



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:28 PM UTC

PDB ID : 1B8J / pdb_00001b8j
Title : ALKALINE PHOSPHATASE COMPLEXED WITH VANADATE
Authors : Holtz, K.M.; Stec, B.; Kantrowitz, E.R.
Deposited on : 1999-02-01
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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

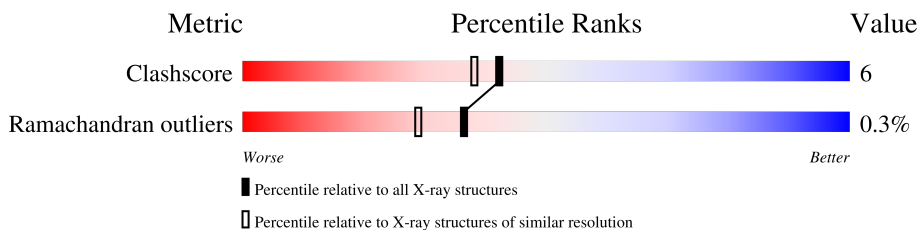
1 Overall quality at a glance

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))
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	449	 81% 18% .
1	B	449	 81% 18% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7579 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 PROTEIN (ALKALINE PHOSPHATASE).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	449	Total	C	N	O	S	V	0	0	0
			3309	2042	581	673	12	1			
1	B	449	Total	C	N	O	S	V	0	0	0
			3309	2042	581	673	12	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	SVA	SER	modified residue	UNP P00634
B	102	SVA	SER	modified residue	UNP P00634

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is water.

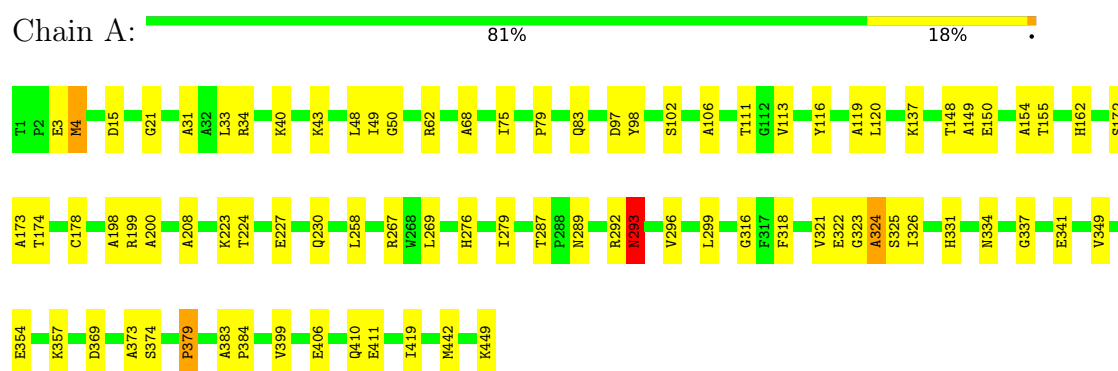
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	505	Total	O	0	0
			505	505		
5	B	440	Total	O	0	0
			440	440		

3 Residue-property plots

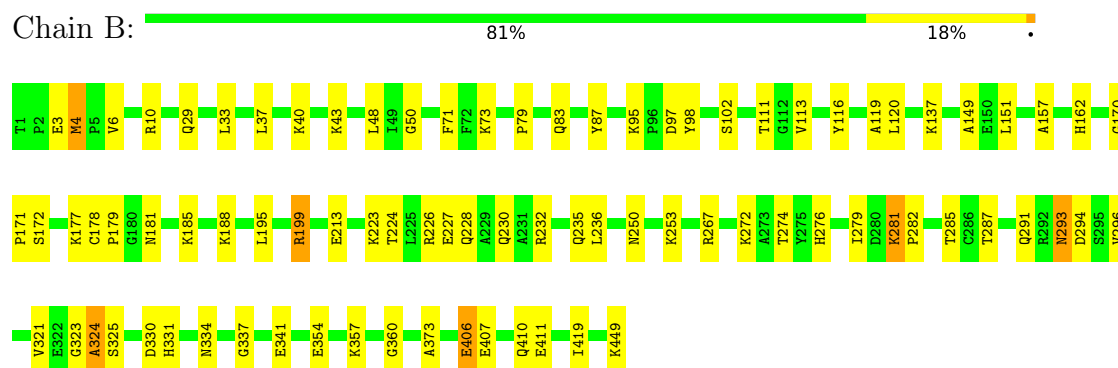
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

• Molecule 1: PROTEIN (ALKALINE PHOSPHATASE)



• Molecule 1: PROTEIN (ALKALINE PHOSPHATASE)



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	195.41 Å 167.81 Å 76.52 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 1.90	Depositor
% Data completeness (in resolution range)	98.4 (8.00-1.90)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.177 , 0.196	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7579	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	1.17	7/3352 (0.2%)	1.31	31/4549 (0.7%)
1	B	1.04	1/3352 (0.0%)	1.28	30/4549 (0.7%)
All	All	1.11	8/6704 (0.1%)	1.30	61/9098 (0.7%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	442	MET	SD-CE	-8.36	1.58	1.79
1	A	198	ALA	CA-CB	6.71	1.64	1.53
1	B	4	MET	SD-CE	6.55	1.96	1.79
1	A	31	ALA	CA-CB	-6.46	1.43	1.53
1	A	419	ILE	CA-CB	5.94	1.62	1.54
1	A	155	THR	CA-CB	5.82	1.61	1.53
1	A	106	ALA	CA-CB	-5.52	1.44	1.53
1	A	4	MET	SD-CE	5.44	1.93	1.79

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	293	ASN	N-CA-C	10.17	132.46	110.80
1	B	293	ASN	N-CA-C	9.90	131.90	110.80
1	A	324	ALA	N-CA-C	9.67	121.90	111.36
1	B	324	ALA	N-CA-C	9.59	121.81	111.36
1	B	170	GLY	CA-C-N	7.95	127.51	119.24
1	B	170	GLY	C-N-CA	7.95	127.51	119.24
1	B	325	SER	N-CA-C	7.77	122.60	113.20
1	A	325	SER	N-CA-C	7.47	122.24	113.20
1	A	79	PRO	N-CA-C	7.31	124.34	114.18
1	A	373	ALA	N-CA-C	7.18	120.53	111.69
1	B	172	SER	N-CA-C	7.16	119.16	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	411	GLU	N-CA-C	6.95	121.19	110.42
1	A	98	TYR	N-CA-C	6.87	120.76	112.38
1	B	373	ALA	N-CA-C	6.60	121.63	112.45
1	A	154	ALA	N-CA-C	6.59	119.31	111.33
1	A	419	ILE	N-CA-C	-6.50	98.67	108.17
1	B	79	PRO	N-CA-C	6.44	123.13	114.18
1	B	330	ASP	N-CA-C	-6.34	104.45	111.36
1	B	98	TYR	N-CA-C	6.14	119.87	112.38
1	A	323	GLY	N-CA-C	-6.12	98.67	113.18
1	A	172	SER	N-CA-C	6.08	117.91	111.28
1	B	97	ASP	N-CA-C	-6.01	97.34	107.99
1	A	4	MET	CA-C-N	5.85	126.34	120.14
1	A	4	MET	C-N-CA	5.85	126.34	120.14
1	A	318	PHE	N-CA-C	-5.76	100.01	109.40
1	A	173	ALA	N-CA-C	-5.73	105.17	111.82
1	B	419	ILE	N-CA-C	-5.67	100.02	108.46
1	B	285	THR	N-CA-C	-5.62	99.75	108.90
1	A	50	GLY	N-CA-C	-5.59	100.32	111.34
1	B	323	GLY	N-CA-C	-5.47	100.23	113.18
1	B	360	GLY	N-CA-C	5.46	122.61	115.40
1	B	411	GLU	N-CA-C	5.45	118.86	110.42
1	B	334	ASN	CA-C-N	5.41	124.86	119.24
1	B	334	ASN	C-N-CA	5.41	124.86	119.24
1	A	49	ILE	CA-C-N	-5.40	117.48	122.29
1	A	49	ILE	C-N-CA	-5.40	117.48	122.29
1	A	150	GLU	N-CA-C	-5.39	102.31	110.28
1	A	379	PRO	N-CA-C	5.36	120.88	113.65
1	B	185	LYS	N-CA-C	-5.29	106.05	112.88
1	B	406	GLU	N-CA-C	-5.29	106.83	113.28
1	B	235	GLN	N-CA-C	-5.24	99.97	108.41
1	A	326	ILE	N-CA-C	-5.18	105.44	110.42
1	B	50	GLY	N-CA-C	-5.17	100.92	113.18
1	A	200	ALA	N-CA-C	-5.17	103.71	110.53
1	A	97	ASP	N-CA-C	-5.15	98.87	107.99
1	A	374	SER	N-CA-C	5.15	117.81	110.24
1	A	75	ILE	N-CA-C	5.15	117.48	111.05
1	A	322	GLU	N-CA-C	5.15	116.80	108.41
1	B	6	VAL	N-CA-C	-5.14	101.08	108.53
1	B	281	LYS	CA-C-N	5.13	125.11	120.03
1	B	281	LYS	C-N-CA	5.13	125.11	120.03
1	A	334	ASN	CA-C-N	5.13	124.57	119.24
1	A	334	ASN	C-N-CA	5.13	124.57	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	199	ARG	N-CA-C	5.13	118.20	111.28
1	B	95	LYS	CA-C-N	-5.11	114.49	119.76
1	B	95	LYS	C-N-CA	-5.11	114.49	119.76
1	A	406	GLU	N-CA-C	-5.10	107.06	113.28
1	A	43	LYS	N-CA-C	-5.09	105.42	111.69
1	B	43	LYS	N-CA-C	-5.09	106.23	112.90
1	B	267	ARG	N-CA-C	5.09	116.91	111.36
1	A	316	GLY	N-CA-C	-5.08	105.75	111.85

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	3309	0	3248	39	0
1	B	3309	0	3248	49	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	505	0	0	3	1
5	B	440	0	0	15	0
All	All	7579	0	6496	83	1

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 (83) 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:294:ASP:HB3	5:B:1310:HOH:O	1.77	0.83
1:A:34:ARG:HD3	5:B:1231:HOH:O	1.79	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:ARG:HB2	5:A:1077:HOH:O	1.83	0.79
1:B:48:LEU:HD13	1:B:321:VAL:HB	1.66	0.77
1:B:195:LEU:HD11	5:B:1059:HOH:O	1.83	0.77
1:B:224:THR:OG1	1:B:227:GLU:HG3	1.84	0.76
1:B:151:LEU:HD13	5:B:1059:HOH:O	1.85	0.76
1:B:73:LYS:NZ	5:B:1408:HOH:O	2.23	0.71
1:A:137:LYS:HE3	1:A:199:ARG:O	1.93	0.68
1:A:48:LEU:HD13	1:A:321:VAL:HB	1.78	0.65
1:A:224:THR:OG1	1:A:227:GLU:HG3	1.96	0.65
1:A:293:ASN:HB3	1:A:296:VAL:HG23	1.80	0.64
1:A:111:THR:OG1	1:A:113:VAL:HG12	1.98	0.62
1:A:287:THR:HG22	5:A:1423:HOH:O	1.99	0.62
1:B:195:LEU:HD12	5:B:1049:HOH:O	1.99	0.62
1:B:287:THR:HG22	5:B:1373:HOH:O	1.98	0.62
1:B:73:LYS:HE3	5:B:1194:HOH:O	2.00	0.61
1:B:137:LYS:HE3	1:B:199:ARG:O	2.00	0.61
1:B:272:LYS:HE2	5:B:1012:HOH:O	2.01	0.59
1:B:188:LYS:HE3	5:B:1179:HOH:O	2.01	0.59
1:B:293:ASN:HB3	1:B:296:VAL:HG23	1.84	0.59
1:A:267:ARG:HG2	1:A:292:ARG:NH1	2.19	0.57
1:A:83:GLN:HE21	1:B:83:GLN:HE21	1.55	0.55
1:A:116:TYR:CZ	1:A:119:ALA:HB2	2.41	0.55
1:B:291:GLN:HG3	5:B:1322:HOH:O	2.09	0.53
1:A:379:PRO:HA	1:A:399:VAL:HG21	1.90	0.53
1:A:208:ALA:HB2	1:A:258:LEU:HB3	1.91	0.53
1:B:111:THR:OG1	1:B:113:VAL:HG12	2.10	0.52
1:A:34:ARG:HG2	1:B:37:LEU:HD12	1.92	0.51
1:A:120:LEU:O	1:A:162:HIS:HA	2.09	0.51
1:B:10:ARG:HB2	1:B:71:PHE:CE1	2.46	0.51
1:A:276:HIS:O	1:A:279:ILE:HG12	2.12	0.50
1:B:116:TYR:CZ	1:B:119:ALA:HB2	2.47	0.49
1:B:120:LEU:O	1:B:162:HIS:HA	2.12	0.49
1:A:174:THR:HG23	1:A:178:CYS:HB2	1.95	0.49
1:B:228:GLN:O	1:B:232:ARG:HG3	2.13	0.49
1:A:354:GLU:HA	1:A:357:LYS:HE2	1.95	0.48
1:A:269:LEU:HD22	5:A:1002:HOH:O	2.13	0.48
1:B:354:GLU:HA	1:B:357:LYS:HE2	1.96	0.47
1:B:178:CYS:HB3	1:B:181:ASN:OD1	2.14	0.47
1:A:369:ASP:OD1	1:A:369:ASP:N	2.47	0.47
1:B:10:ARG:HB2	1:B:71:PHE:CD1	2.50	0.47
1:B:274:THR:HG21	5:B:1088:HOH:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:ALA:HB2	1:B:324:ALA:CB	2.45	0.46
1:B:281:LYS:HB3	1:B:282:PRO:HD2	1.98	0.46
1:B:226:ARG:O	1:B:230:GLN:HG2	2.16	0.46
1:A:449:LYS:HA	1:A:449:LYS:HD3	1.73	0.46
1:A:33:LEU:HD12	1:A:33:LEU:HA	1.76	0.45
1:B:230:GLN:HG2	1:B:230:GLN:H	1.59	0.45
1:B:276:HIS:O	1:B:279:ILE:HG12	2.17	0.45
1:A:149:ALA:HB2	1:A:324:ALA:CB	2.47	0.44
1:B:331:HIS:ND1	1:B:410:GLN:O	2.47	0.44
1:A:40:LYS:HA	1:A:40:LYS:HD2	1.61	0.44
1:B:40:LYS:HA	1:B:40:LYS:HD2	1.59	0.44
1:A:3:GLU:HG2	1:A:4:MET:N	2.33	0.44
1:A:223:LYS:HB3	1:A:223:LYS:HE2	1.61	0.44
1:A:289:ASN:O	1:A:292:ARG:HD3	2.17	0.44
1:B:29:GLN:HG3	5:B:1340:HOH:O	2.17	0.44
1:A:331:HIS:ND1	1:A:410:GLN:O	2.48	0.44
1:B:3:GLU:HG2	1:B:4:MET:N	2.31	0.44
1:A:48:LEU:HG	1:A:349:VAL:HG22	2.00	0.43
1:B:33:LEU:HD12	1:B:33:LEU:HA	1.81	0.43
1:A:383:ALA:C	1:B:406:GLU:HG3	2.43	0.43
1:A:379:PRO:HA	1:A:399:VAL:CG2	2.48	0.42
1:B:291:GLN:NE2	5:B:1246:HOH:O	2.52	0.42
1:B:250:ASN:OD1	1:B:253:LYS:HD2	2.19	0.42
1:A:148:THR:HG23	1:A:299:LEU:HD13	2.00	0.42
1:A:15:ASP:O	1:A:21:GLY:HA3	2.19	0.42
1:B:48:LEU:CD1	1:B:321:VAL:HB	2.44	0.42
1:B:171:PRO:HD2	1:B:213:GLU:OE1	2.20	0.42
1:A:276:HIS:HE1	1:B:406:GLU:OE1	2.02	0.41
1:B:177:LYS:C	1:B:179:PRO:HD3	2.44	0.41
1:A:383:ALA:HB1	1:A:384:PRO:HD2	2.02	0.41
1:B:149:ALA:HB2	1:B:324:ALA:HB1	2.01	0.41
1:B:337:GLY:O	1:B:341:GLU:HG2	2.20	0.41
1:A:68:ALA:HB1	1:B:87:TYR:CD1	2.55	0.41
1:B:223:LYS:HE2	1:B:223:LYS:HB3	1.63	0.40
1:B:226:ARG:CZ	1:B:236:LEU:HD23	2.51	0.40
1:A:337:GLY:O	1:A:341:GLU:HG2	2.21	0.40
1:B:157:ALA:HA	5:B:1049:HOH:O	2.20	0.40
1:B:449:LYS:HA	1:B:449:LYS:HD3	1.86	0.40
1:A:230:GLN:H	1:A:230:GLN:HG2	1.56	0.40
1:A:62:ARG:HH11	1:A:62:ARG:HD2	1.72	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1198:HOH:O	5:A:1198:HOH:O[3_656]	1.68	0.52

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	356/449 (79%)	349 (98%)	6 (2%)	1 (0%)	36	29
1	B	389/449 (87%)	381 (98%)	7 (2%)	1 (0%)	36	29
All	All	745/898 (83%)	730 (98%)	13 (2%)	2 (0%)	36	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	293	ASN
1	B	407	GLU

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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
1	SVA	B	102	2,1	5,10,11	2.04	2 (40%)	0,16,18	-	-
1	SVA	A	102	2,1	5,10,11	1.44	1 (20%)	0,16,18	-	-

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	SVA	B	102	2,1	-	0/1/9/11	-
1	SVA	A	102	2,1	-	0/1/9/11	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	102	SVA	OG-CB	-3.21	1.34	1.43
1	B	102	SVA	OG-V	-2.60	1.70	1.80
1	A	102	SVA	OG-V	-2.04	1.72	1.80

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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	$\# Z > 2$
3	SO4	B	971	-	4,4,4	0.98	0	6,6,6	0.60	0
3	SO4	A	970	-	4,4,4	0.81	0	6,6,6	0.21	0

There are no bond length outliers.

There are no bond angle outliers.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.