



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 10:42 AM UTC

PDB ID : 4Y3K / pdb_00004y3k
Title : Structure of Vaspin mutant E379S
Authors : Pippel, J.; Strater, N.; Ulbricht, D.; Schultz, S.; Meier, R.; Heiker, J.T.
Deposited on : 2015-02-10
Resolution : 2.20 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

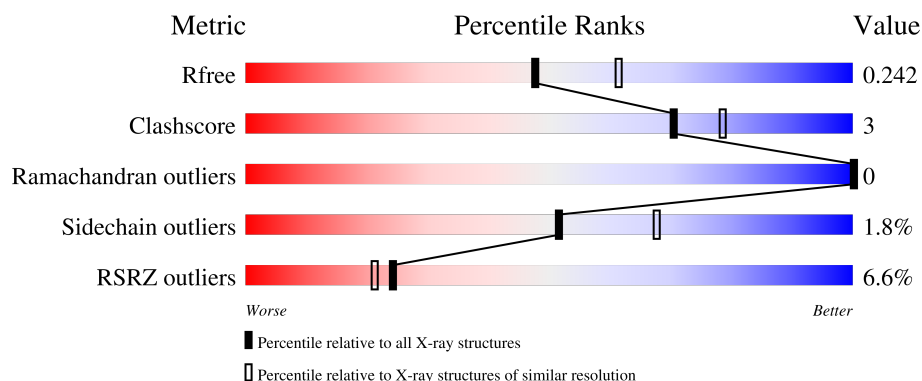
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 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	414	
1	B	414	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6224 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 Serpin A12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	0	6	0
			3008	1942	508	544	14			
1	B	361	Total	C	N	O	S	0	5	0
			2975	1927	500	535	13			

There are 44 discrepancies between the modelled and reference sequences:

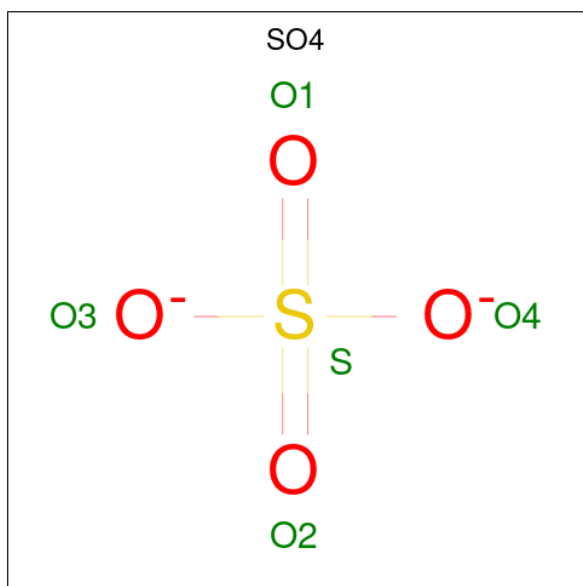
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q8IW75
A	2	HIS	-	expression tag	UNP Q8IW75
A	3	HIS	-	expression tag	UNP Q8IW75
A	4	HIS	-	expression tag	UNP Q8IW75
A	5	HIS	-	expression tag	UNP Q8IW75
A	6	HIS	-	expression tag	UNP Q8IW75
A	7	HIS	-	expression tag	UNP Q8IW75
A	8	HIS	-	expression tag	UNP Q8IW75
A	9	HIS	-	expression tag	UNP Q8IW75
A	10	HIS	-	expression tag	UNP Q8IW75
A	11	HIS	-	expression tag	UNP Q8IW75
A	12	SER	-	expression tag	UNP Q8IW75
A	13	SER	-	expression tag	UNP Q8IW75
A	14	GLY	-	expression tag	UNP Q8IW75
A	15	HIS	-	expression tag	UNP Q8IW75
A	16	ILE	-	expression tag	UNP Q8IW75
A	17	GLU	-	expression tag	UNP Q8IW75
A	18	GLY	-	expression tag	UNP Q8IW75
A	19	ARG	-	expression tag	UNP Q8IW75
A	20	HIS	-	expression tag	UNP Q8IW75
A	21	MET	-	expression tag	UNP Q8IW75
A	379	SER	GLU	engineered mutation	UNP Q8IW75
B	1	GLY	-	expression tag	UNP Q8IW75
B	2	HIS	-	expression tag	UNP Q8IW75
B	3	HIS	-	expression tag	UNP Q8IW75

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Chain	Residue	Modelled	Actual	Comment	Reference
B	4	HIS	-	expression tag	UNP Q8IW75
B	5	HIS	-	expression tag	UNP Q8IW75
B	6	HIS	-	expression tag	UNP Q8IW75
B	7	HIS	-	expression tag	UNP Q8IW75
B	8	HIS	-	expression tag	UNP Q8IW75
B	9	HIS	-	expression tag	UNP Q8IW75
B	10	HIS	-	expression tag	UNP Q8IW75
B	11	HIS	-	expression tag	UNP Q8IW75
B	12	SER	-	expression tag	UNP Q8IW75
B	13	SER	-	expression tag	UNP Q8IW75
B	14	GLY	-	expression tag	UNP Q8IW75
B	15	HIS	-	expression tag	UNP Q8IW75
B	16	ILE	-	expression tag	UNP Q8IW75
B	17	GLU	-	expression tag	UNP Q8IW75
B	18	GLY	-	expression tag	UNP Q8IW75
B	19	ARG	-	expression tag	UNP Q8IW75
B	20	HIS	-	expression tag	UNP Q8IW75
B	21	MET	-	expression tag	UNP Q8IW75
B	379	SER	GLU	engineered mutation	UNP Q8IW75

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



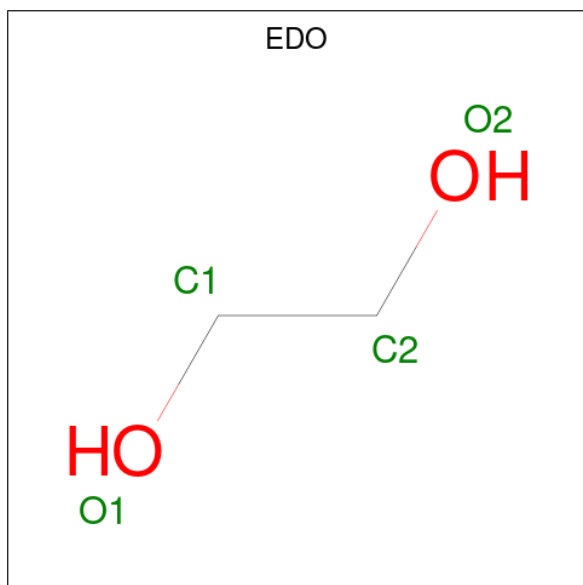
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	134	Total 134	O 134	0	0
4	B	54	Total 54	O 54	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	132.96Å 151.35Å 61.60Å 90.00° 98.43° 90.00°	Depositor
Resolution (Å)	48.37 – 2.20 48.37 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.37-2.20) 99.8 (48.37-2.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.20Å)	Xtriage
Refinement program	BUSTER 2.10.1	Depositor
R, R_{free}	0.186 , 0.219 0.206 , 0.242	Depositor DCC
R_{free} test set	3083 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 73.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6224	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.31% 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

5.1 Standard geometry

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

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	0.89	0/3093	1.30	12/4158 (0.3%)
1	B	0.80	0/3052	1.27	9/4103 (0.2%)
All	All	0.85	0/6145	1.29	21/8261 (0.3%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	52	ASN	CA-CB-CG	7.75	120.35	112.60
1	B	83	PHE	CA-CB-CG	-6.86	106.94	113.80
1	B	104	PHE	CA-CB-CG	6.33	120.13	113.80
1	B	104	PHE	CA-C-N	6.26	128.96	120.38
1	B	104	PHE	C-N-CA	6.26	128.96	120.38
1	A	104	PHE	CA-CB-CG	6.24	120.04	113.80
1	B	35	GLU	CA-C-N	6.06	129.08	120.53
1	B	35	GLU	C-N-CA	6.06	129.08	120.53
1	A	290[A]	ASP	CA-C-N	5.58	128.02	120.38
1	A	290[A]	ASP	C-N-CA	5.58	128.02	120.38
1	A	290[B]	ASP	CA-C-N	5.58	128.02	120.38
1	A	290[B]	ASP	C-N-CA	5.58	128.02	120.38
1	A	127	THR	CA-C-N	5.54	127.96	120.38
1	A	127	THR	C-N-CA	5.54	127.96	120.38
1	A	52	ASN	CA-CB-CG	5.38	117.98	112.60
1	A	394	ILE	N-CA-CB	5.25	117.97	111.41
1	B	147	GLN	CA-C-N	5.14	127.49	120.54
1	B	147	GLN	C-N-CA	5.14	127.49	120.54
1	A	281	LYS	CA-C-N	5.05	127.30	120.38
1	A	281	LYS	C-N-CA	5.05	127.30	120.38
1	A	83	PHE	CA-CB-CG	-5.05	108.75	113.80

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	3008	0	3083	21	0
1	B	2975	0	3049	20	0
2	A	15	0	0	0	0
2	B	10	0	0	0	0
3	A	20	0	30	0	0
3	B	8	0	12	3	0
4	A	134	0	0	0	0
4	B	54	0	0	0	0
All	All	6224	0	6174	41	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 3.

All (41) 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:335:HIS:HB2	3:B:501:EDO:H12	1.76	0.66
1:B:82:ALA:HB1	1:B:204:LEU:HD21	1.78	0.65
1:B:223:THR:HG22	1:B:241:MET:HA	1.82	0.62
1:A:49:ALA:O	1:A:53[A]:MET:HG3	2.00	0.61
1:B:337:ASP:HB2	3:B:501:EDO:H11	1.82	0.61
1:B:151:LEU:HD23	1:B:162:THR:HB	1.83	0.61
1:A:131:LYS:HB2	1:A:211:ARG:HB2	1.84	0.59
1:A:294:ARG:O	1:A:297:THR:HG22	2.04	0.57
1:B:71:ASN:H	1:B:414:LYS:HG2	1.72	0.55
1:A:77:LEU:HD11	1:A:119:ILE:HG21	1.90	0.54
1:A:395:TYR:OH	1:A:400:PRO:HB2	2.09	0.52
1:B:135:GLY:HA3	1:B:207[B]:TYR:CE1	2.45	0.51
1:A:42:ARG:HG2	1:A:118:TYR:CZ	2.46	0.50
1:A:334:GLU:HG2	1:A:346:SER:O	2.13	0.48
1:B:215:LYS:HD2	1:B:265:GLN:HA	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:334:GLU:HG3	1:B:346:SER:HA	1.97	0.46
1:B:70:ARG:HD2	1:B:414:LYS:HD2	1.97	0.46
1:B:142:GLN:HA	1:B:164:LEU:HD22	1.96	0.46
1:A:303:VAL:HG12	1:A:381:PRO:HB3	1.98	0.46
1:A:146:PRO:HB2	1:A:151:LEU:HD11	1.98	0.45
1:B:293:SER:HA	1:B:296:LYS:HD2	1.97	0.45
1:B:308:VAL:HG12	1:B:386:ILE:HG13	1.98	0.45
1:B:177:ILE:O	1:B:181:ILE:HD12	2.17	0.45
1:B:304:VAL:HG13	1:B:383:VAL:HA	1.99	0.45
1:A:287:LEU:HD23	1:A:291:THR:HG21	1.99	0.45
1:B:99:LYS:HA	1:B:104:PHE:HB2	2.00	0.43
1:B:71:ASN:HD22	1:B:414:LYS:HB2	1.83	0.43
1:A:49:ALA:HB1	1:A:400:PRO:O	2.19	0.42
1:A:82:ALA:HB1	1:A:204:LEU:HD21	2.01	0.42
1:B:89:GLY:HA3	1:B:341:ILE:HG13	2.01	0.42
1:A:73:PHE:O	1:A:315:GLY:HA3	2.19	0.42
1:A:38:GLY:O	1:A:42:ARG:HG3	2.18	0.42
1:A:395:TYR:CZ	1:A:400:PRO:HB2	2.55	0.42
1:A:50:ARG:HA	1:A:53[A]:MET:HE3	2.01	0.41
1:A:137:THR:HG23	1:A:161:GLU:HG2	2.02	0.41
1:A:252:ASP:HB2	1:A:259:ILE:HD11	2.02	0.41
1:A:61:LYS:HE3	1:A:285:LYS:HG2	2.01	0.41
1:B:71:ASN:ND2	1:B:414:LYS:HB2	2.35	0.41
1:B:333:GLU:O	3:B:501:EDO:H21	2.21	0.41
1:A:133:SER:HB2	1:A:209:PHE:CZ	2.56	0.40
1:A:71:ASN:HB2	1:A:414:LYS:HG3	2.03	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	369/414 (89%)	356 (96%)	13 (4%)	0	100	100
1	B	360/414 (87%)	343 (95%)	17 (5%)	0	100	100
All	All	729/828 (88%)	699 (96%)	30 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	337/369 (91%)	333 (99%)	4 (1%)	63	78
1	B	332/369 (90%)	324 (98%)	8 (2%)	43	58
All	All	669/738 (91%)	657 (98%)	12 (2%)	51	68

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	SER
1	A	161	GLU
1	A	308	VAL
1	A	393	LEU
1	B	72	ILE
1	B	161	GLU
1	B	181	ILE
1	B	280	LEU
1	B	287	LEU
1	B	321	LYS
1	B	391	LEU
1	B	412	ILE

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

Mol	Chain	Res	Type
1	A	67	ASN

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Mol	Chain	Res	Type
1	A	174	GLN
1	A	183	GLN
1	A	191	ASN
1	A	195	ASN
1	A	216	HIS
1	A	233	ASN
1	A	265	GLN
1	A	288	GLN
1	B	51	GLN
1	B	117	HIS
1	B	121	HIS
1	B	267	ASN

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](#)

12 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	507	-	4,4,4	0.44	0	6,6,6	0.25	0
3	EDO	A	506	-	3,3,3	0.49	0	2,2,2	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	B	504	-	3,3,3	0.46	0	2,2,2	0.33	0
2	SO4	A	501	-	4,4,4	0.22	0	6,6,6	0.33	0
3	EDO	A	505	-	3,3,3	0.48	0	2,2,2	0.38	0
2	SO4	A	508	-	4,4,4	0.30	0	6,6,6	0.17	0
3	EDO	A	503	-	3,3,3	0.53	0	2,2,2	0.38	0
3	EDO	A	502	-	3,3,3	0.69	0	2,2,2	0.14	0
3	EDO	B	501	-	3,3,3	0.33	0	2,2,2	0.39	0
2	SO4	B	502	-	4,4,4	0.29	0	6,6,6	0.14	0
3	EDO	A	504	-	3,3,3	0.45	0	2,2,2	0.72	0
2	SO4	B	503	-	4,4,4	0.14	0	6,6,6	0.31	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
3	EDO	A	506	-	-	0/1/1/1	-
3	EDO	B	504	-	-	1/1/1/1	-
3	EDO	A	505	-	-	0/1/1/1	-
3	EDO	A	503	-	-	1/1/1/1	-
3	EDO	A	502	-	-	0/1/1/1	-
3	EDO	B	501	-	-	0/1/1/1	-
3	EDO	A	504	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	504	EDO	O1-C1-C2-O2
3	A	503	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	501	EDO	3	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	367/414 (88%)	-0.06	8 (2%) 62 59	31, 52, 81, 118	6 (1%)
1	B	361/414 (87%)	0.65	40 (11%) 10 8	27, 78, 146, 184	5 (1%)
All	All	728/828 (87%)	0.29	48 (6%) 24 21	27, 61, 138, 184	11 (1%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	384	VAL	4.2
1	B	365	THR	4.0
1	B	280	LEU	3.8
1	A	399	ILE	3.7
1	B	304	VAL	3.6
1	B	395	TYR	3.5
1	A	367	GLY	3.2
1	B	251	TYR	3.0
1	A	106	LYS	3.0
1	B	306	VAL	2.9
1	A	36	VAL	2.9
1	B	247	TYR	2.8
1	B	299	LEU	2.7
1	B	209[A]	PHE	2.7
1	B	274	LEU	2.7
1	B	236	VAL	2.7
1	B	303	VAL	2.7
1	A	400	PRO	2.7
1	B	36	VAL	2.6
1	B	297	THR	2.6
1	B	228	PHE	2.6
1	B	383	VAL	2.4
1	B	223	THR	2.4
1	A	344[A]	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	398	LYS	2.4
1	B	232	LYS	2.4
1	B	295	TRP	2.3
1	B	233	ASN	2.3
1	B	293	SER	2.3
1	B	216	HIS	2.3
1	B	413	GLY	2.3
1	B	399	ILE	2.3
1	B	37	GLN	2.2
1	B	256	SER	2.2
1	B	246	ILE	2.2
1	B	287	LEU	2.2
1	B	234	SER	2.2
1	B	255	LEU	2.2
1	A	414	LYS	2.1
1	B	296	LYS	2.1
1	B	298	LEU	2.1
1	B	242	PHE	2.1
1	A	233	ASN	2.1
1	B	292	PHE	2.1
1	B	310	ARG	2.1
1	B	41	GLN	2.0
1	B	51	GLN	2.0
1	B	309	PRO	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
3	EDO	B	504	4/4	0.78	0.20	55,56,56,57	4
3	EDO	A	503	4/4	0.86	0.15	45,45,46,49	4
2	SO4	A	507	5/5	0.87	0.13	56,59,61,64	5
3	EDO	A	506	4/4	0.87	0.17	52,56,60,62	0
3	EDO	A	502	4/4	0.87	0.25	56,61,63,64	0
2	SO4	B	503	5/5	0.89	0.17	49,53,55,56	5
2	SO4	B	502	5/5	0.91	0.09	71,71,72,73	5
2	SO4	A	508	5/5	0.91	0.14	71,71,72,73	5
2	SO4	A	501	5/5	0.92	0.12	58,62,64,68	5
3	EDO	A	505	4/4	0.92	0.13	63,63,63,64	0
3	EDO	B	501	4/4	0.94	0.25	66,66,67,70	0
3	EDO	A	504	4/4	0.98	0.07	35,36,39,40	0

6.5 Other polymers ⓘ

There are no such residues in this entry.