



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:35 PM GMT

PDB ID : 1PQA
Title : Trypsin with PMSF at atomic resolution
Authors : Schmidt, A.; Jelsch, C.; Rypniewski, W.; Lamzin, V.S.
Deposited on : 2003-06-18
Resolution : 1.23 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

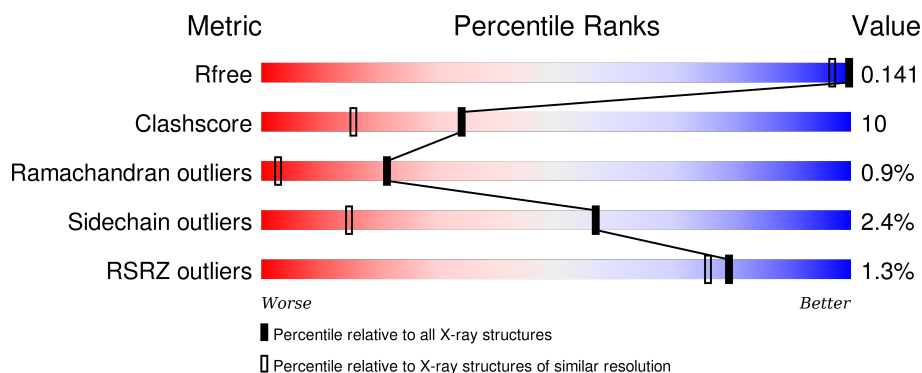
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.23 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	1229 (1.28-1.20)
Clashscore	102246	1327 (1.28-1.20)
Ramachandran outliers	100387	1274 (1.28-1.20)
Sidechain outliers	100360	1272 (1.28-1.20)
RSRZ outliers	91569	1233 (1.28-1.20)

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	224	<div> <div></div> <div>71%</div> <div>23%</div> <div>6%</div> </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
1	SEB	A	195	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	401	-	X	-	X
2	SO4	A	402	-	X	-	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3511 atoms, of which 1576 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 Trypsin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	224	Total	C	H	N	O	S	0	16	0
			3231	1010	1576	296	341	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	195	SEB	SER	MODIFIED RESIDUE	UNP P35049

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

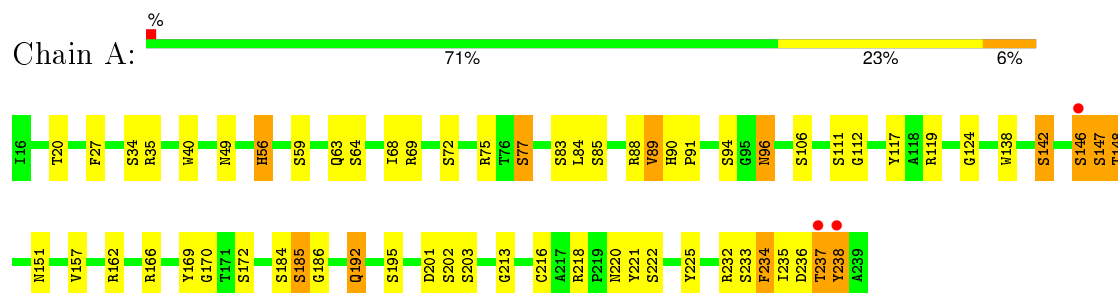
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	270	Total 270	O 270	0	0

3 Residue-property plots

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: Trypsin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	33.07Å 66.76Å 39.27Å 90.00° 108.02° 90.00°	Depositor
Resolution (Å)	25.00 – 1.23 24.89 – 1.23	Depositor EDS
% Data completeness (in resolution range)	93.8 (25.00-1.23) 90.1 (24.89-1.23)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 1.23Å)	Xtriage
Refinement program	SHELXL-96	Depositor
R, R_{free}	0.141 , (Not available) 0.145 , 0.141	Depositor DCC
R_{free} test set	2232 reflections (5.55%)	DCC
Wilson B-factor (Å ²)	11.0	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 59.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 44082 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	3511	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.71% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, SEB

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	1.86	30/1704 (1.8%)	2.24	75/2318 (3.2%)

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	233	SER	CB-OG	8.04	1.52	1.42
1	A	72	SER	CA-CB	7.44	1.64	1.52
1	A	88	ARG	CZ-NH2	6.84	1.42	1.33
1	A	96	ASN	CG-OD1	6.78	1.38	1.24
1	A	83	SER	CB-OG	6.47	1.50	1.42
1	A	34	SER	CB-OG	6.39	1.50	1.42
1	A	111	SER	CB-OG	-6.27	1.34	1.42
1	A	112	GLY	N-CA	5.91	1.54	1.46
1	A	119	ARG	CZ-NH1	5.72	1.40	1.33
1	A	184	SER	CB-OG	5.70	1.49	1.42
1	A	162	ARG	CZ-NH1	5.68	1.40	1.33
1	A	56[A]	HIS	CB-CG	-5.62	1.40	1.50
1	A	56[B]	HIS	CB-CG	-5.62	1.40	1.50
1	A	112	GLY	C-N	5.48	1.43	1.33
1	A	172	SER	CB-OG	5.48	1.49	1.42
1	A	142	SER	CB-OG	5.46	1.49	1.42
1	A	170	GLY	N-CA	5.38	1.54	1.46
1	A	186	GLY	N-CA	5.37	1.54	1.46
1	A	64	SER	CB-OG	5.35	1.49	1.42
1	A	119	ARG	CZ-NH2	5.30	1.40	1.33
1	A	20	THR	CB-OG1	5.25	1.53	1.43
1	A	56[A]	HIS	CE1-NE2	5.24	1.44	1.32
1	A	56[B]	HIS	CE1-NE2	5.24	1.44	1.32
1	A	186	GLY	C-O	5.21	1.31	1.23
1	A	185	SER	C-O	5.16	1.33	1.23
1	A	238	TYR	CG-CD1	-5.07	1.32	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	157	VAL	CB-CG1	5.06	1.63	1.52
1	A	59	SER	CB-OG	5.05	1.48	1.42
1	A	202	SER	CB-OG	5.04	1.48	1.42
1	A	222	SER	C-N	5.01	1.42	1.33

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	236	ASP	CB-CG-OD2	-23.19	97.43	118.30
1	A	236	ASP	CB-CG-OD1	20.99	137.19	118.30
1	A	119	ARG	NE-CZ-NH1	17.32	128.96	120.30
1	A	166	ARG	NE-CZ-NH2	-16.50	112.05	120.30
1	A	232	ARG	NE-CZ-NH1	14.07	127.33	120.30
1	A	119	ARG	CD-NE-CZ	13.46	142.45	123.60
1	A	238	TYR	CB-CG-CD1	13.27	128.96	121.00
1	A	88	ARG	NE-CZ-NH1	12.65	126.62	120.30
1	A	35	ARG	NE-CZ-NH1	-12.55	114.03	120.30
1	A	213	GLY	O-C-N	-11.09	104.95	122.70
1	A	238	TYR	CG-CD1-CE1	11.08	130.16	121.30
1	A	213	GLY	C-N-CA	9.28	144.91	121.70
1	A	201	ASP	CB-CG-OD2	-9.26	109.96	118.30
1	A	218	ARG	NE-CZ-NH2	9.23	124.91	120.30
1	A	192[A]	GLN	CG-CD-OE1	9.17	139.94	121.60
1	A	192[B]	GLN	CG-CD-OE1	9.17	139.94	121.60
1	A	238	TYR	CD1-CE1-CZ	-9.04	111.66	119.80
1	A	234	PHE	CD1-CE1-CZ	9.01	130.91	120.10
1	A	232	ARG	NE-CZ-NH2	-8.95	115.82	120.30
1	A	238	TYR	CA-CB-CG	-8.65	96.96	113.40
1	A	69	ARG	NE-CZ-NH2	-8.39	116.10	120.30
1	A	218	ARG	NE-CZ-NH1	-8.29	116.16	120.30
1	A	27	PHE	CB-CG-CD2	-8.16	115.09	120.80
1	A	27	PHE	CB-CG-CD1	8.14	126.50	120.80
1	A	85	SER	CA-CB-OG	8.12	133.14	111.20
1	A	40	TRP	CH2-CZ2-CE2	-7.62	109.78	117.40
1	A	88	ARG	NH1-CZ-NH2	-7.51	111.14	119.40
1	A	56[A]	HIS	CG-ND1-CE1	7.47	118.65	108.20
1	A	56[B]	HIS	CG-ND1-CE1	7.47	118.65	108.20
1	A	148[A]	THR	N-CA-CB	7.31	124.19	110.30
1	A	148[B]	THR	N-CA-CB	7.31	124.19	110.30
1	A	119	ARG	CG-CD-NE	-6.97	97.16	111.80
1	A	119	ARG	NH1-CZ-NH2	-6.94	111.77	119.40
1	A	56[A]	HIS	ND1-CE1-NE2	-6.92	94.69	109.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	56[B]	HIS	ND1-CE1-NE2	-6.92	94.69	109.90
1	A	77	SER	N-CA-CB	-6.87	100.20	110.50
1	A	35	ARG	NE-CZ-NH2	6.76	123.68	120.30
1	A	148[A]	THR	CA-CB-OG1	-6.64	95.06	109.00
1	A	148[B]	THR	CA-CB-OG1	-6.64	95.06	109.00
1	A	234	PHE	CG-CD1-CE1	-6.44	113.71	120.80
1	A	233	SER	CA-C-N	6.43	131.35	117.20
1	A	89	VAL	CA-CB-CG2	6.37	120.46	110.90
1	A	75	ARG	NE-CZ-NH2	-6.37	117.12	120.30
1	A	77	SER	CB-CA-C	6.34	122.14	110.10
1	A	237	THR	O-C-N	6.26	132.71	122.70
1	A	220	ASN	CA-C-N	6.19	130.82	117.20
1	A	89	VAL	CA-CB-CG1	-6.14	101.69	110.90
1	A	213	GLY	CA-C-N	6.12	130.66	117.20
1	A	63[A]	GLN	CB-CG-CD	6.09	127.42	111.60
1	A	63[B]	GLN	CB-CG-CD	6.09	127.42	111.60
1	A	221	TYR	CG-CD2-CE2	6.04	126.13	121.30
1	A	192[A]	GLN	OE1-CD-NE2	-5.96	108.19	121.90
1	A	192[B]	GLN	OE1-CD-NE2	-5.96	108.19	121.90
1	A	117	TYR	CB-CG-CD1	5.87	124.52	121.00
1	A	220	ASN	CA-CB-CG	-5.73	100.79	113.40
1	A	235	ILE	O-C-N	-5.68	113.61	122.70
1	A	221	TYR	CZ-CE2-CD2	-5.56	114.79	119.80
1	A	56[A]	HIS	CG-CD2-NE2	-5.52	98.71	109.20
1	A	56[B]	HIS	CG-CD2-NE2	-5.52	98.71	109.20
1	A	138	TRP	CB-CG-CD2	5.52	133.78	126.60
1	A	236	ASP	CA-CB-CG	-5.46	101.38	113.40
1	A	220	ASN	O-C-N	-5.42	114.03	122.70
1	A	203	SER	N-CA-CB	-5.40	102.39	110.50
1	A	166	ARG	NH1-CZ-NH2	5.33	125.27	119.40
1	A	225	TYR	CB-CG-CD1	5.31	124.19	121.00
1	A	185	SER	CB-CA-C	-5.28	100.06	110.10
1	A	234	PHE	CE1-CZ-CE2	-5.23	110.58	120.00
1	A	169	TYR	CG-CD2-CE2	5.22	125.48	121.30
1	A	56[A]	HIS	CE1-NE2-CD2	5.15	119.48	106.60
1	A	56[B]	HIS	CE1-NE2-CD2	5.15	119.48	106.60
1	A	49	ASN	CB-CG-ND2	-5.12	104.42	116.70
1	A	75	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	A	237	THR	CA-C-O	-5.07	109.45	120.10
1	A	35	ARG	CD-NE-CZ	5.05	130.67	123.60
1	A	216	CYS	CB-CA-C	-5.00	100.39	110.40

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	1655	1576	1577	30	0
2	A	10	0	0	0	0
3	A	270	0	0	12	0
All	All	1935	1576	1577	30	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 10.

All (30) 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:56[A]:HIS:CD2	1:A:195:SEB:HH1	1.36	1.54
1:A:56[A]:HIS:NE2	1:A:195:SEB:HH1	1.20	1.48
1:A:192[B]:GLN:NE2	1:A:195:SEB:HH2	1.44	1.31
1:A:56[A]:HIS:HD2	1:A:195:SEB:CI1	1.45	1.19
1:A:56[A]:HIS:CD2	1:A:195:SEB:HI1	1.67	1.18
1:A:192[B]:GLN:NE2	1:A:195:SEB:CH2	2.25	0.98
1:A:91[B]:PRO:HA	3:A:587:HOH:O	1.65	0.97
1:A:192[B]:GLN:HE21	1:A:195:SEB:HH2	1.15	0.88
1:A:192[B]:GLN:HE22	1:A:195:SEB:CH2	1.97	0.77
1:A:56[A]:HIS:HD2	1:A:195:SEB:CH1	1.55	0.73
1:A:89:VAL:HG23	1:A:90[B]:HIS:O	2.00	0.61
1:A:91[B]:PRO:HB3	3:A:592:HOH:O	2.01	0.60
1:A:192[B]:GLN:HE22	1:A:195:SEB:HE2	1.69	0.58
1:A:91[B]:PRO:CA	3:A:587:HOH:O	2.35	0.57
1:A:151[A]:ASN:ND2	3:A:652:HOH:O	2.39	0.54
1:A:124:GLY:HA3	3:A:670:HOH:O	2.08	0.53
1:A:90[A]:HIS:HA	1:A:234:PHE:CE1	2.44	0.52
1:A:89:VAL:HG23	3:A:587:HOH:O	2.10	0.52
1:A:192[B]:GLN:NE2	1:A:195:SEB:HE2	2.24	0.51
1:A:94[B]:SER:HA	3:A:552:HOH:O	2.11	0.51
1:A:237:THR:HG22	1:A:238:TYR:CE1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192[B]:GLN:HE22	1:A:195:SEB:CE	2.27	0.46
1:A:192[B]:GLN:HE22	1:A:195:SEB:CZ	2.29	0.45
1:A:96:ASN:ND2	3:A:629:HOH:O	2.50	0.44
1:A:148[A]:THR:HG23	3:A:551:HOH:O	2.18	0.44
1:A:147[B]:SER:HB2	3:A:608:HOH:O	2.17	0.43
1:A:56[B]:HIS:HD2	3:A:604:HOH:O	2.00	0.43
1:A:68:ILE:HD13	1:A:84:LEU:HD21	2.02	0.42
1:A:146[B]:SER:O	1:A:146[B]:SER:OG	2.30	0.42
1:A:91[B]:PRO:C	3:A:587:HOH:O	2.56	0.41

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	237/224 (106%)	227 (96%)	6 (2%)	4 (2%)	11 1

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	146[A]	SER
1	A	146[B]	SER
1	A	147[A]	SER
1	A	147[B]	SER

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	180/165 (109%)	176 (98%)	4 (2%)	60 18

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	SER
1	A	106	SER
1	A	142	SER
1	A	185	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	SEB	A	195	1	14,16,17	2.62	3 (21%)	15,21,23	3.23	2 (13%)

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	SEB	A	195	1	-	0/10/13/15	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	195	SEB	CE-CZ	3.40	1.57	1.50
1	A	195	SEB	OD1-SD	6.26	1.60	1.44
1	A	195	SEB	OD2-SD	6.41	1.61	1.44

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	195	SEB	CZ-CE-SD	-11.91	99.03	112.58
1	A	195	SEB	CB-OG-SD	3.08	126.12	119.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	195	SEB	13	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	401	-	4,4,4	3.22	4 (100%)	6,6,6	2.73	1 (16%)
2	SO4	A	402	-	4,4,4	3.54	4 (100%)	6,6,6	3.65	2 (33%)

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	SO4	A	401	-	-	0/0/0/0	0/0/0/0
2	SO4	A	402	-	-	0/0/0/0	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	SO4	O1-S	2.88	1.57	1.47
2	A	401	SO4	O3-S	3.17	1.58	1.47
2	A	401	SO4	O2-S	3.22	1.58	1.47
2	A	402	SO4	O1-S	3.28	1.58	1.47
2	A	402	SO4	O4-S	3.42	1.59	1.47
2	A	402	SO4	O3-S	3.57	1.60	1.47
2	A	401	SO4	O4-S	3.57	1.60	1.47
2	A	402	SO4	O2-S	3.87	1.60	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	402	SO4	O2-S-O1	-3.37	98.83	109.50
2	A	401	SO4	O4-S-O3	6.39	134.97	108.98
2	A	402	SO4	O4-S-O3	8.07	141.79	108.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	223/224 (99%)	-0.30	3 (1%) 79 76	7, 14, 26, 39	2 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	146[A]	SER	3.6
1	A	238	TYR	2.6
1	A	237	THR	2.1

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
1	SEB	A	195	16/17	0.96	0.09	-	7,17,23,29	7

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
2	SO4	A	401	5/5	0.95	0.13	2.62	25,28,40,41	5
2	SO4	A	402	5/5	0.94	0.14	1.23	20,21,30,36	5

6.5 Other polymers [i](#)

There are no such residues in this entry.