



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:06 PM UTC

PDB ID : 4QTO / pdb_00004qto
Title : 1.65 Angstrom resolution crystal structure of betaine aldehyde dehydrogenase (betB) from Staphylococcus aureus with BME-modified Cys289 and PEG molecule in active site
Authors : Halavaty, A.S.; Minasov, G.; Dubrovskaya, I.; Winsor, J.; Shuvalova, L.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2014-07-08
Resolution : 1.65 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

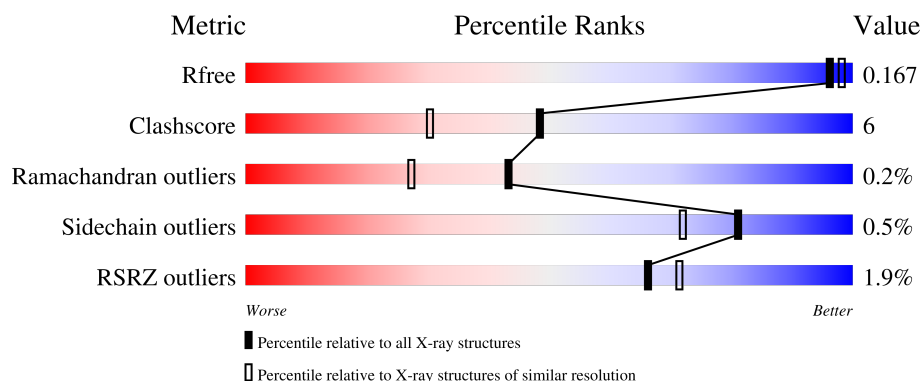
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	2% 86% 9% .
1	B	520	3% 86% 9% .
1	C	520	2% 87% 8% 5% .
1	D	520	0% 87% 8% 5% .

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 19380 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 Betaine aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	497	Total	C	N	O	S	0	34	0
			4145	2611	699	815	20			
1	B	497	Total	C	N	O	S	0	33	0
			4124	2596	695	814	19			
1	C	496	Total	C	N	O	S	0	38	0
			4161	2620	707	814	20			
1	D	495	Total	C	N	O	S	0	28	0
			4079	2568	690	803	18			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	expression tag	UNP Q5HCU0
A	-22	HIS	-	expression tag	UNP Q5HCU0
A	-21	HIS	-	expression tag	UNP Q5HCU0
A	-20	HIS	-	expression tag	UNP Q5HCU0
A	-19	HIS	-	expression tag	UNP Q5HCU0
A	-18	HIS	-	expression tag	UNP Q5HCU0
A	-17	HIS	-	expression tag	UNP Q5HCU0
A	-16	SER	-	expression tag	UNP Q5HCU0
A	-15	SER	-	expression tag	UNP Q5HCU0
A	-14	GLY	-	expression tag	UNP Q5HCU0
A	-13	VAL	-	expression tag	UNP Q5HCU0
A	-12	ASP	-	expression tag	UNP Q5HCU0
A	-11	LEU	-	expression tag	UNP Q5HCU0
A	-10	GLY	-	expression tag	UNP Q5HCU0
A	-9	THR	-	expression tag	UNP Q5HCU0
A	-8	GLU	-	expression tag	UNP Q5HCU0
A	-7	ASN	-	expression tag	UNP Q5HCU0
A	-6	LEU	-	expression tag	UNP Q5HCU0
A	-5	TYR	-	expression tag	UNP Q5HCU0
A	-4	PHE	-	expression tag	UNP Q5HCU0
A	-3	GLN	-	expression tag	UNP Q5HCU0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q5HCU0
A	-1	ASN	-	expression tag	UNP Q5HCU0
A	0	ALA	-	expression tag	UNP Q5HCU0
B	-23	MET	-	expression tag	UNP Q5HCU0
B	-22	HIS	-	expression tag	UNP Q5HCU0
B	-21	HIS	-	expression tag	UNP Q5HCU0
B	-20	HIS	-	expression tag	UNP Q5HCU0
B	-19	HIS	-	expression tag	UNP Q5HCU0
B	-18	HIS	-	expression tag	UNP Q5HCU0
B	-17	HIS	-	expression tag	UNP Q5HCU0
B	-16	SER	-	expression tag	UNP Q5HCU0
B	-15	SER	-	expression tag	UNP Q5HCU0
B	-14	GLY	-	expression tag	UNP Q5HCU0
B	-13	VAL	-	expression tag	UNP Q5HCU0
B	-12	ASP	-	expression tag	UNP Q5HCU0
B	-11	LEU	-	expression tag	UNP Q5HCU0
B	-10	GLY	-	expression tag	UNP Q5HCU0
B	-9	THR	-	expression tag	UNP Q5HCU0
B	-8	GLU	-	expression tag	UNP Q5HCU0
B	-7	ASN	-	expression tag	UNP Q5HCU0
B	-6	LEU	-	expression tag	UNP Q5HCU0
B	-5	TYR	-	expression tag	UNP Q5HCU0
B	-4	PHE	-	expression tag	UNP Q5HCU0
B	-3	GLN	-	expression tag	UNP Q5HCU0
B	-2	SER	-	expression tag	UNP Q5HCU0
B	-1	ASN	-	expression tag	UNP Q5HCU0
B	0	ALA	-	expression tag	UNP Q5HCU0
C	-23	MET	-	expression tag	UNP Q5HCU0
C	-22	HIS	-	expression tag	UNP Q5HCU0
C	-21	HIS	-	expression tag	UNP Q5HCU0
C	-20	HIS	-	expression tag	UNP Q5HCU0
C	-19	HIS	-	expression tag	UNP Q5HCU0
C	-18	HIS	-	expression tag	UNP Q5HCU0
C	-17	HIS	-	expression tag	UNP Q5HCU0
C	-16	SER	-	expression tag	UNP Q5HCU0
C	-15	SER	-	expression tag	UNP Q5HCU0
C	-14	GLY	-	expression tag	UNP Q5HCU0
C	-13	VAL	-	expression tag	UNP Q5HCU0
C	-12	ASP	-	expression tag	UNP Q5HCU0
C	-11	LEU	-	expression tag	UNP Q5HCU0
C	-10	GLY	-	expression tag	UNP Q5HCU0
C	-9	THR	-	expression tag	UNP Q5HCU0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	GLU	-	expression tag	UNP Q5HCU0
C	-7	ASN	-	expression tag	UNP Q5HCU0
C	-6	LEU	-	expression tag	UNP Q5HCU0
C	-5	TYR	-	expression tag	UNP Q5HCU0
C	-4	PHE	-	expression tag	UNP Q5HCU0
C	-3	GLN	-	expression tag	UNP Q5HCU0
C	-2	SER	-	expression tag	UNP Q5HCU0
C	-1	ASN	-	expression tag	UNP Q5HCU0
C	0	ALA	-	expression tag	UNP Q5HCU0
D	-23	MET	-	expression tag	UNP Q5HCU0
D	-22	HIS	-	expression tag	UNP Q5HCU0
D	-21	HIS	-	expression tag	UNP Q5HCU0
D	-20	HIS	-	expression tag	UNP Q5HCU0
D	-19	HIS	-	expression tag	UNP Q5HCU0
D	-18	HIS	-	expression tag	UNP Q5HCU0
D	-17	HIS	-	expression tag	UNP Q5HCU0
D	-16	SER	-	expression tag	UNP Q5HCU0
D	-15	SER	-	expression tag	UNP Q5HCU0
D	-14	GLY	-	expression tag	UNP Q5HCU0
D	-13	VAL	-	expression tag	UNP Q5HCU0
D	-12	ASP	-	expression tag	UNP Q5HCU0
D	-11	LEU	-	expression tag	UNP Q5HCU0
D	-10	GLY	-	expression tag	UNP Q5HCU0
D	-9	THR	-	expression tag	UNP Q5HCU0
D	-8	GLU	-	expression tag	UNP Q5HCU0
D	-7	ASN	-	expression tag	UNP Q5HCU0
D	-6	LEU	-	expression tag	UNP Q5HCU0
D	-5	TYR	-	expression tag	UNP Q5HCU0
D	-4	PHE	-	expression tag	UNP Q5HCU0
D	-3	GLN	-	expression tag	UNP Q5HCU0
D	-2	SER	-	expression tag	UNP Q5HCU0
D	-1	ASN	-	expression tag	UNP Q5HCU0
D	0	ALA	-	expression tag	UNP Q5HCU0

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Na	0	0
			1	1		

- 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		
3	A	1	Total	C	O	0	0
			10	6	4		
3	B	1	Total	C	O	0	0
			10	6	4		
3	C	1	Total	C	O	0	0
			10	6	4		
3	D	1	Total	C	O	0	0
			10	6	4		

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



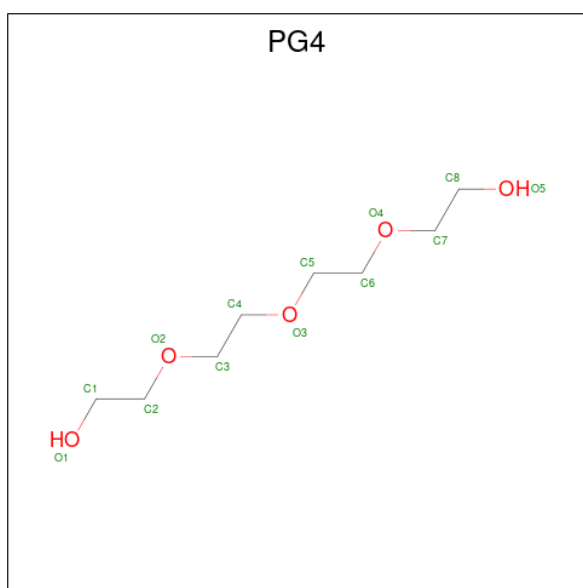
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	A	1	Total	C	O	0	0
			7	4	3		
4	A	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			8	4	1	3		
5	B	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			13	8	5		

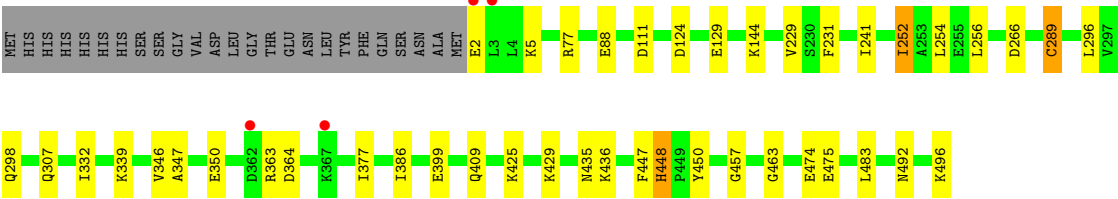
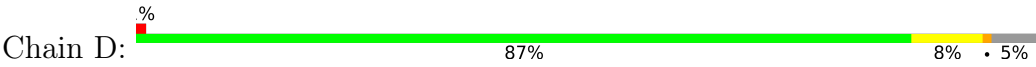
- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	677	Total	O	0	36
			700	700		
8	B	646	Total	O	0	29
			662	662		
8	C	664	Total	O	0	31
			685	685		
8	D	676	Total	O	0	25
			695	695		



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	224.25Å 102.47Å 118.18Å 90.00° 104.45° 90.00°	Depositor
Resolution (Å)	29.71 – 1.65 29.71 – 1.65	Depositor EDS
% Data completeness (in resolution range)	98.6 (29.71-1.65) 98.6 (29.71-1.65)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.136 , 0.156 0.149 , 0.167	Depositor DCC
R_{free} test set	15433 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	15.0	Xtriage
Anisotropy	0.228	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19380	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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, CME, PG4, PGE, TRS, NA, 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	0.70	0/4196	0.92	9/5667 (0.2%)
1	B	0.69	1/4177 (0.0%)	0.92	10/5645 (0.2%)
1	C	0.69	0/4212	0.91	8/5688 (0.1%)
1	D	0.68	0/4132	0.91	7/5585 (0.1%)
All	All	0.69	1/16717 (0.0%)	0.91	34/22585 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	258	GLY	C-O	-5.14	1.18	1.23

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	457	GLY	CA-C-N	-8.33	116.43	122.18
1	B	457	GLY	C-N-CA	-8.33	116.43	122.18
1	C	457	GLY	CA-C-N	-8.24	116.50	122.18
1	C	457	GLY	C-N-CA	-8.24	116.50	122.18
1	A	457	GLY	CA-C-N	-7.19	117.22	122.18
1	A	457	GLY	C-N-CA	-7.19	117.22	122.18
1	D	457	GLY	CA-C-N	-7.08	117.29	122.18
1	D	457	GLY	C-N-CA	-7.08	117.29	122.18
1	D	256	LEU	N-CA-C	6.65	121.06	112.41
1	C	256	LEU	N-CA-C	6.61	121.01	112.41
1	B	332	ILE	N-CA-C	6.43	120.06	111.44
1	D	332	ILE	N-CA-C	6.04	119.53	111.44
1	B	187	LEU	N-CA-C	6.01	117.52	110.97
1	A	332	ILE	N-CA-C	5.99	119.46	111.44
1	C	449	PRO	N-CA-C	5.69	119.79	111.03
1	A	256	LEU	N-CA-C	5.69	119.80	112.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	269	ASP	N-CA-C	-5.56	101.14	109.15
1	A	269	ASP	N-CA-C	-5.41	101.36	109.15
1	A	458	GLY	N-CA-C	5.36	120.56	111.19
1	A	10	ARG	N-CA-C	5.33	117.97	110.14
1	C	10	ARG	N-CA-C	5.30	117.94	110.14
1	D	363	ARG	N-CA-C	-5.23	100.51	109.24
1	C	332	ILE	N-CA-C	5.21	118.42	111.44
1	B	448	HIS	N-CA-C	5.20	121.30	109.81
1	C	448	HIS	N-CA-C	5.17	121.24	109.81
1	B	256	LEU	N-CA-C	5.15	119.11	112.41
1	B	449	PRO	N-CA-C	5.15	119.12	111.09
1	D	463	GLY	N-CA-C	5.14	117.98	110.63
1	B	364	ASP	N-CA-C	5.13	117.53	111.33
1	A	187	LEU	N-CA-C	5.11	116.54	111.07
1	B	108	ASP	CB-CA-C	-5.07	102.91	110.88
1	D	364	ASP	N-CA-C	5.05	117.17	111.11
1	B	10	ARG	N-CA-C	5.03	117.53	110.14
1	A	364	ASP	N-CA-C	5.01	117.12	111.11

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	4145	0	4075	43	0
1	B	4124	0	4047	55	0
1	C	4161	0	4097	48	0
1	D	4079	0	3991	41	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	20	0	28	0	0
3	B	10	0	14	0	0
3	C	10	0	14	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	10	0	14	1	0
4	A	21	0	30	2	0
4	B	14	0	20	1	0
4	D	7	0	10	0	0
5	A	8	0	12	0	0
5	B	8	0	12	0	0
6	C	13	0	18	0	0
7	C	4	0	6	3	0
8	A	700	0	0	25	0
8	B	662	0	0	24	0
8	C	685	0	0	28	0
8	D	695	0	0	18	0
All	All	19380	0	16388	183	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 6.

All (183) 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:450[B]:TYR:HE2	8:B:951:HOH:O	1.32	1.10
8:C:1237:HOH:O	1:D:474:GLU:HG3	1.51	1.09
1:D:450[A]:TYR:HE2	8:D:1252:HOH:O	1.38	1.07
1:A:450[B]:TYR:HE1	8:A:729:HOH:O	1.39	1.05
1:B:77[B]:ARG:NH1	1:C:117[B]:MET:CE	2.25	1.00
1:C:29[B]:ILE:HD11	8:C:890:HOH:O	1.65	0.96
1:A:212:ALA:HB3	1:A:215[B]:GLU:HG2	1.52	0.91
1:B:77[B]:ARG:NH1	1:C:117[B]:MET:HE2	1.87	0.88
1:A:211:GLY:HA3	8:A:1099[B]:HOH:O	1.76	0.84
1:C:475:GLU:HG3	8:C:914:HOH:O	1.75	0.84
1:D:229[B]:VAL:HG13	1:D:252[B]:ILE:HD12	1.59	0.83
1:A:126[A]:ASP:HB3	8:A:1277[A]:HOH:O	1.79	0.82
1:C:211:GLY:HA3	8:C:880:HOH:O	1.80	0.81
1:A:311:ASP:HB3	8:A:1213:HOH:O	1.80	0.80
1:B:126[A]:ASP:HB3	8:B:1246[A]:HOH:O	1.82	0.79
1:C:129:GLU:HG3	8:C:1220:HOH:O	1.82	0.78
1:B:77[B]:ARG:NH1	1:C:117[B]:MET:HE1	1.99	0.77
1:D:2:GLU:O	1:D:5:LYS:HG2	1.85	0.76
1:A:212:ALA:HB3	1:A:215[B]:GLU:CG	2.15	0.76
1:B:2:GLU:O	1:B:5:LYS:HG2	1.84	0.76
1:B:307[A]:GLN:HG2	8:B:959:HOH:O	1.85	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:475:GLU:HG3	8:A:844:HOH:O	1.86	0.75
1:A:2:GLU:O	1:A:5:LYS:HG2	1.86	0.75
1:D:347:ALA:HB2	1:D:386[B]:ILE:HG21	1.71	0.72
1:D:77[B]:ARG:NH2	8:D:849:HOH:O	2.22	0.72
1:A:431[B]:GLN:NE2	8:A:1210:HOH:O	2.24	0.71
1:D:129:GLU:HG3	8:D:1269:HOH:O	1.90	0.71
1:B:307[B]:GLN:NE2	1:B:311[B]:ASP:OD1	2.24	0.70
1:C:147[B]:VAL:HG13	1:C:476:TYR:C	2.16	0.70
1:B:351:GLY:C	8:B:954:HOH:O	2.32	0.70
1:A:77[A]:ARG:NH2	8:A:902:HOH:O	2.22	0.70
1:A:347:ALA:HA	1:A:386[B]:ILE:HG21	1.74	0.69
1:D:307:GLN:HG2	8:D:893:HOH:O	1.91	0.69
1:B:475:GLU:HG3	8:B:896:HOH:O	1.93	0.69
1:A:347:ALA:CA	1:A:386[B]:ILE:HG21	2.24	0.68
1:A:84:LYS:NZ	8:A:891:HOH:O	2.26	0.68
1:B:311[B]:ASP:HB3	8:B:920[B]:HOH:O	1.93	0.68
1:C:244[A]:ASN:ND2	8:C:941:HOH:O	2.25	0.68
1:B:409[A]:GLN:HG2	8:B:888:HOH:O	1.93	0.68
1:D:475:GLU:HG3	8:D:1133:HOH:O	1.94	0.67
1:C:229[B]:VAL:HG13	1:C:252[B]:ILE:HD12	1.78	0.66
1:C:347:ALA:HB2	1:C:386[B]:ILE:HG21	1.78	0.66
1:B:77[B]:ARG:CZ	1:C:117[B]:MET:CE	2.74	0.66
1:C:93[B]:LEU:HD12	8:C:888:HOH:O	1.95	0.65
1:C:3:LEU:HB3	8:C:958:HOH:O	1.97	0.65
1:B:77[B]:ARG:CZ	1:C:117[B]:MET:HE1	2.26	0.65
1:D:347:ALA:CB	1:D:386[B]:ILE:HG21	2.27	0.63
1:C:289[A]:CME:HZ3	8:C:1261[A]:HOH:O	1.97	0.63
4:A:504:PEG:H22	8:A:1235:HOH:O	1.97	0.63
8:A:992:HOH:O	1:C:124:ASP:HB2	1.98	0.63
1:D:229[B]:VAL:HG13	1:D:252[B]:ILE:CD1	2.29	0.62
1:D:425[A]:LYS:NZ	8:D:946:HOH:O	2.23	0.62
1:A:350:GLU:OE1	1:A:386[B]:ILE:HG23	2.00	0.61
1:A:77[B]:ARG:HD3	8:A:784[B]:HOH:O	2.00	0.61
1:C:224:LYS:HE2	8:C:1116[B]:HOH:O	2.00	0.61
1:A:450[B]:TYR:C	1:A:450[B]:TYR:CD1	2.78	0.61
1:A:347:ALA:HB2	1:A:386[B]:ILE:HG21	1.84	0.60
1:A:450[B]:TYR:CE1	8:A:729:HOH:O	2.27	0.60
1:C:93[B]:LEU:CD1	8:C:888:HOH:O	2.48	0.60
1:C:1:MET:SD	1:C:3:LEU:HD12	2.42	0.60
1:B:307[B]:GLN:HE21	1:B:311[B]:ASP:CG	2.10	0.59
1:A:347:ALA:CB	1:A:386[B]:ILE:HG21	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ASP:HB2	8:C:1054:HOH:O	2.01	0.59
1:A:117[B]:MET:CE	1:D:77[B]:ARG:NH1	2.66	0.59
1:A:129:GLU:CG	1:A:142[B]:ILE:HD12	2.33	0.58
1:A:304[A]:LYS:NZ	8:A:1272:HOH:O	2.32	0.58
1:D:88:GLU:HG2	8:D:1242:HOH:O	2.02	0.58
1:B:409[A]:GLN:NE2	8:B:1012:HOH:O	2.35	0.58
3:D:502:PGE:H3	8:D:1085[B]:HOH:O	2.03	0.57
1:A:126[A]:ASP:CG	8:A:1277[A]:HOH:O	2.46	0.57
1:B:351:GLY:HA2	8:B:954:HOH:O	2.04	0.57
1:D:450[A]:TYR:CE2	8:D:1252:HOH:O	2.27	0.56
1:B:450[B]:TYR:CE2	8:B:951:HOH:O	2.21	0.56
1:D:347:ALA:CA	1:D:386[B]:ILE:HG21	2.35	0.56
8:B:1167:HOH:O	1:D:124:ASP:HB2	2.05	0.56
1:B:289[B]:CME:SG	1:B:418:LEU:HD21	2.45	0.56
1:C:224:LYS:CE	8:C:1116[B]:HOH:O	2.54	0.56
7:C:504[B]:EDO:H11	8:C:803:HOH:O	2.06	0.56
1:B:144[B]:LYS:HE2	1:B:474[B]:GLU:OE2	2.05	0.56
1:B:347:ALA:HB2	1:B:386[B]:ILE:HG21	1.87	0.55
1:D:436:LYS:NZ	8:D:1223:HOH:O	2.25	0.55
1:A:126[A]:ASP:CB	8:A:1277[A]:HOH:O	2.45	0.55
7:C:504[B]:EDO:C1	8:C:803:HOH:O	2.55	0.55
1:D:429:LYS:NZ	8:D:1244[A]:HOH:O	2.39	0.54
1:C:1:MET:HG2	1:C:3:LEU:H	1.72	0.54
1:C:449:PRO:HB2	1:D:483[A]:LEU:CD2	2.38	0.54
1:A:339:LYS:HD2	8:A:1119:HOH:O	2.08	0.53
1:A:241[B]:ILE:HD13	8:A:918:HOH:O	2.07	0.53
1:D:409[A]:GLN:HG2	8:D:876:HOH:O	2.07	0.53
1:D:409[A]:GLN:HG3	8:D:831:HOH:O	2.08	0.53
1:D:266:ASP:OD1	1:D:298[B]:GLN:OE1	2.27	0.52
1:C:347:ALA:CA	1:C:386[B]:ILE:HG21	2.40	0.52
7:C:504[B]:EDO:C2	8:C:951:HOH:O	2.58	0.52
1:A:409[A]:GLN:HG3	8:A:803:HOH:O	2.09	0.51
1:B:409[A]:GLN:HG3	8:B:738:HOH:O	2.11	0.51
1:A:244[A]:ASN:ND2	8:A:918:HOH:O	2.44	0.51
1:D:347:ALA:HA	1:D:386[B]:ILE:HG21	1.92	0.51
1:B:241[B]:ILE:HD13	8:B:1218:HOH:O	2.11	0.51
1:B:347:ALA:CA	1:B:386[B]:ILE:HG21	2.41	0.51
1:C:147[B]:VAL:HG13	1:C:477:LEU:N	2.26	0.51
1:C:289[A]:CME:CZ	8:C:1261[A]:HOH:O	2.56	0.50
1:C:409[A]:GLN:HG3	8:C:789:HOH:O	2.11	0.50
1:C:347:ALA:CB	1:C:386[B]:ILE:HG21	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:ASP:O	8:B:1167:HOH:O	2.20	0.50
1:B:351:GLY:CA	8:B:954:HOH:O	2.57	0.50
1:B:289[A]:CME:HZ2	4:B:504:PEG:C4	2.42	0.49
1:A:346:VAL:O	1:A:350:GLU:HG3	2.13	0.49
1:B:150[B]:VAL:HG12	1:B:228[B]:LEU:HB3	1.94	0.49
8:A:1207:HOH:O	1:B:431:GLN:HG2	2.12	0.49
1:A:405:GLN:HG3	8:A:945:HOH:O	2.13	0.48
1:A:296:LEU:HD23	1:A:399:GLU:HB2	1.95	0.48
1:B:307[A]:GLN:HG3	8:B:1117:HOH:O	2.12	0.48
1:D:435[B]:ASN:ND2	8:D:800:HOH:O	2.43	0.48
1:C:431[B]:GLN:NE2	1:C:435[B]:ASN:OD1	2.42	0.48
1:C:270:PHE:CE1	1:C:304[B]:LYS:HE3	2.49	0.47
1:A:77[A]:ARG:HG3	1:A:117[A]:MET:SD	2.54	0.47
1:D:346:VAL:O	1:D:350:GLU:HG3	2.14	0.47
1:C:132:ASP:O	8:C:1054:HOH:O	2.21	0.47
1:C:346:VAL:O	1:C:350:GLU:HG3	2.15	0.47
1:B:483[B]:LEU:HD23	1:B:483[B]:LEU:C	2.39	0.47
1:C:296:LEU:HD23	1:C:399:GLU:HB2	1.97	0.47
1:D:289[B]:CME:HB3	8:D:1275[B]:HOH:O	2.14	0.47
1:B:255[B]:GLU:HG3	8:B:730:HOH:O	2.15	0.46
1:B:347:ALA:HA	1:B:386[B]:ILE:HG21	1.97	0.46
1:B:347:ALA:CB	1:B:386[B]:ILE:HG21	2.44	0.46
1:D:229[B]:VAL:CG1	1:D:252[B]:ILE:CD1	2.92	0.46
1:C:3:LEU:HD21	8:C:979:HOH:O	2.14	0.46
1:D:296:LEU:HD23	1:D:399:GLU:HB2	1.97	0.46
1:B:346:VAL:O	1:B:350:GLU:HG3	2.15	0.46
1:A:144:LYS:NZ	8:A:884:HOH:O	2.49	0.46
1:C:314:LYS:NZ	8:C:942:HOH:O	2.48	0.46
1:C:289[B]:CME:SD	8:C:1240:HOH:O	2.61	0.45
1:D:229[B]:VAL:CG1	1:D:252[B]:ILE:HD12	2.38	0.45
1:C:252[B]:ILE:HD11	1:C:254:LEU:HD21	1.99	0.45
1:A:80:ALA:HB1	1:A:117[B]:MET:HG2	1.99	0.45
1:B:126[A]:ASP:CG	8:B:1246[A]:HOH:O	2.60	0.45
1:D:377:ILE:HG21	1:D:386[B]:ILE:HD12	1.99	0.45
1:D:450[A]:TYR:CD2	1:D:450[A]:TYR:C	2.95	0.45
1:B:350:GLU:OE1	1:B:386[B]:ILE:HG23	2.17	0.45
1:B:471:GLU:OE1	8:B:990:HOH:O	2.21	0.45
1:C:425[B]:LYS:CE	8:C:893:HOH:O	2.64	0.45
1:A:409[A]:GLN:HG2	8:A:901:HOH:O	2.17	0.44
1:B:126[A]:ASP:CB	8:B:1246[A]:HOH:O	2.52	0.44
1:D:350:GLU:OE1	1:D:386[B]:ILE:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:ASP:N	8:A:1006:HOH:O	2.41	0.44
1:B:296:LEU:HD23	1:B:399:GLU:HB2	1.99	0.44
1:C:229[B]:VAL:HG13	1:C:252[B]:ILE:CD1	2.45	0.44
1:B:77[B]:ARG:HG3	1:B:117:MET:SD	2.58	0.44
1:D:231:PHE:CD1	1:D:241:ILE:CD1	3.00	0.43
1:D:339:LYS:HD2	8:D:1092:HOH:O	2.18	0.43
1:A:431[B]:GLN:NE2	1:A:435[B]:ASN:OD1	2.51	0.43
1:A:270:PHE:CE1	1:A:304[B]:LYS:HE3	2.53	0.43
1:C:144[B]:LYS:HE2	8:C:1220:HOH:O	2.17	0.43
1:C:386[B]:ILE:HD12	1:C:387:VAL:HG23	2.01	0.43
1:A:117[B]:MET:HE1	1:D:77[B]:ARG:NH1	2.34	0.43
1:B:231:PHE:CD1	1:B:241[A]:ILE:CD1	3.02	0.42
1:A:449:PRO:HB2	1:B:483[B]:LEU:HD13	2.00	0.42
1:B:144[B]:LYS:HE3	1:B:477:LEU:HB3	2.01	0.42
1:C:409[A]:GLN:HG2	8:C:873:HOH:O	2.18	0.42
1:B:450[B]:TYR:CD2	1:B:450[B]:TYR:C	2.97	0.42
1:B:144[B]:LYS:NZ	1:B:474[B]:GLU:OE2	2.52	0.42
1:B:270:PHE:CE1	1:B:304[B]:LYS:HE3	2.55	0.42
1:D:144[A]:LYS:HE2	8:D:1269:HOH:O	2.20	0.42
1:B:74:LYS:HE2	8:B:795:HOH:O	2.20	0.42
1:D:447:PHE:O	1:D:448:HIS:HB2	2.20	0.42
1:B:77[B]:ARG:HH12	1:C:117[B]:MET:HE2	1.81	0.41
1:B:16:GLU:HG3	8:B:930:HOH:O	2.19	0.41
1:B:77[B]:ARG:HH11	1:C:117[B]:MET:CE	2.26	0.41
1:C:447:PHE:O	1:C:448:HIS:HB2	2.19	0.41
1:D:111[B]:ASP:OD1	8:D:987:HOH:O	2.21	0.41
1:A:252[B]:ILE:HD11	1:A:254:LEU:HD21	2.03	0.41
1:B:350:GLU:HG3	8:B:875:HOH:O	2.20	0.41
1:B:447:PHE:O	1:B:448:HIS:HB2	2.21	0.41
4:A:504:PEG:C2	8:A:1235:HOH:O	2.61	0.41
1:C:77[A]:ARG:HG3	1:C:117[A]:MET:SD	2.61	0.41
1:D:492:ASN:OD1	1:D:496:LYS:CE	2.69	0.41
1:B:307[B]:GLN:HG2	8:B:1117:HOH:O	2.20	0.41
1:D:252[B]:ILE:HD11	1:D:254:LEU:HD21	2.03	0.41
1:C:425[B]:LYS:NZ	8:C:893:HOH:O	2.48	0.40
3:C:502:PGE:H42	8:C:1208:HOH:O	2.20	0.40
1:B:144[B]:LYS:CE	1:B:474[B]:GLU:OE2	2.69	0.40
1:A:307:GLN:NE2	1:A:311:ASP:OD1	2.49	0.40
1:C:304[B]:LYS:NZ	8:C:1090:HOH:O	2.33	0.40

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	528/520 (102%)	519 (98%)	8 (2%)	1 (0%)	43	27
1	B	526/520 (101%)	517 (98%)	8 (2%)	1 (0%)	43	27
1	C	530/520 (102%)	521 (98%)	8 (2%)	1 (0%)	43	27
1	D	519/520 (100%)	509 (98%)	9 (2%)	1 (0%)	43	27
All	All	2103/2080 (101%)	2066 (98%)	33 (2%)	4 (0%)	43	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	448	HIS
1	B	448	HIS
1	C	448	HIS
1	D	448	HIS

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	443/431 (103%)	439 (99%)	4 (1%)	70	56
1	B	441/431 (102%)	436 (99%)	5 (1%)	65	48
1	C	446/431 (104%)	441 (99%)	5 (1%)	65	48
1	D	435/431 (101%)	433 (100%)	2 (0%)	81	72
All	All	1765/1724 (102%)	1749 (99%)	16 (1%)	81	56

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	252[A]	ILE
1	A	252[B]	ILE
1	A	255[A]	GLU
1	A	255[B]	GLU
1	B	129	GLU
1	B	252[A]	ILE
1	B	252[B]	ILE
1	B	255[A]	GLU
1	B	255[B]	GLU
1	C	84	LYS
1	C	252[A]	ILE
1	C	252[B]	ILE
1	C	380[A]	CYS
1	C	380[B]	CYS
1	D	252[A]	ILE
1	D	252[B]	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	248	ASN
1	A	251	ASN
1	B	34	GLN
1	B	162	GLN
1	B	248	ASN
1	B	251	ASN
1	C	34	GLN
1	C	251	ASN
1	D	34	GLN
1	D	251	ASN
1	D	260	ASN
1	D	307	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues 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
1	CME	C	289[B]	1	8,9,10	0.87	0	6,9,11	1.89	1 (16%)
1	CME	A	289[A]	1	8,9,10	0.75	0	6,9,11	1.82	1 (16%)
1	CME	D	289[A]	1	8,9,10	0.81	0	6,9,11	1.75	1 (16%)
1	CME	D	289[B]	1	8,9,10	0.85	0	6,9,11	2.66	1 (16%)
1	CME	A	289[B]	1	8,9,10	0.96	0	6,9,11	4.17	3 (50%)
1	CME	B	289[A]	1	8,9,10	0.77	0	6,9,11	1.68	1 (16%)
1	CME	B	289[B]	1	8,9,10	1.00	0	6,9,11	2.03	1 (16%)
1	CME	C	289[A]	1	8,9,10	0.92	1 (12%)	6,9,11	1.89	3 (50%)

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
1	CME	C	289[B]	1	-	3/5/8/10	-
1	CME	A	289[A]	1	-	1/5/8/10	-
1	CME	D	289[A]	1	-	3/5/8/10	-
1	CME	D	289[B]	1	-	2/5/8/10	-
1	CME	A	289[B]	1	-	0/5/8/10	-
1	CME	B	289[A]	1	-	1/5/8/10	-
1	CME	B	289[B]	1	-	1/5/8/10	-
1	CME	C	289[A]	1	-	3/5/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	289[A]	CME	CE-SD	-2.02	1.73	1.82

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	289[B]	CME	CB-SG-SD	9.44	128.28	103.86
1	D	289[B]	CME	CB-SG-SD	-6.21	87.78	103.86
1	A	289[A]	CME	CB-SG-SD	-4.23	92.92	103.86
1	B	289[B]	CME	CB-SG-SD	3.82	113.74	103.86
1	C	289[B]	CME	CB-SG-SD	3.81	113.73	103.86
1	B	289[A]	CME	CB-SG-SD	-3.73	94.20	103.86
1	D	289[A]	CME	CB-SG-SD	3.51	112.94	103.86
1	C	289[A]	CME	CA-CB-SG	-3.20	101.32	114.45
1	A	289[B]	CME	CZ-CE-SD	-2.70	104.35	113.39
1	A	289[B]	CME	CA-CB-SG	2.38	124.22	114.45
1	C	289[A]	CME	CB-CA-C	2.20	116.78	110.80
1	C	289[A]	CME	CE-SD-SG	-2.17	93.92	103.46

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	289[A]	CME	SD-CE-CZ-OH
1	C	289[B]	CME	CE-SD-SG-CB
1	D	289[A]	CME	CE-SD-SG-CB
1	C	289[A]	CME	CE-SD-SG-CB
1	A	289[A]	CME	SD-CE-CZ-OH
1	D	289[B]	CME	SD-CE-CZ-OH
1	C	289[A]	CME	CA-CB-SG-SD
1	C	289[B]	CME	SD-CE-CZ-OH
1	D	289[A]	CME	SD-CE-CZ-OH
1	C	289[A]	CME	SD-CE-CZ-OH
1	D	289[B]	CME	CE-SD-SG-CB
1	B	289[B]	CME	CZ-CE-SD-SG
1	D	289[A]	CME	CA-CB-SG-SD
1	C	289[B]	CME	CA-CB-SG-SD

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	289[B]	CME	1	0
1	D	289[B]	CME	1	0
1	B	289[A]	CME	1	0
1	B	289[B]	CME	1	0
1	C	289[A]	CME	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 4 are 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	$\# Z > 2$
4	PEG	A	504	-	6,6,6	0.60	0	5,5,5	0.51	0
3	PGE	C	502	-	9,9,9	0.57	0	8,8,8	0.31	0
4	PEG	B	504	-	6,6,6	0.40	0	5,5,5	0.18	0
4	PEG	B	503	-	6,6,6	0.46	0	5,5,5	0.36	0
5	TRS	B	505	-	7,7,7	0.29	0	9,9,9	0.48	0
5	TRS	A	507	-	7,7,7	0.33	0	9,9,9	0.35	0
3	PGE	B	502	-	9,9,9	0.42	0	8,8,8	0.24	0
3	PGE	A	502	-	9,9,9	0.49	0	8,8,8	0.51	0
6	PG4	C	503	-	12,12,12	0.46	0	11,11,11	0.55	0
3	PGE	A	503	-	9,9,9	0.52	0	8,8,8	0.15	0
4	PEG	D	503	-	6,6,6	0.42	0	5,5,5	0.69	0
7	EDO	C	504[B]	-	3,3,3	0.51	0	2,2,2	0.14	0
4	PEG	A	505	-	6,6,6	0.47	0	5,5,5	0.17	0
4	PEG	A	506	-	6,6,6	0.44	0	5,5,5	0.24	0
3	PGE	D	502	-	9,9,9	0.52	0	8,8,8	0.56	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	504	-	-	1/4/4/4	-
3	PGE	C	502	-	-	5/7/7/7	-
4	PEG	B	504	-	-	3/4/4/4	-
4	PEG	B	503	-	-	3/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	TRS	B	505	-	-	4/9/9/9	-
5	TRS	A	507	-	-	5/9/9/9	-
3	PGE	B	502	-	-	4/7/7/7	-
3	PGE	A	502	-	-	4/7/7/7	-
6	PG4	C	503	-	-	4/10/10/10	-
3	PGE	A	503	-	-	1/7/7/7	-
4	PEG	D	503	-	-	1/4/4/4	-
7	EDO	C	504[B]	-	-	0/1/1/1	-
4	PEG	A	505	-	-	2/4/4/4	-
4	PEG	A	506	-	-	2/4/4/4	-
3	PGE	D	502	-	-	3/7/7/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	507	TRS	C2-C-C1-O1
5	A	507	TRS	C3-C-C1-O1
5	A	507	TRS	N-C-C1-O1
5	B	505	TRS	N-C-C3-O3
3	A	502	PGE	C6-C5-O3-C4
6	C	503	PG4	O2-C3-C4-O3
4	A	506	PEG	O2-C3-C4-O4
3	C	502	PGE	O3-C5-C6-O4
3	C	502	PGE	O2-C3-C4-O3
5	B	505	TRS	C2-C-C3-O3
4	A	504	PEG	O2-C3-C4-O4
3	B	502	PGE	O1-C1-C2-O2
3	B	502	PGE	O3-C5-C6-O4
6	C	503	PG4	O4-C7-C8-O5
3	A	502	PGE	O3-C5-C6-O4
3	C	502	PGE	O1-C1-C2-O2
6	C	503	PG4	O1-C1-C2-O2
4	A	505	PEG	O1-C1-C2-O2
4	B	504	PEG	O1-C1-C2-O2
4	D	503	PEG	O1-C1-C2-O2
3	A	502	PGE	C4-C3-O2-C2

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Mol	Chain	Res	Type	Atoms
3	B	502	PGE	C6-C5-O3-C4
3	D	502	PGE	C1-C2-O2-C3
3	C	502	PGE	C6-C5-O3-C4
3	A	502	PGE	C3-C4-O3-C5
4	B	503	PEG	O2-C3-C4-O4
4	B	504	PEG	O2-C3-C4-O4
4	B	504	PEG	C4-C3-O2-C2
4	B	503	PEG	C4-C3-O2-C2
4	B	503	PEG	O1-C1-C2-O2
6	C	503	PG4	C1-C2-O2-C3
4	A	505	PEG	C4-C3-O2-C2
4	A	506	PEG	O1-C1-C2-O2
3	B	502	PGE	C3-C4-O3-C5
5	A	507	TRS	C1-C-C3-O3
5	B	505	TRS	C1-C-C3-O3
3	A	503	PGE	C4-C3-O2-C2
3	D	502	PGE	C3-C4-O3-C5
3	D	502	PGE	O3-C5-C6-O4
5	A	507	TRS	N-C-C3-O3
5	B	505	TRS	C1-C-C2-O2
3	C	502	PGE	C3-C4-O3-C5

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	504	PEG	2	0
3	C	502	PGE	1	0
4	B	504	PEG	1	0
7	C	504[B]	EDO	3	0
3	D	502	PGE	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	496/520 (95%)	-0.43	8 (1%) 70 75	5, 14, 28, 69	33 (6%)
1	B	496/520 (95%)	-0.36	15 (3%) 52 57	5, 14, 32, 71	32 (6%)
1	C	495/520 (95%)	-0.43	10 (2%) 65 70	5, 15, 28, 100	37 (7%)
1	D	494/520 (95%)	-0.40	4 (0%) 82 87	5, 16, 30, 84	27 (5%)
All	All	1981/2080 (95%)	-0.41	37 (1%) 66 71	5, 15, 30, 100	129 (6%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	3	LEU	6.8
1	D	3	LEU	6.7
1	A	3	LEU	5.2
1	B	3	LEU	5.0
1	B	213	GLY	4.2
1	B	0	ALA	4.1
1	B	348	LYS	3.9
1	A	362	ASP	3.7
1	B	351	GLY	3.6
1	A	0	ALA	3.6
1	C	213	GLY	3.5
1	C	214	SER	3.5
1	C	1	MET	3.5
1	B	212	ALA	3.3
1	B	343	TYR	3.2
1	B	349	ALA	3.2
1	A	212	ALA	3.2
1	C	212	ALA	3.2
1	B	362	ASP	3.2
1	C	215	GLU	3.1
1	B	345	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	2	GLU	2.9
1	D	2	GLU	2.8
1	A	215[A]	GLU	2.8
1	A	213	GLY	2.4
1	D	362	ASP	2.4
1	D	367	LYS	2.3
1	C	409[A]	GLN	2.3
1	C	362	ASP	2.3
1	B	1	MET	2.3
1	B	2	GLU	2.2
1	B	215	GLU	2.2
1	B	214	SER	2.1
1	C	126[A]	ASP	2.1
1	A	1	MET	2.1
1	A	2	GLU	2.1
1	B	335	GLU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CME	D	289[A]	10/11	0.78	0.19	19,24,47,48	10
1	CME	D	289[B]	10/11	0.78	0.19	18,20,27,31	10
1	CME	A	289[A]	10/11	0.87	0.15	18,21,34,34	10
1	CME	A	289[B]	10/11	0.87	0.15	18,23,36,37	10
1	CME	B	289[A]	10/11	0.91	0.14	18,22,34,36	10
1	CME	B	289[B]	10/11	0.91	0.14	19,25,38,40	10
1	CME	C	289[A]	10/11	0.92	0.11	15,19,27,34	10
1	CME	C	289[B]	10/11	0.92	0.11	16,21,40,41	10

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	TRS	B	505	8/8	0.74	0.20	42,64,74,86	0
4	PEG	A	505	7/7	0.75	0.19	48,56,61,61	0
4	PEG	D	503	7/7	0.77	0.19	40,48,54,54	0
3	PGE	C	502	10/10	0.78	0.17	42,48,56,56	0
4	PEG	A	504	7/7	0.80	0.19	41,42,51,54	0
4	PEG	B	504	7/7	0.80	0.17	49,53,59,60	0
3	PGE	A	503	10/10	0.82	0.15	40,47,55,59	0
5	TRS	A	507	8/8	0.83	0.14	41,50,53,59	0
4	PEG	B	503	7/7	0.84	0.15	47,49,55,57	0
3	PGE	D	502	10/10	0.86	0.14	45,48,55,57	0
4	PEG	A	506	7/7	0.87	0.15	48,53,62,73	0
6	PG4	C	503	13/13	0.88	0.14	31,42,55,60	0
3	PGE	B	502	10/10	0.90	0.12	38,44,46,55	0
7	EDO	C	504[B]	4/4	0.90	0.12	20,21,22,22	4
3	PGE	A	502	10/10	0.93	0.10	34,36,48,50	0
2	NA	D	501	1/1	0.99	0.05	15,15,15,15	0
2	NA	A	501	1/1	0.99	0.04	11,11,11,11	0
2	NA	B	501	1/1	0.99	0.06	10,10,10,10	0
2	NA	C	501	1/1	0.99	0.03	11,11,11,11	0

6.5 Other polymers [i](#)

There are no such residues in this entry.