44,44,44	1

6.5 Other polymers [i](#)

There are no such residues in this entry.