



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:07 PM GMT

PDB ID : 3SVW
Title : Crystal Structure of the P107V-MauG/pre-Methylamine Dehydrogenase Complex
Authors : Jensen, L.M.R.; Wilmot, C.M.
Deposited on : 2011-07-12
Resolution : 1.86 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

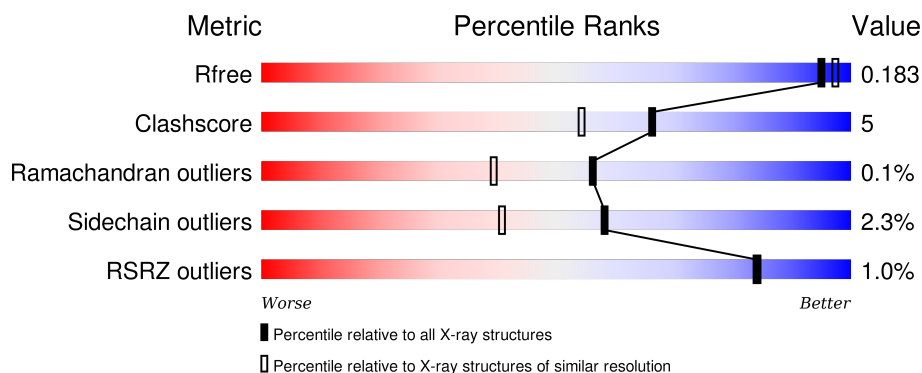
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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}	91344	1745 (1.86-1.86)
Clashscore	102246	1898 (1.86-1.86)
Ramachandran outliers	100387	1875 (1.86-1.86)
Sidechain outliers	100360	1875 (1.86-1.86)
RSRZ outliers	91569	1747 (1.86-1.86)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	373	<div> <div>83%</div> <div>7% • 10%</div> </div>
1	B	373	<div> <div>85%</div> <div>10% • 5%</div> </div>
2	C	137	<div> <div>80%</div> <div>15% • •</div> </div>
2	E	137	<div> <div>76%</div> <div>14% • 9%</div> </div>
3	D	386	<div> <div>89%</div> <div>8% • •</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	386	 <div>85% 12%</div>

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
11	MES	F	387	-	-	-	X
7	PGE	A	374	-	-	-	X
8	PEG	B	376	-	-	X	X
8	PEG	F	388	-	-	X	X
9	EDO	D	387	-	-	X	-
9	EDO	F	389	-	-	X	X

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 15652 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 Methylamine utilization protein MauG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	5	0
			2648	1663	472	500	13			
1	B	355	Total	C	N	O	S	0	5	0
			2774	1733	498	531	12			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	VAL	PRO	ENGINEERED MUTATION	UNP Q51658
A	368	HIS	-	EXPRESSION TAG	UNP Q51658
A	369	HIS	-	EXPRESSION TAG	UNP Q51658
A	370	HIS	-	EXPRESSION TAG	UNP Q51658
A	371	HIS	-	EXPRESSION TAG	UNP Q51658
A	372	HIS	-	EXPRESSION TAG	UNP Q51658
A	373	HIS	-	EXPRESSION TAG	UNP Q51658
B	107	VAL	PRO	ENGINEERED MUTATION	UNP Q51658
B	368	HIS	-	EXPRESSION TAG	UNP Q51658
B	369	HIS	-	EXPRESSION TAG	UNP Q51658
B	370	HIS	-	EXPRESSION TAG	UNP Q51658
B	371	HIS	-	EXPRESSION TAG	UNP Q51658
B	372	HIS	-	EXPRESSION TAG	UNP Q51658
B	373	HIS	-	EXPRESSION TAG	UNP Q51658

- Molecule 2 is a protein called Methylamine dehydrogenase light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	131	Total	C	N	O	S	0	3	0
			1027	634	179	200	14			
2	E	125	Total	C	N	O	S	0	3	0
			974	606	162	192	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	132	HIS	-	EXPRESSION TAG	UNP P22619
C	133	HIS	-	EXPRESSION TAG	UNP P22619
C	134	HIS	-	EXPRESSION TAG	UNP P22619
C	135	HIS	-	EXPRESSION TAG	UNP P22619
C	136	HIS	-	EXPRESSION TAG	UNP P22619
C	137	HIS	-	EXPRESSION TAG	UNP P22619
E	132	HIS	-	EXPRESSION TAG	UNP P22619
E	133	HIS	-	EXPRESSION TAG	UNP P22619
E	134	HIS	-	EXPRESSION TAG	UNP P22619
E	135	HIS	-	EXPRESSION TAG	UNP P22619
E	136	HIS	-	EXPRESSION TAG	UNP P22619
E	137	HIS	-	EXPRESSION TAG	UNP P22619

- Molecule 3 is a protein called Methylamine dehydrogenase heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	376	Total	C	N	O	S	0	2	0
			2932	1862	502	560	8			
3	F	376	Total	C	N	O	S	0	5	0
			2953	1876	506	563	8			

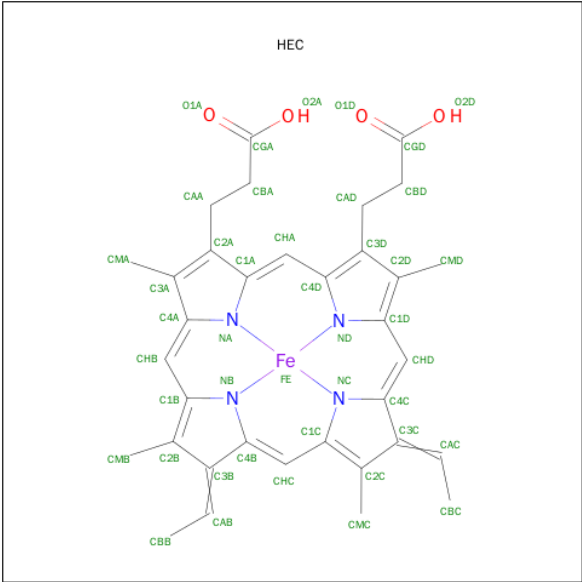
- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

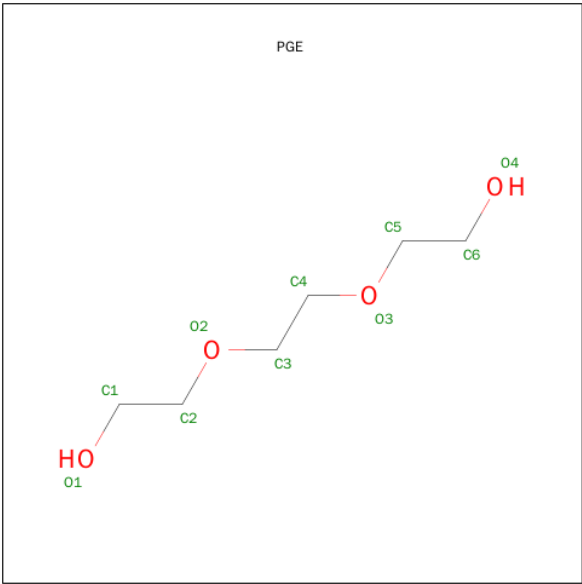
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Na	0	0
			2	2		
5	A	2	Total	Na	0	0
			2	2		

- Molecule 6 is HEME C (three-letter code: HEC) (formula: C₃₄H₃₄FeN₄O₄).



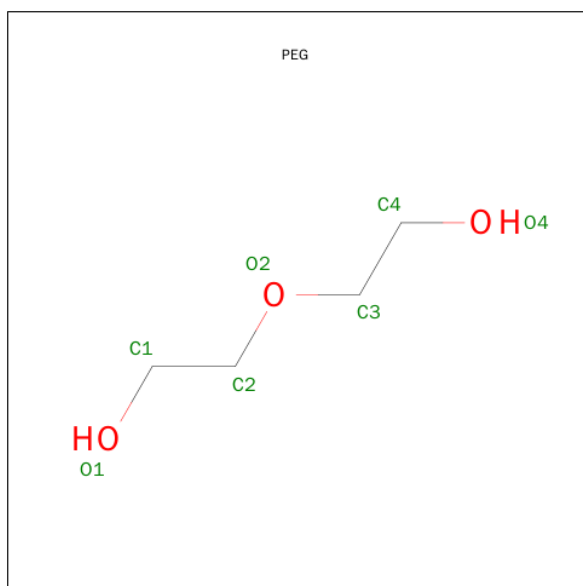
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
6	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
6	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
6	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



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

- Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



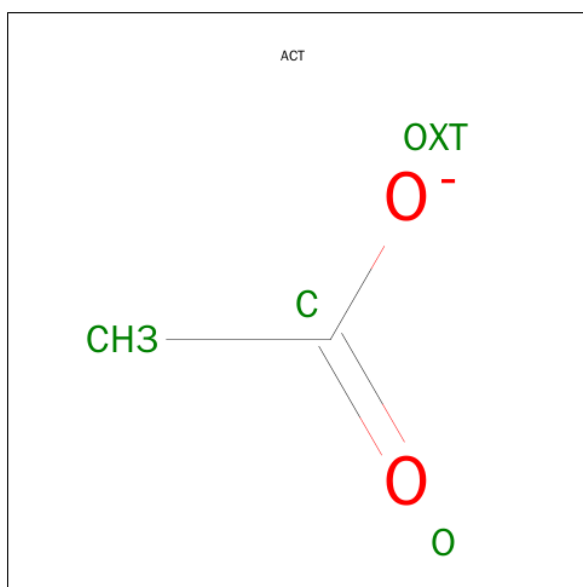
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			7	4	3		
8	B	1	Total	C	O	0	0
			7	4	3		
8	F	1	Total	C	O	0	0
			7	4	3		

- Molecule 9 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



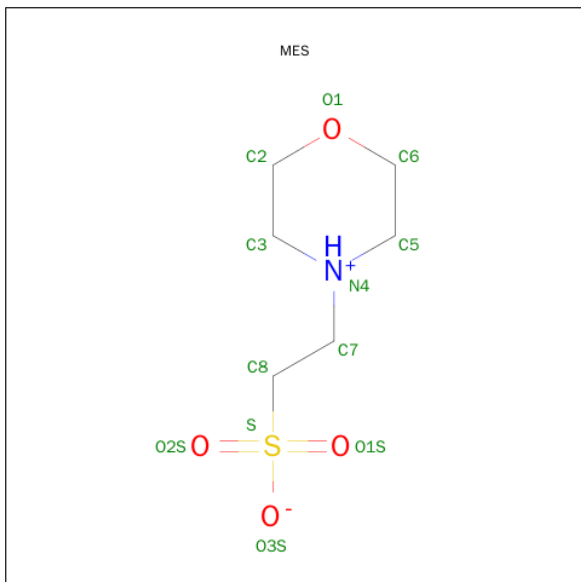
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			4	2	2		
9	D	1	Total	C	O	0	0
			4	2	2		
9	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 10 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 11 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	F	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

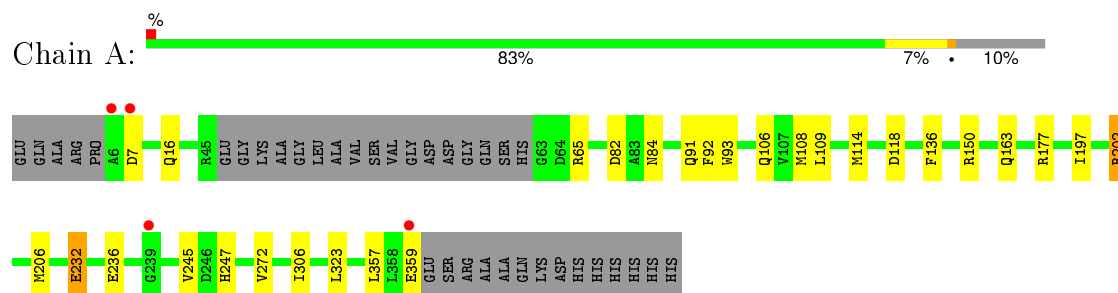
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	431	Total	O	0	0
			431	431		
12	B	438	Total	O	0	0
			438	438		
12	C	141	Total	O	0	0
			141	141		
12	D	410	Total	O	0	0
			410	410		
12	E	146	Total	O	0	0
			146	146		
12	F	541	Total	O	0	0
			541	541		

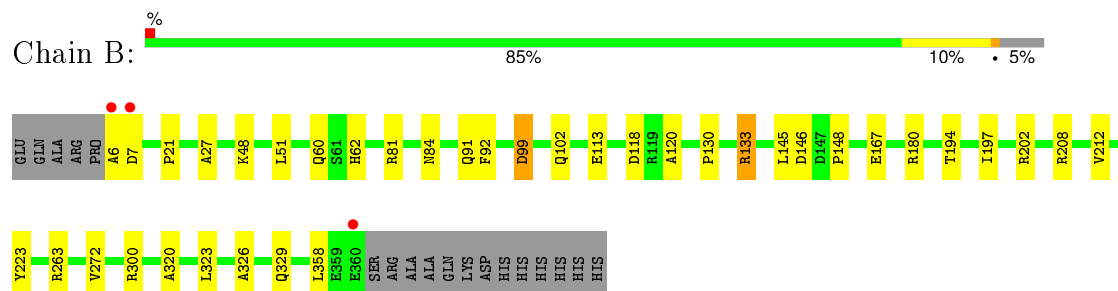
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

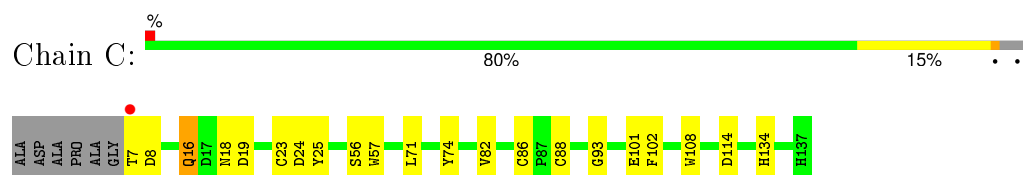
- Molecule 1: Methylamine utilization protein MauG



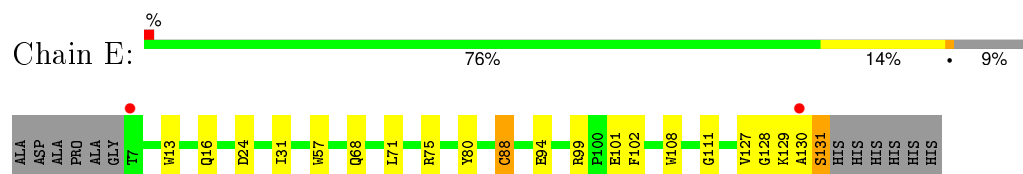
- Molecule 1: Methylamine utilization protein MauG



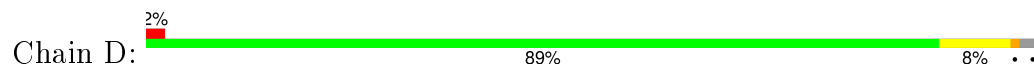
- Molecule 2: Methylamine dehydrogenase light chain

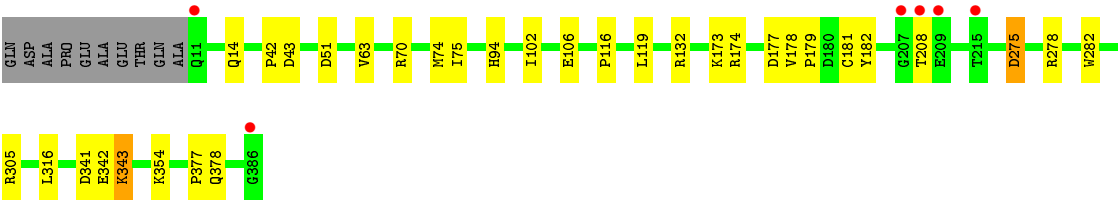


- Molecule 2: Methylamine dehydrogenase light chain

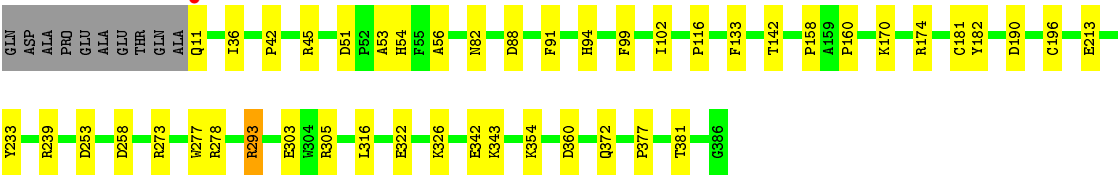
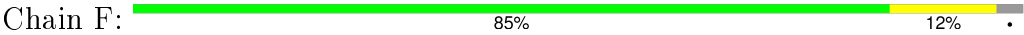


- Molecule 3: Methylamine dehydrogenase heavy chain





● Molecule 3: Methylamine dehydrogenase heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.61Å 89.00Å 104.81Å 67.05° 79.51° 79.72°	Depositor
Resolution (Å)	33.89 – 1.86 33.89 – 1.86	Depositor EDS
% Data completeness (in resolution range)	96.2 (33.89-1.86) 84.4 (33.89-1.86)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.99 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.131 , 0.174 0.141 , 0.183	Depositor DCC
R_{free} test set	7327 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	16.5	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 145457 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	15652	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 4.47% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, 0AF, NA, CA, HEC, EDO, MES, ACT, PEG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.13	1/2724 (0.0%)	0.97	6/3694 (0.2%)
1	B	1.16	2/2852 (0.1%)	0.97	6/3867 (0.2%)
2	C	1.22	2/1050 (0.2%)	0.98	2/1430 (0.1%)
2	E	1.19	3/994 (0.3%)	1.05	2/1358 (0.1%)
3	D	1.10	1/3015 (0.0%)	0.97	5/4109 (0.1%)
3	F	1.27	9/3045 (0.3%)	1.06	12/4148 (0.3%)
All	All	1.18	18/13680 (0.1%)	1.00	33/18606 (0.2%)

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	133	PHE	CE2-CZ	6.55	1.49	1.37
2	E	88[A]	CYS	CB-SG	-6.21	1.71	1.82
2	E	88[B]	CYS	CB-SG	-6.21	1.71	1.82
1	A	136	PHE	CE2-CZ	6.11	1.49	1.37
3	F	213	GLU	CD-OE1	5.90	1.32	1.25
3	F	277	TRP	CE3-CZ3	5.75	1.48	1.38
3	F	322	GLU	CG-CD	5.68	1.60	1.51
3	F	99	PHE	CE2-CZ	5.58	1.48	1.37
1	B	167	GLU	CB-CG	5.58	1.62	1.52
2	C	25	TYR	CD1-CE1	5.42	1.47	1.39
2	E	94	GLU	CB-CG	5.41	1.62	1.52
3	F	326	LYS	CE-NZ	5.30	1.62	1.49
1	B	223	TYR	CD1-CE1	5.28	1.47	1.39
3	F	342	GLU	C-O	5.26	1.33	1.23
3	D	63	VAL	CB-CG1	5.15	1.63	1.52
3	F	303	GLU	CG-CD	5.13	1.59	1.51
2	C	82	VAL	CB-CG1	5.05	1.63	1.52
3	F	91	PHE	CE2-CZ	5.04	1.47	1.37

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	81	ARG	NE-CZ-NH1	7.59	124.09	120.30
1	A	177	ARG	NE-CZ-NH1	7.58	124.09	120.30
1	A	82	ASP	CB-CG-OD2	7.54	125.09	118.30
3	F	278	ARG	NE-CZ-NH1	7.26	123.93	120.30
3	F	239	ARG	NE-CZ-NH1	7.08	123.84	120.30
2	C	24	ASP	CB-CG-OD1	6.97	124.57	118.30
1	B	99	ASP	CB-CG-OD2	6.80	124.42	118.30
1	A	150	ARG	NE-CZ-NH1	6.29	123.44	120.30
3	F	88	ASP	CB-CG-OD1	6.20	123.88	118.30
2	E	99	ARG	NE-CZ-NH1	6.17	123.39	120.30
1	B	263	ARG	NE-CZ-NH2	-6.14	117.23	120.30
3	F	190	ASP	CB-CG-OD1	6.13	123.82	118.30
3	F	360	ASP	CB-CG-OD1	6.06	123.75	118.30
3	F	170	LYS	CD-CE-NZ	-6.03	97.82	111.70
3	F	305	ARG	NE-CZ-NH2	-5.97	117.32	120.30
3	D	305	ARG	NE-CZ-NH2	-5.94	117.33	120.30
1	A	118	ASP	CB-CG-OD1	5.86	123.57	118.30
1	B	118	ASP	CB-CG-OD1	5.84	123.56	118.30
3	F	278	ARG	NE-CZ-NH2	-5.82	117.39	120.30
3	F	174	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	B	180	ARG	NE-CZ-NH2	5.72	123.16	120.30
3	D	275	ASP	CB-CG-OD2	-5.53	113.33	118.30
3	F	253	ASP	CB-CG-OD1	5.46	123.22	118.30
3	D	173	LYS	CD-CE-NZ	-5.45	99.17	111.70
1	A	82	ASP	CB-CG-OD1	-5.41	113.43	118.30
1	B	146	ASP	CB-CG-OD1	-5.39	113.45	118.30
3	D	43	ASP	CB-CG-OD1	5.38	123.14	118.30
3	F	293	ARG	NE-CZ-NH1	-5.19	117.70	120.30
1	A	65	ARG	NE-CZ-NH1	-5.18	117.71	120.30
2	C	8	ASP	CB-CG-OD1	5.17	122.96	118.30
3	F	133	PHE	CB-CG-CD2	-5.11	117.22	120.80
3	D	132	ARG	NE-CZ-NH2	-5.07	117.76	120.30
2	E	24	ASP	CB-CG-OD1	5.04	122.83	118.30

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	2648	0	2556	18	0
1	B	2774	0	2665	23	0
2	C	1027	0	916	21	0
2	E	974	0	880	18	0
3	D	2932	0	2828	16	0
3	F	2953	0	2857	24	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	86	0	60	5	0
6	B	86	0	60	3	0
7	A	10	0	14	3	0
8	B	14	0	20	12	0
8	F	7	0	10	9	0
9	B	4	0	6	0	0
9	D	4	0	6	5	0
9	F	4	0	6	6	0
10	D	4	0	3	0	0
11	F	12	0	12	0	0
12	A	431	0	0	6	0
12	B	438	0	0	5	0
12	C	141	0	0	4	0
12	D	410	0	0	6	0
12	E	146	0	0	5	0
12	F	541	0	0	13	0
All	All	15652	0	12899	129	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 (129) 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)
3:F:372[B]:GLN:NE2	12:F:1487:HOH:O	1.80	1.12
3:F:372[A]:GLN:OE1	12:F:2086:HOH:O	1.76	1.03
2:C:101[A]:GLU:HG3	12:C:2537:HOH:O	0.80	0.97
8:F:388:PEG:H22	12:F:479:HOH:O	1.68	0.91
9:D:387:EDO:H12	12:D:1241:HOH:O	1.75	0.86
8:F:388:PEG:H21	12:F:457:HOH:O	1.74	0.85
2:C:16:GLN:NE2	2:C:18:ASN:H	1.73	0.85
2:C:16:GLN:HE21	2:C:18:ASN:H	1.25	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:ALA:HB1	12:B:2480:HOH:O	1.78	0.82
8:F:388:PEG:H12	12:F:475:HOH:O	1.84	0.77
8:B:376:PEG:H22	2:E:111:GLY:O	1.86	0.76
3:F:233:TYR:O	8:F:388:PEG:H42	1.89	0.73
1:B:208[B]:ARG:NH1	12:B:1695:HOH:O	2.21	0.71
1:A:245[A]:VAL:HG13	1:A:247:HIS:NE2	2.06	0.71
1:A:93:TRP:HZ2	12:A:2279:HOH:O	1.74	0.70
8:F:388:PEG:H31	12:F:438:HOH:O	1.90	0.69
8:B:376:PEG:H32	2:E:111:GLY:O	1.93	0.68
8:B:376:PEG:H41	12:B:1767:HOH:O	1.94	0.67
1:B:197:ILE:O	1:B:202:ARG:HD2	1.95	0.67
2:C:16:GLN:HE22	2:C:19:ASP:H	1.41	0.67
1:A:306:ILE:HD11	12:A:1640:HOH:O	1.95	0.66
2:C:57:0AF:CE3	2:C:108:TRP:CD1	2.78	0.66
2:E:57:0AF:CE3	2:E:108:TRP:CD1	2.81	0.64
12:B:1704:HOH:O	2:E:127:VAL:HG12	1.98	0.64
2:C:93:GLY:O	2:C:134:HIS:HE1	1.80	0.63
1:B:358:LEU:O	8:B:374:PEG:H22	1.98	0.63
3:F:54:HIS:HE1	12:F:618:HOH:O	1.80	0.63
2:E:130:ALA:O	2:E:131:SER:HB2	1.99	0.62
2:C:86:CYS:H	3:F:372[A]:GLN:HE22	1.46	0.62
3:D:14:GLN:NE2	12:D:1592:HOH:O	2.33	0.61
1:A:16:GLN:NE2	12:A:1861:HOH:O	2.33	0.60
3:F:53:ALA:O	9:F:389:EDO:H11	2.01	0.60
8:F:388:PEG:C3	12:F:438:HOH:O	2.48	0.60
3:F:56:ALA:HB2	9:F:389:EDO:H11	1.84	0.60
2:C:101[B]:GLU:OE2	12:C:2002:HOH:O	2.16	0.60
12:D:1668:HOH:O	9:F:389:EDO:H22	2.00	0.60
3:D:378:GLN:HE22	9:D:387:EDO:C1	2.15	0.60
1:A:202:ARG:HB2	1:A:206[B]:MET:HG3	1.84	0.59
1:A:306:ILE:HD12	1:A:306:ILE:N	2.19	0.57
1:A:232:GLU:O	1:A:236:GLU:HG3	2.06	0.56
1:A:245[A]:VAL:CG1	1:A:247:HIS:NE2	2.69	0.55
3:F:273[A]:ARG:NH1	12:F:2361:HOH:O	2.36	0.55
3:D:278:ARG:NH1	12:D:1526:HOH:O	2.38	0.55
3:F:53:ALA:O	9:F:389:EDO:C1	2.55	0.55
3:D:343:LYS:HD2	3:D:343:LYS:N	2.21	0.54
1:A:106:GLN:HA	7:A:374:PGE:H2	1.89	0.54
12:A:1700:HOH:O	2:C:134:HIS:HD2	1.90	0.54
3:D:74:MET:O	3:D:75[A]:ILE:HD13	2.08	0.53
3:F:82:ASN:HB3	3:F:142:THR:HB	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:D:1668:HOH:O	9:F:389:EDO:C2	2.57	0.53
3:D:378:GLN:HE22	9:D:387:EDO:H11	1.73	0.53
8:B:376:PEG:C3	2:E:111:GLY:O	2.56	0.53
3:D:378:GLN:NE2	9:D:387:EDO:H11	2.23	0.53
1:B:130:PRO:HA	1:B:133:ARG:HH21	1.73	0.52
1:B:326:ALA:O	8:B:376:PEG:H31	2.10	0.52
8:B:376:PEG:C2	2:E:111:GLY:O	2.56	0.52
3:D:51:ASP:HA	3:D:377:PRO:HA	1.93	0.51
2:C:101[B]:GLU:HG2	2:C:102:PHE:CD2	2.46	0.51
3:F:381:THR:HG23	8:F:388:PEG:H11	1.94	0.50
1:B:329:GLN:O	8:B:376:PEG:H42	2.11	0.50
3:D:282:TRP:CH2	9:D:387:EDO:H22	2.47	0.50
1:B:194:THR:HG21	2:E:101[B]:GLU:HG3	1.93	0.50
6:A:600:HEC:HBC3	6:A:600:HEC:HMC1	1.93	0.50
3:D:75[A]:ILE:HD11	3:D:119:LEU:HB3	1.93	0.49
1:A:91:GLN:O	1:A:92:PHE:HB2	2.13	0.49
2:E:31:ILE:CG2	2:E:88[B]:CYS:SG	3.01	0.49
2:C:16:GLN:NE2	2:C:18:ASN:N	2.52	0.49
1:B:197:ILE:O	1:B:202:ARG:CD	2.61	0.49
2:C:57:0AF:HE3	2:C:108:TRP:CD1	2.47	0.49
3:D:42:PRO:HG3	3:D:116:PRO:HB2	1.95	0.49
12:C:1671:HOH:O	3:F:36[A]:ILE:HD11	2.12	0.49
3:F:51:ASP:HA	3:F:377:PRO:HA	1.95	0.48
1:A:323[B]:LEU:HD12	12:A:1525:HOH:O	2.12	0.48
2:C:114:ASP:O	12:C:1564:HOH:O	2.20	0.48
2:C:86:CYS:H	3:F:372[A]:GLN:NE2	2.12	0.47
1:A:114[A]:MET:SD	6:A:500:HEC:HMC2	2.55	0.47
8:F:388:PEG:H11	12:F:606:HOH:O	2.14	0.47
1:A:93:TRP:CZ2	12:A:2279:HOH:O	2.54	0.47
1:B:48:LYS:HE2	1:B:62:HIS:CE1	2.49	0.47
1:B:113:GLU:HG2	6:B:500:HEC:HBC2	1.98	0.46
1:B:99:ASP:OD2	1:B:102[A]:GLN:HG2	2.15	0.46
1:A:109:LEU:HD12	7:A:374:PGE:H22	1.98	0.46
1:A:272:VAL:HG21	6:A:600:HEC:HMA3	1.96	0.46
1:B:120:ALA:HA	1:B:148:PRO:HB3	1.96	0.46
2:C:23:CYS:SG	2:C:88[B]:CYS:HB2	2.55	0.46
2:E:75:ARG:HD2	12:E:1628:HOH:O	2.14	0.46
2:E:101[B]:GLU:HG2	2:E:102:PHE:CD2	2.50	0.46
2:C:23:CYS:HB3	2:C:88[B]:CYS:SG	2.54	0.46
2:E:68:GLN:OE1	2:E:129:LYS:HE2	2.16	0.46
1:A:108:MET:HA	1:A:114[A]:MET:HB2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:57:0AF:HE3	2:E:108:TRP:CD1	2.51	0.45
3:F:293:ARG:HG3	12:F:1175:HOH:O	2.16	0.44
1:B:326:ALA:O	8:B:376:PEG:C3	2.65	0.44
8:B:376:PEG:H11	12:E:884:HOH:O	2.17	0.44
3:F:181:CYS:HA	3:F:196:CYS:HA	1.99	0.44
3:F:42:PRO:HG3	3:F:116:PRO:HB2	1.99	0.44
1:B:21:PRO:O	1:B:27:ALA:HA	2.17	0.44
1:B:272:VAL:HG21	6:B:600:HEC:HMA3	2.00	0.44
1:A:163:GLN:HE22	6:A:500:HEC:HMA1	1.83	0.44
1:B:300[A]:ARG:HD3	3:F:158:PRO:HG3	2.00	0.44
3:D:14:GLN:HE21	3:D:70:ARG:HH11	1.65	0.44
3:F:53:ALA:HB1	9:F:389:EDO:H12	1.99	0.43
3:F:258:ASP:OD1	12:F:2107:HOH:O	2.21	0.43
1:B:91:GLN:O	1:B:92:PHE:HB2	2.17	0.43
2:E:75:ARG:HA	12:E:625:HOH:O	2.19	0.43
3:D:178:VAL:HB	3:D:179:PRO:CD	2.48	0.43
6:A:500:HEC:HMC1	6:A:500:HEC:HBC3	2.01	0.42
3:F:45:ARG:NH2	3:F:343:LYS:O	2.52	0.42
2:E:13[B]:TRP:HB3	12:E:1442:HOH:O	2.18	0.42
1:B:212:VAL:HG12	2:E:128:GLY:HA2	2.02	0.42
1:B:323:LEU:HD23	12:B:410:HOH:O	2.18	0.42
2:C:23:CYS:CB	2:C:88[B]:CYS:SG	3.08	0.41
7:A:374:PGE:H32	7:A:374:PGE:H12	1.78	0.41
6:B:600:HEC:HBC3	6:B:600:HEC:HMC1	2.02	0.41
2:C:93:GLY:O	2:C:134:HIS:CE1	2.67	0.41
3:F:381:THR:HG23	8:F:388:PEG:C1	2.50	0.41
1:A:197:ILE:HD13	2:C:71:LEU:HD13	2.03	0.41
1:B:329:GLN:O	8:B:376:PEG:C4	2.68	0.41
2:C:56[A]:SER:HB2	2:C:74:TYR:O	2.21	0.41
1:B:133:ARG:HG3	1:B:145:LEU:HD12	2.03	0.41
3:D:341:ASP:N	3:D:341:ASP:OD1	2.51	0.40
3:D:181:CYS:C	3:D:182:TYR:CD1	2.95	0.40
3:F:181:CYS:C	3:F:182:TYR:CD1	2.95	0.40
2:E:13[B]:TRP:HD1	12:E:1442:HOH:O	2.04	0.40
2:C:57:0AF:CE3	2:C:108:TRP:HD1	2.32	0.40
1:B:320:ALA:O	1:B:323:LEU:CD2	2.69	0.40
3:D:174:ARG:NH2	12:D:1746:HOH:O	2.49	0.40
3:F:372[A]:GLN:NE2	12:F:1487:HOH:O	2.47	0.40
8:B:376:PEG:H32	2:E:111:GLY:C	2.42	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	338/373 (91%)	330 (98%)	8 (2%)	0	100	100
1	B	358/373 (96%)	350 (98%)	8 (2%)	0	100	100
2	C	131/137 (96%)	127 (97%)	4 (3%)	0	100	100
2	E	125/137 (91%)	123 (98%)	2 (2%)	0	100	100
3	D	376/386 (97%)	365 (97%)	10 (3%)	1 (0%)	46	29
3	F	379/386 (98%)	366 (97%)	12 (3%)	1 (0%)	46	29
All	All	1707/1792 (95%)	1661 (97%)	44 (3%)	2 (0%)	56	39

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	D	102	ILE
3	F	102	ILE

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	270/292 (92%)	264 (98%)	6 (2%)	60	43
1	B	282/292 (97%)	277 (98%)	5 (2%)	66	52
2	C	113/112 (101%)	111 (98%)	2 (2%)	66	52
2	E	107/112 (96%)	103 (96%)	4 (4%)	41	20
3	D	306/311 (98%)	297 (97%)	9 (3%)	50	31

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	F	309/311 (99%)	304 (98%)	5 (2%)	70	57
All	All	1387/1430 (97%)	1356 (98%)	31 (2%)	58	43

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	84	ASN
1	A	202	ARG
1	A	232	GLU
1	A	357	LEU
1	A	359	GLU
1	B	7	ASP
1	B	51	LEU
1	B	60	GLN
1	B	84	ASN
1	B	133	ARG
2	C	7	THR
2	C	16	GLN
3	D	94	HIS
3	D	106	GLU
3	D	177	ASP
3	D	208	THR
3	D	275	ASP
3	D	316	LEU
3	D	342	GLU
3	D	343	LYS
3	D	354	LYS
2	E	16	GLN
2	E	71	LEU
2	E	80	TYR
2	E	131	SER
3	F	11	GLN
3	F	94	HIS
3	F	160	PRO
3	F	316	LEU
3	F	354	LYS

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

Mol	Chain	Res	Type
1	A	84	ASN
1	A	163	GLN
1	B	16	GLN
1	B	26	ASN
1	B	60	GLN
1	B	91	GLN
2	C	16	GLN
2	C	134	HIS
2	C	136	HIS
3	D	14	GLN
3	D	30	GLN
3	F	54	HIS
3	F	300	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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$
2	0AF	C	57	2	13,16,17	1.60	2 (15%)	10,22,24	1.41	1 (10%)
2	0AF	E	57	2	13,16,17	1.73	3 (23%)	10,22,24	1.41	1 (10%)

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	0AF	C	57	2	-	0/3/6/8	0/2/2/2
2	0AF	E	57	2	-	0/3/6/8	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	57	0AF	CB-CA	-3.84	1.45	1.53
2	E	57	0AF	CB-CA	-2.89	1.47	1.53
2	C	57	0AF	CZ3-CE3	2.71	1.42	1.36
2	E	57	0AF	CD1-CG	2.78	1.43	1.38
2	E	57	0AF	CZ3-CE3	2.89	1.43	1.36

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	57	0AF	O-C-CA	-4.12	114.77	125.49
2	C	57	0AF	O-C-CA	-3.57	116.19	125.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	57	0AF	3	0
2	E	57	0AF	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 6 are monoatomic - leaving 13 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	$\# Z > 2$
7	PGE	A	374	-	9,9,9	0.39	0	8,8,8	0.93	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	HEC	A	500	1	24,50,50	1.69	9 (37%)	19,82,82	3.27	7 (36%)
6	HEC	A	600	1	24,50,50	1.51	5 (20%)	19,82,82	3.88	9 (47%)
8	PEG	B	374	-	6,6,6	0.58	0	5,5,5	0.85	0
9	EDO	B	375	-	3,3,3	0.37	0	2,2,2	0.27	0
8	PEG	B	376	-	6,6,6	0.29	0	5,5,5	0.92	0
6	HEC	B	500	1	24,50,50	1.75	9 (37%)	19,82,82	3.09	6 (31%)
6	HEC	B	600	1	24,50,50	1.83	10 (41%)	19,82,82	3.73	11 (57%)
9	EDO	D	387	-	3,3,3	1.58	0	2,2,2	1.78	1 (50%)
10	ACT	D	388	-	1,3,3	1.43	0	0,3,3	0.00	-
11	MES	F	387	-	11,12,12	0.70	0	14,16,16	2.95	5 (35%)
8	PEG	F	388	-	6,6,6	0.51	0	5,5,5	1.34	1 (20%)
9	EDO	F	389	-	3,3,3	0.28	0	2,2,2	1.45	1 (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
7	PGE	A	374	-	-	0/7/7/7	0/0/0/0
6	HEC	A	500	1	-	0/6/54/54	0/0/8/8
6	HEC	A	600	1	-	0/6/54/54	0/0/8/8
8	PEG	B	374	-	-	0/4/4/4	0/0/0/0
9	EDO	B	375	-	-	0/1/1/1	0/0/0/0
8	PEG	B	376	-	-	0/4/4/4	0/0/0/0
6	HEC	B	500	1	-	0/6/54/54	0/0/8/8
6	HEC	B	600	1	-	0/6/54/54	0/0/8/8
9	EDO	D	387	-	-	0/1/1/1	0/0/0/0
10	ACT	D	388	-	-	0/0/0/0	0/0/0/0
11	MES	F	387	-	-	0/6/14/14	0/1/1/1
8	PEG	F	388	-	-	0/4/4/4	0/0/0/0
9	EDO	F	389	-	-	0/1/1/1	0/0/0/0

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	600	HEC	C4B-NB	-4.16	1.31	1.36
6	A	500	HEC	C1A-NA	-3.52	1.31	1.36
6	B	500	HEC	C4B-NB	-3.38	1.32	1.36
6	B	500	HEC	C4A-NA	-2.92	1.32	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	600	HEC	C4A-NA	-2.89	1.32	1.36
6	A	500	HEC	C4A-NA	-2.60	1.33	1.36
6	B	600	HEC	C4C-NC	-2.53	1.33	1.36
6	B	600	HEC	C3B-C2B	-2.50	1.38	1.40
6	B	600	HEC	C3C-C2C	-2.43	1.38	1.40
6	B	500	HEC	C3C-C2C	-2.38	1.38	1.40
6	B	500	HEC	C4C-NC	-2.37	1.33	1.36
6	A	600	HEC	C1A-NA	-2.36	1.33	1.36
6	A	500	HEC	C4C-NC	-2.25	1.33	1.36
6	A	500	HEC	C4B-NB	-2.16	1.33	1.36
6	A	500	HEC	C1D-CHD	2.07	1.45	1.39
6	B	500	HEC	C1D-CHD	2.11	1.45	1.39
6	B	600	HEC	C1B-CHB	2.25	1.46	1.39
6	A	500	HEC	C3B-C4B	2.29	1.48	1.42
6	A	500	HEC	C1C-CHC	2.34	1.46	1.39
6	A	600	HEC	C1D-CHD	2.35	1.46	1.39
6	B	500	HEC	C3B-C4B	2.38	1.48	1.42
6	B	600	HEC	C1D-CHD	2.41	1.46	1.39
6	B	600	HEC	C1C-CHC	2.45	1.46	1.39
6	B	600	HEC	CAD-C3D	2.46	1.56	1.52
6	A	500	HEC	C3C-C4C	2.47	1.48	1.42
6	A	600	HEC	C1C-CHC	2.48	1.46	1.39
6	A	600	HEC	C1B-CHB	2.54	1.46	1.39
6	B	600	HEC	C4D-CHA	2.58	1.46	1.39
6	B	500	HEC	C4D-CHA	2.66	1.47	1.39
6	B	500	HEC	C3C-C4C	2.68	1.48	1.42
6	B	500	HEC	C1C-CHC	2.75	1.47	1.39
6	A	600	HEC	C4D-CHA	2.83	1.47	1.39
6	A	500	HEC	C1B-CHB	2.94	1.47	1.39

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	600	HEC	CBB-CAB-C3B	-10.43	104.18	127.35
6	B	500	HEC	CBB-CAB-C3B	-9.37	106.53	127.35
6	A	500	HEC	CBB-CAB-C3B	-8.83	107.73	127.35
6	B	600	HEC	CBB-CAB-C3B	-8.00	109.58	127.35
6	B	600	HEC	CBD-CAD-C3D	-7.66	98.81	112.53
6	A	500	HEC	CBC-CAC-C3C	-7.61	110.44	127.35
6	B	600	HEC	CBC-CAC-C3C	-7.57	110.53	127.35
6	A	600	HEC	CBD-CAD-C3D	-7.53	99.03	112.53
6	A	600	HEC	CBC-CAC-C3C	-7.12	111.53	127.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	600	HEC	CBA-CAA-C2A	-5.33	102.98	112.53
11	F	387	MES	O1S-S-C8	-5.15	102.51	106.91
6	B	500	HEC	CBC-CAC-C3C	-4.92	116.41	127.35
6	B	500	HEC	CBD-CAD-C3D	-4.23	104.94	112.53
6	A	500	HEC	CBD-CAD-C3D	-4.17	105.05	112.53
6	B	500	HEC	CAA-CBA-CGA	-3.92	105.56	112.75
6	A	600	HEC	C4C-C3C-C2C	-3.92	102.12	106.35
11	F	387	MES	C2-C3-N4	-3.92	104.19	110.12
6	A	600	HEC	CBA-CAA-C2A	-3.82	105.68	112.53
6	B	600	HEC	C4B-C3B-C2B	-3.53	102.54	106.35
6	A	600	HEC	CAD-CBD-CGD	-3.26	106.77	112.75
6	B	500	HEC	CBA-CAA-C2A	-3.15	106.89	112.53
6	A	500	HEC	C4B-C3B-C2B	-2.81	103.31	106.35
6	B	600	HEC	C4C-C3C-C2C	-2.73	103.40	106.35
6	B	600	HEC	CMC-C2C-C1C	-2.70	123.90	128.36
6	A	600	HEC	CMB-C2B-C1B	-2.24	124.65	128.36
6	A	500	HEC	C4C-C3C-C2C	-2.14	104.04	106.35
6	A	600	HEC	CMC-C2C-C1C	-2.13	124.83	128.36
6	B	600	HEC	CAD-C3D-C4D	-2.06	124.77	127.01
9	F	389	EDO	O2-C2-C1	-2.01	98.16	112.54
6	B	600	HEC	CMB-C2B-C1B	-2.01	125.05	128.36
6	B	600	HEC	CMA-C3A-C2A	2.19	129.81	125.24
6	A	500	HEC	CMA-C3A-C2A	2.28	130.00	125.24
8	F	388	PEG	O2-C2-C1	2.32	121.13	110.43
9	D	387	EDO	O1-C1-C2	2.46	130.14	112.54
6	B	600	HEC	C3C-C4C-NC	2.66	115.97	110.94
11	F	387	MES	O2S-S-C8	2.75	109.25	106.91
6	A	600	HEC	CMD-C2D-C3D	2.93	131.37	125.24
6	B	500	HEC	CAA-C2A-C1A	3.32	130.61	127.01
11	F	387	MES	C7-N4-C5	3.91	121.30	111.27
6	A	500	HEC	CAD-C3D-C4D	4.13	131.49	127.01
11	F	387	MES	C5-N4-C3	6.64	123.29	108.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 43 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	374	PGE	3	0
6	A	500	HEC	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	600	HEC	2	0
8	B	374	PEG	1	0
8	B	376	PEG	11	0
6	B	500	HEC	1	0
6	B	600	HEC	2	0
9	D	387	EDO	5	0
8	F	388	PEG	9	0
9	F	389	EDO	6	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 ⓘ

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	337/373 (90%)	-0.54	4 (1%) 81 81	11, 19, 34, 50	0
1	B	355/373 (95%)	-0.60	3 (0%) 87 87	9, 17, 33, 56	0
2	C	130/137 (94%)	-0.57	1 (0%) 87 87	10, 16, 30, 42	0
2	E	124/137 (90%)	-0.63	2 (1%) 74 74	10, 14, 25, 52	0
3	D	376/386 (97%)	-0.34	6 (1%) 74 74	11, 21, 37, 64	0
3	F	376/386 (97%)	-0.56	1 (0%) 94 93	9, 13, 26, 49	0
All	All	1698/1792 (94%)	-0.52	17 (1%) 84 84	9, 17, 33, 64	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	6	ALA	5.6
1	A	6	ALA	5.5
1	B	7	ASP	4.0
3	D	208	THR	4.0
3	D	207	GLY	3.8
3	F	11	GLN	3.8
1	A	7	ASP	3.8
3	D	11	GLN	3.5
3	D	209	GLU	3.1
1	A	239	GLY	2.8
3	D	386	GLY	2.6
2	C	7	THR	2.5
1	A	359	GLU	2.5
1	B	360	GLU	2.4
2	E	130	ALA	2.3
3	D	215	THR	2.3
2	E	7	THR	2.0

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	0AF	E	57	15/16	0.97	0.08	-	13,16,26,27	0
2	0AF	C	57	15/16	0.96	0.09	-	16,18,25,25	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
9	EDO	F	389	4/4	0.94	0.26	23.79	27,33,38,47	0
8	PEG	B	376	7/7	0.92	0.28	13.22	25,33,43,50	0
7	PGE	A	374	10/10	0.94	0.14	6.59	34,39,44,45	0
11	MES	F	387	12/12	0.98	0.10	4.52	23,37,42,43	0
8	PEG	F	388	7/7	0.91	0.15	2.08	31,37,39,41	0
9	EDO	B	375	4/4	0.93	0.09	1.55	28,33,35,40	0
9	EDO	D	387	4/4	0.92	0.14	0.88	26,28,31,33	0
5	NA	A	402	1/1	0.98	0.07	0.50	22,22,22,22	0
6	HEC	B	500	43/43	0.99	0.09	0.41	7,12,14,17	0
6	HEC	A	600	43/43	0.99	0.09	0.40	10,14,15,15	0
6	HEC	A	500	43/43	0.99	0.10	0.31	9,12,16,18	0
6	HEC	B	600	43/43	0.99	0.10	0.22	4,10,11,14	0
4	CA	B	400	1/1	1.00	0.08	0.19	13,13,13,13	0
4	CA	A	400	1/1	1.00	0.02	-2.36	17,17,17,17	0
5	NA	B	402	1/1	0.99	0.03	-3.31	23,23,23,23	0
5	NA	A	401	1/1	0.97	0.15	-	30,30,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
10	ACT	D	388	4/4	0.98	0.07	-	20,21,24,25	0
8	PEG	B	374	7/7	0.86	0.11	-	39,40,42,44	0
5	NA	B	401	1/1	0.98	0.10	-	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.