



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 06:14 AM UTC

PDB ID : 3EIF / pdb_00003eif
Title : 1.9 angstrom crystal structure of the active form of the C5a peptidase from Streptococcus pyogenes (ScpA)
Authors : Cooney, J.C.; Kagawa, T.F.; O'Connell, M.R.; Paoli, M.; Mouat, P.; O'Toole, P.W.
Deposited on : 2008-09-15
Resolution : 1.90 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

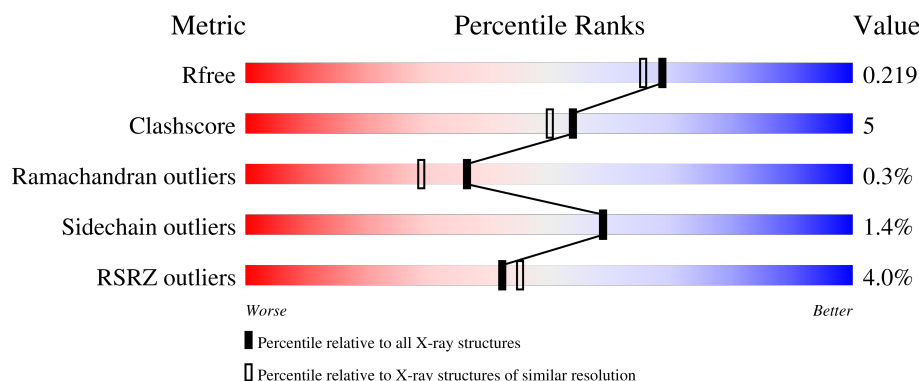
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	936	

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
4	SO4	A	7	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8105 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 C5a peptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	936	Total	C	N	O	S	0	5	0
			7129	4482	1194	1434	19			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	119	SER	GLY	engineered mutation	UNP B6ETQ5

- 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		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



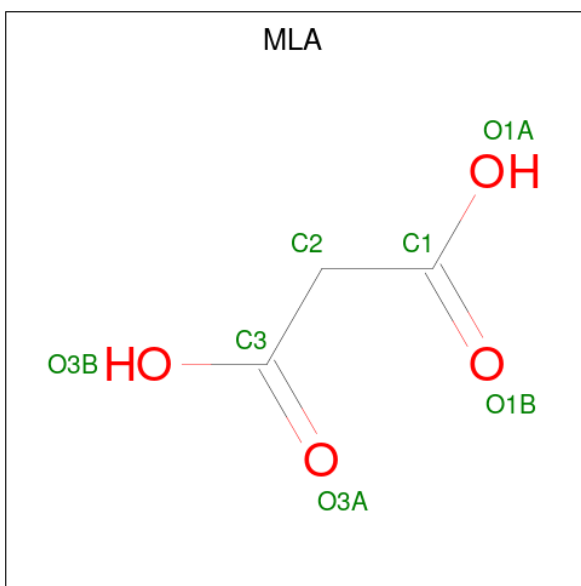
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



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

- Molecule 6 is MALONIC ACID (CCD ID: MLA) (formula: $C_3H_4O_4$).



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

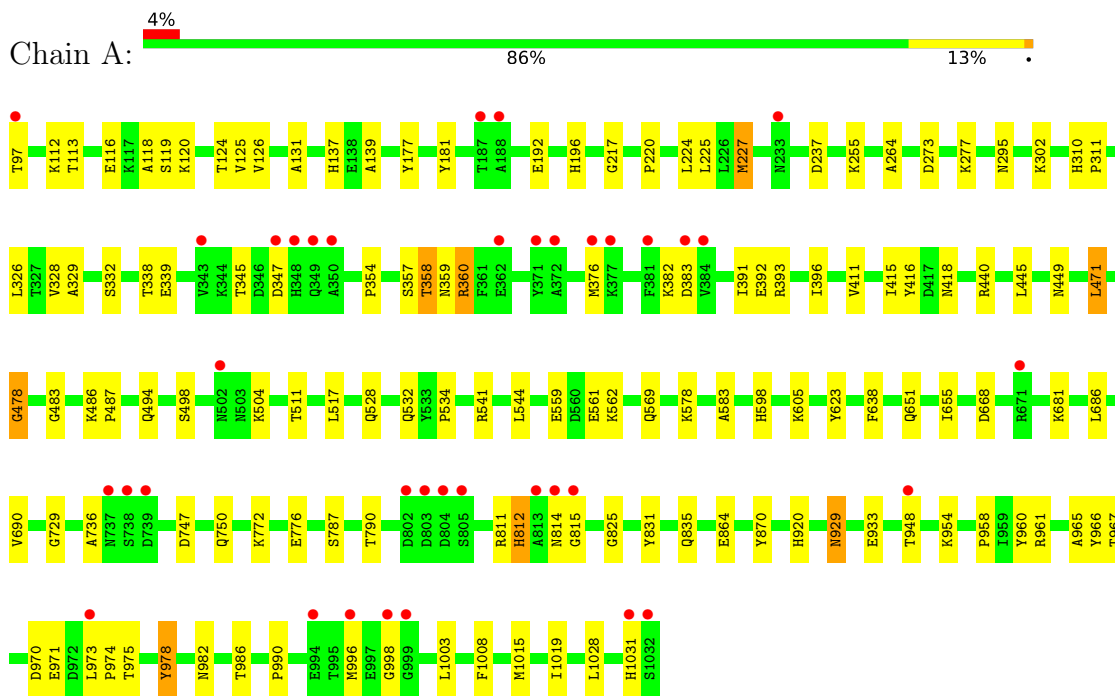
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	917	Total 917	O 917	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: C5a peptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	167.56Å 167.56Å 141.76Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	51.15 – 1.90 51.15 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (51.15-1.90) 99.9 (51.15-1.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 1.90Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.191 , 0.223 0.189 , 0.219	Depositor DCC
R_{free} test set	4619 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8105	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% 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: EPE, CA, MLA, SO4, NA

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	0.45	0/7298	0.96	30/9909 (0.3%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	681	LYS	N-CA-C	6.92	119.67	111.71
1	A	339	GLU	N-CA-C	-6.86	101.47	110.53
1	A	864	GLU	N-CA-C	6.59	118.46	111.28
1	A	690	VAL	N-CA-C	-6.58	98.38	107.99
1	A	638	PHE	N-CA-C	-6.50	99.92	110.20
1	A	478	GLY	N-CA-C	-6.45	102.74	112.41
1	A	790	THR	N-CA-C	-6.42	99.55	109.76
1	A	310	HIS	CA-C-N	6.25	125.94	119.56
1	A	310	HIS	C-N-CA	6.25	125.94	119.56
1	A	360	ARG	N-CA-C	6.17	119.19	108.76
1	A	511	THR	N-CA-C	-6.00	105.80	113.01
1	A	328	VAL	N-CA-C	5.82	116.31	108.17
1	A	982	ASN	N-CA-C	-5.62	103.28	110.53
1	A	870	TYR	N-CA-C	5.59	119.85	113.19
1	A	264	ALA	N-CA-C	5.58	117.49	110.24
1	A	358	THR	N-CA-C	-5.52	100.19	109.07
1	A	357	SER	N-CA-C	5.43	118.67	109.76
1	A	131	ALA	N-CA-C	5.41	119.73	113.18
1	A	686	LEU	N-CA-C	-5.30	98.74	108.02
1	A	812	HIS	N-CA-C	-5.28	103.91	110.41
1	A	311	PRO	N-CA-C	5.26	120.68	113.84
1	A	583	ALA	N-CA-C	5.22	117.88	110.10
1	A	192	GLU	N-CA-C	5.19	116.63	110.97
1	A	1019	ILE	N-CA-C	5.17	115.82	108.42

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	623	TYR	N-CA-C	-5.06	101.44	109.59
1	A	835	GLN	CA-C-N	-5.05	117.72	121.61
1	A	835	GLN	C-N-CA	-5.05	117.72	121.61
1	A	120	LYS	N-CA-C	5.03	118.95	112.41
1	A	978	TYR	N-CA-C	5.03	117.31	109.52
1	A	825	GLY	N-CA-C	5.02	122.53	115.30

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	7129	0	6827	69	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	35	0	0	2	0
5	A	15	0	18	0	0
6	A	7	0	2	0	0
7	A	917	0	0	5	0
All	All	8105	0	6847	69	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 (69) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:811:ARG:HD3	1:A:815:GLY:O	1.56	1.05
1:A:137:HIS:HD2	1:A:139:ALA:H	1.28	0.79
1:A:338:THR:HG21	1:A:354:PRO:HB3	1.66	0.78
1:A:948:THR:HG22	1:A:986:THR:HG22	1.74	0.70
1:A:948:THR:HG23	7:A:1232:HOH:O	1.94	0.68
1:A:471:LEU:HD22	1:A:569:GLN:HG2	1.79	0.63
1:A:295:ASN:HB3	4:A:7:SO4:O3	1.99	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:747:ASP:O	1:A:750:GLN:HG2	2.00	0.62
1:A:445:LEU:C	1:A:445:LEU:HD23	2.25	0.61
1:A:181:TYR:HE1	1:A:227[B]:MET:HG2	1.68	0.58
1:A:920:HIS:HD2	7:A:1388:HOH:O	1.87	0.58
1:A:973:LEU:C	1:A:973:LEU:HD13	2.29	0.58
1:A:137:HIS:CD2	1:A:139:ALA:H	2.16	0.57
1:A:332:SER:HB2	1:A:494:GLN:HG3	1.88	0.55
1:A:787:SER:HB3	1:A:973:LEU:HD11	1.89	0.55
1:A:358:THR:O	1:A:359:ASN:HB2	2.07	0.53
1:A:605:LYS:HE2	1:A:668:ASP:OD2	2.09	0.53
1:A:534:PRO:HA	1:A:541:ARG:HH21	1.74	0.52
1:A:112:LYS:O	1:A:116:GLU:HG3	2.11	0.50
1:A:528:GLN:HE21	1:A:532:GLN:NE2	2.09	0.50
1:A:295:ASN:HB3	4:A:7:SO4:S	2.51	0.50
1:A:392:GLU:HB2	1:A:416:TYR:CZ	2.47	0.50
1:A:948:THR:HG22	1:A:986:THR:CG2	2.40	0.50
1:A:965:ALA:HB2	1:A:978:TYR:CE2	2.46	0.49
1:A:359:ASN:O	1:A:360:ARG:HD2	2.13	0.49
1:A:376:MET:HE3	1:A:396:ILE:HG23	1.94	0.48
1:A:391:ILE:O	1:A:415:ILE:HA	2.14	0.48
1:A:137:HIS:HE1	1:A:504:LYS:O	1.97	0.47
1:A:273:ASP:O	1:A:277:LYS:HG2	2.14	0.47
1:A:967:THR:HA	1:A:975:THR:O	2.14	0.47
1:A:812:HIS:HB2	1:A:814:ASN:OD1	2.15	0.46
1:A:124:THR:HG22	1:A:125:VAL:N	2.30	0.45
1:A:376:MET:HE3	1:A:396:ILE:CG2	2.46	0.45
1:A:217:GLY:O	1:A:220:PRO:HD3	2.16	0.45
1:A:345:THR:HA	1:A:449:ASN:OD1	2.16	0.45
1:A:960:TYR:CD1	1:A:961:ARG:HG3	2.52	0.44
1:A:532:GLN:NE2	7:A:1332:HOH:O	2.50	0.44
1:A:958:PRO:O	1:A:1015[A]:MET:HG2	2.17	0.44
1:A:478:GLY:HA3	1:A:486:LYS:O	2.17	0.44
1:A:177:TYR:HB3	1:A:225:LEU:HD22	2.00	0.43
1:A:772:LYS:O	1:A:776:GLU:HG3	2.18	0.43
1:A:483:GLY:O	1:A:598:HIS:CE1	2.71	0.43
1:A:196:HIS:CD2	1:A:498:SER:HB3	2.52	0.43
1:A:929:ASN:HD22	1:A:929:ASN:HA	1.58	0.43
1:A:970:ASP:O	1:A:971:GLU:HB2	2.19	0.43
1:A:561:GLU:O	1:A:562:LYS:HB2	2.17	0.43
1:A:996:MET:C	1:A:998:GLY:H	2.27	0.43
1:A:655:ILE:HD12	1:A:655:ILE:N	2.34	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:831:TYR:CD1	1:A:831:TYR:C	2.96	0.42
1:A:812:HIS:HE1	7:A:1074:HOH:O	2.02	0.42
1:A:255:LYS:HE2	7:A:1182:HOH:O	2.20	0.42
1:A:729:GLY:C	1:A:736:ALA:HB2	2.44	0.42
1:A:996:MET:C	1:A:998:GLY:N	2.77	0.42
1:A:118:ALA:O	1:A:119:SER:HB2	2.20	0.41
1:A:126:VAL:HG23	1:A:224:LEU:HD23	2.03	0.41
1:A:113:THR:HG21	1:A:578:LYS:HG3	2.02	0.41
1:A:137:HIS:HD2	1:A:139:ALA:N	2.07	0.41
1:A:411:VAL:O	1:A:411:VAL:HG22	2.19	0.41
1:A:329:ALA:HB2	1:A:487:PRO:HG3	2.03	0.41
1:A:393:ARG:HD3	1:A:418:ASN:OD1	2.20	0.41
1:A:559:GLU:OE2	1:A:559:GLU:HA	2.20	0.41
1:A:973:LEU:HD13	1:A:974:PRO:C	2.46	0.41
1:A:483:GLY:O	1:A:598:HIS:HE1	2.03	0.41
1:A:966:TYR:HB3	1:A:1008:PHE:CD2	2.56	0.41
1:A:990:PRO:HG2	1:A:1003:LEU:HD12	2.03	0.41
1:A:1028:LEU:O	1:A:1031:HIS:HD2	2.03	0.41
1:A:440:ARG:NH1	1:A:440:ARG:HG3	2.37	0.40
1:A:445:LEU:C	1:A:445:LEU:CD2	2.94	0.40
1:A:933:GLU:HB2	1:A:954:LYS:HB2	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	939/936 (100%)	898 (96%)	38 (4%)	3 (0%)	36 29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	383	ASP
1	A	382	LYS
1	A	347	ASP

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	745/786 (95%)	734 (98%)	11 (2%)	57 56

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	THR
1	A	227[A]	MET
1	A	227[B]	MET
1	A	237	ASP
1	A	302	LYS
1	A	326	LEU
1	A	471	LEU
1	A	517	LEU
1	A	544	LEU
1	A	651	GLN
1	A	929	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	HIS
1	A	258	ASN
1	A	373	ASN
1	A	419	GLN
1	A	432	GLN
1	A	494	GLN
1	A	528	GLN
1	A	592	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	598	HIS
1	A	625	GLN
1	A	651	GLN
1	A	658	ASN
1	A	679	GLN
1	A	694	GLN
1	A	756	ASN
1	A	812	HIS
1	A	920	HIS
1	A	929	ASN
1	A	982	ASN
1	A	1031	HIS

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 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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	SO4	A	6	-	4,4,4	0.40	0	6,6,6	0.10	0
4	SO4	A	2	-	4,4,4	0.41	0	6,6,6	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	A	1034	-	4,4,4	0.40	0	6,6,6	0.16	0
4	SO4	A	7	-	4,4,4	0.45	0	6,6,6	0.24	0
5	EPE	A	1035	-	15,15,15	1.03	1 (6%)	19,20,20	1.37	3 (15%)
6	MLA	A	1036	-	6,6,6	1.01	0	7,7,7	1.17	0
4	SO4	A	5	-	4,4,4	0.45	0	6,6,6	0.09	0
4	SO4	A	3	-	4,4,4	0.46	0	6,6,6	0.08	0
4	SO4	A	4	-	4,4,4	0.44	0	6,6,6	0.11	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
5	EPE	A	1035	-	-	1/9/19/19	0/1/1/1
6	MLA	A	1036	-	-	0/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1035	EPE	C10-S	2.25	1.80	1.77

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1035	EPE	O1S-S-C10	4.00	112.78	106.73
5	A	1035	EPE	O3S-S-O2S	-2.22	105.85	111.40
5	A	1035	EPE	O3S-S-O1S	-2.09	106.16	111.40

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1035	EPE	N4-C7-C8-O8

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	7	SO4	2	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	936/936 (100%)	0.03	37 (3%)	42 45	14, 24, 47, 66	5 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	383	ASP	4.2
1	A	188	ALA	4.1
1	A	996	MET	3.8
1	A	813	ALA	3.5
1	A	1032	SER	3.4
1	A	97	THR	3.4
1	A	350	ALA	3.2
1	A	737	ASN	3.1
1	A	372	ALA	3.1
1	A	362	GLU	2.9
1	A	815	GLY	2.9
1	A	347	ASP	2.8
1	A	1031	HIS	2.8
1	A	738	SER	2.8
1	A	343	VAL	2.8
1	A	384	VAL	2.8
1	A	187	THR	2.7
1	A	804	ASP	2.6
1	A	381	PHE	2.5
1	A	805	SER	2.5
1	A	502	ASN	2.5
1	A	348	HIS	2.4
1	A	376	MET	2.4
1	A	803	ASP	2.4
1	A	371	TYR	2.4
1	A	233	ASN	2.3
1	A	802	ASP	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	739	ASP	2.3
1	A	998	GLY	2.3
1	A	994	GLU	2.2
1	A	349	GLN	2.2
1	A	671	ARG	2.2
1	A	948	THR	2.2
1	A	814	ASN	2.2
1	A	377	LYS	2.1
1	A	973	LEU	2.1
1	A	999	GLY	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($\text{\AA}^2$)	Q<0.9
4	SO4	A	7	5/5	0.75	0.29	41,41,46,48	0
4	SO4	A	4	5/5	0.81	0.13	53,54,57,58	0
4	SO4	A	1034	5/5	0.86	0.12	49,49,52,52	0
4	SO4	A	6	5/5	0.87	0.09	63,64,65,65	0
6	MLA	A	1036	7/7	0.87	0.11	41,44,46,47	0
5	EPE	A	1035	15/15	0.93	0.12	32,37,47,50	0
4	SO4	A	3	5/5	0.95	0.16	34,37,39,40	0
4	SO4	A	2	5/5	0.96	0.15	44,44,45,47	0
4	SO4	A	5	5/5	0.97	0.07	35,36,37,38	0
3	NA	A	1033	1/1	0.99	0.02	18,18,18,18	0
2	CA	A	1	1/1	0.99	0.02	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.