



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2026 – 01:44 PM UTC

PDB ID : 8DY2 / pdb_00008dy2
Title : Crystal Structure of spFv GLK1
Authors : Luo, J.; Boucher, L.E.
Deposited on : 2022-08-03
Resolution : 1.65 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

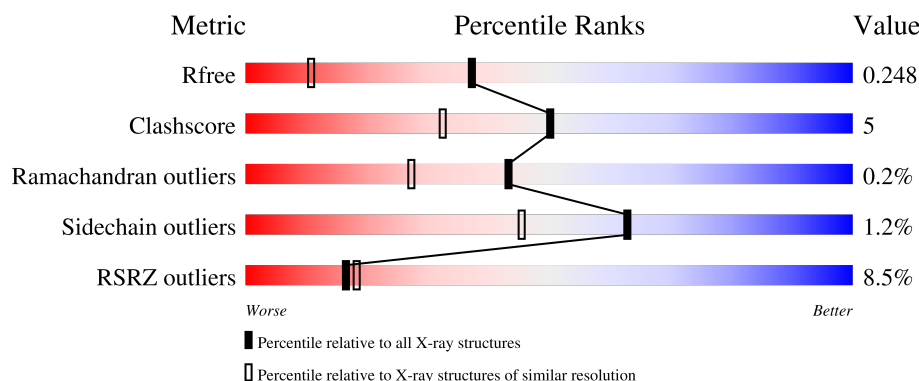
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	252	6% 88% 8%
1	B	252	4% 86% 11%
1	C	252	20% 82% 14%
1	D	252	3% 85% 12%

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
2	SO4	A	301	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7592 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 spFv GLK1 LH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1768	1104	294	359	11			
1	B	245	Total	C	N	O	S	0	0	0
			1781	1110	297	363	11			
1	C	244	Total	C	N	O	S	0	0	0
			1782	1111	297	363	11			
1	D	243	Total	C	N	O	S	0	0	0
			1783	1112	299	361	11			

- 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		

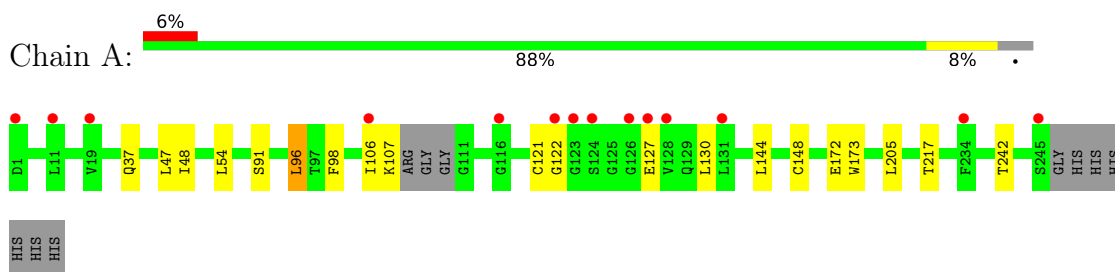
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	137	Total 137	O 137	0	0
3	B	126	Total 126	O 126	0	0
3	C	87	Total 87	O 87	0	0
3	D	123	Total 123	O 123	0	0

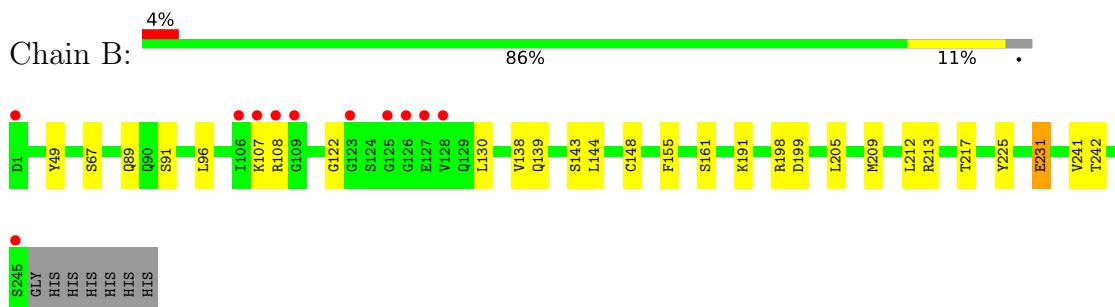
3 Residue-property plots [i](#)

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.

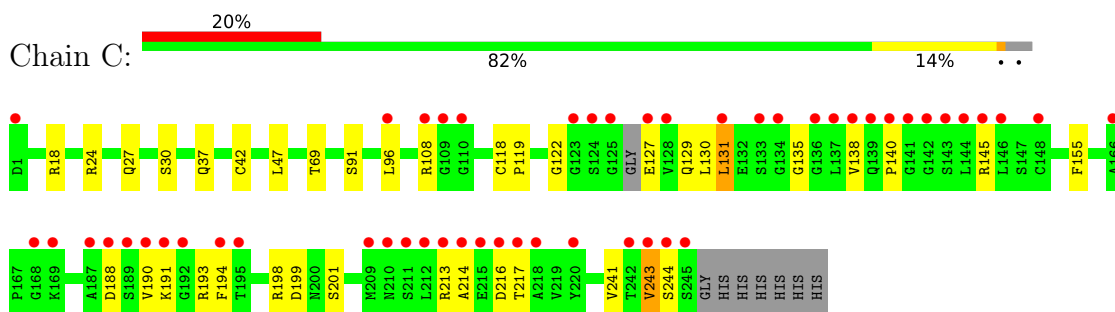
• Molecule 1: spFv GLK1 LH



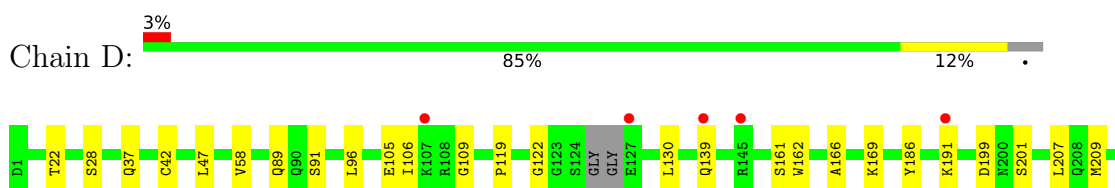
• Molecule 1: spFv GLK1 LH

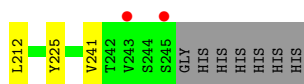


• Molecule 1: spFv GLK1 LH



• Molecule 1: spFv GLK1 LH





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.63Å 78.94Å 257.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.86 – 1.65 35.86 – 1.65	Depositor EDS
% Data completeness (in resolution range)	47.8 (35.86-1.65) 44.8 (35.86-1.65)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 1.65Å)	Xtriage
Refinement program	PHENIX 1.18_3855	Depositor
R, R_{free}	0.215 , 0.246 0.217 , 0.248	Depositor DCC
R_{free} test set	2675 reflections (2.36%)	wwPDB-VP
Wilson B-factor (Å ²)	20.5	Xtriage
Anisotropy	0.145	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 30.7	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	0.93	EDS
Total number of atoms	7592	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.10% 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: 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	0.12	0/1806	0.38	0/2448
1	B	0.14	0/1820	0.36	0/2469
1	C	0.14	0/1820	0.40	1/2466 (0.0%)
1	D	0.14	0/1821	0.37	0/2467
All	All	0.14	0/7267	0.37	1/9850 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	30	SER	CB-CA-C	-5.67	109.52	117.23

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1768	0	1675	14	0
1	B	1781	0	1679	17	0
1	C	1782	0	1689	24	0
1	D	1783	0	1694	19	0
2	A	5	0	0	2	0
3	A	137	0	0	3	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	126	0	0	4	1
3	C	87	0	0	5	0
3	D	123	0	0	3	0
All	All	7592	0	6737	74	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 5.

All (74) 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:139:GLN:NE2	3:B:302:HOH:O	2.04	0.85
1:C:18:ARG:NH1	3:C:302:HOH:O	2.12	0.83
1:C:108:ARG:O	3:C:301:HOH:O	1.99	0.81
1:B:67:SER:O	3:B:301:HOH:O	1.99	0.79
1:C:217:THR:HG22	1:C:243:VAL:H	1.51	0.74
1:A:172:GLU:OE1	3:A:402:HOH:O	2.07	0.73
1:A:98:PHE:N	2:A:301:SO4:O3	2.24	0.68
1:B:199:ASP:OD2	3:B:303:HOH:O	2.15	0.65
1:D:186:TYR:HB2	1:D:191:LYS:HE3	1.79	0.64
1:A:37:GLN:HB2	1:A:47:LEU:HD11	1.81	0.63
1:B:49:TYR:CB	1:B:231:GLU:HG2	2.29	0.62
1:C:135:GLY:HA2	1:C:241:VAL:HG12	1.82	0.61
1:C:193:ARG:NH2	1:C:216:ASP:OD1	2.35	0.59
1:B:138:VAL:HG21	1:B:144:LEU:HG	1.84	0.58
1:B:49:TYR:HB2	1:B:231:GLU:HG2	1.85	0.58
1:A:122:GLY:HA2	1:A:130:LEU:O	2.04	0.57
1:B:155:PHE:O	1:B:198:ARG:NH2	2.38	0.57
1:B:107:LYS:HD3	1:B:108:ARG:N	2.20	0.57
1:D:89:GLN:OE1	3:D:301:HOH:O	2.17	0.56
1:D:91:SER:HA	1:D:96:LEU:HD22	1.85	0.56
1:D:186:TYR:HB2	1:D:191:LYS:HG2	1.88	0.54
1:C:37:GLN:HB2	1:C:47:LEU:HD11	1.89	0.54
1:C:118:CYS:HB3	3:C:347:HOH:O	2.06	0.54
1:B:148:CYS:HB3	1:B:205:LEU:HB3	1.89	0.54
1:C:213:ARG:HB2	3:C:372:HOH:O	2.08	0.54
1:D:37:GLN:HB2	1:D:47:LEU:HD11	1.90	0.54
1:B:91:SER:HA	1:B:96:LEU:HD22	1.89	0.54
1:A:91:SER:HA	1:A:96:LEU:HD12	1.92	0.52
1:C:190:VAL:HB	1:C:194:PHE:CD2	2.45	0.51
1:D:209:MET:HE1	1:D:241:VAL:HG21	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:122:GLY:HA2	1:C:130:LEU:O	2.10	0.50
1:D:122:GLY:HA2	1:D:130:LEU:O	2.13	0.49
1:C:155:PHE:O	1:C:198:ARG:NH2	2.45	0.49
1:C:188:ASP:HA	1:C:191:LYS:HD2	1.95	0.49
1:A:106:ILE:O	1:A:107:LYS:HB2	2.13	0.48
1:A:217:THR:HG23	1:A:242:THR:HA	1.95	0.48
1:A:121:CYS:O	3:A:403:HOH:O	2.19	0.48
1:A:148:CYS:HB3	1:A:205:LEU:HB3	1.96	0.48
1:D:105:GLU:OE1	1:D:106:ILE:N	2.38	0.47
1:D:209:MET:HB3	1:D:212:LEU:HD21	1.95	0.47
1:C:138:VAL:O	1:C:243:VAL:HA	2.12	0.47
1:C:91:SER:HA	1:C:96:LEU:HD22	1.97	0.46
1:A:173:TRP:N	2:A:301:SO4:O2	2.48	0.46
1:B:122:GLY:O	3:B:305:HOH:O	2.21	0.46
1:C:217:THR:CG2	1:C:243:VAL:H	2.25	0.46
1:C:217:THR:HG22	1:C:243:VAL:N	2.27	0.46
1:C:27:GLN:NE2	3:C:303:HOH:O	2.28	0.45
1:D:28:SER:OG	3:D:302:HOH:O	2.20	0.45
1:B:122:GLY:HA2	1:B:130:LEU:O	2.17	0.45
1:B:143:SER:HA	1:B:212:LEU:HD23	1.99	0.44
1:C:24:ARG:HA	1:C:69:THR:O	2.18	0.44
1:D:139:GLN:H	1:D:139:GLN:CD	2.26	0.43
1:D:166:ALA:HB3	1:D:169:LYS:HB2	2.01	0.43
1:D:161:SER:HB2	1:D:225:TYR:CE2	2.54	0.43
1:A:107:LYS:HD2	1:A:107:LYS:HA	1.81	0.43
1:D:47:LEU:HA	1:D:58:VAL:HG21	2.00	0.43
1:A:48:ILE:HD13	1:A:54:LEU:HA	2.01	0.42
1:C:191:LYS:C	1:C:193:ARG:H	2.27	0.42
1:D:109:GLY:O	3:D:303:HOH:O	2.22	0.41
1:D:162:TRP:CE2	1:D:207:LEU:HB2	2.55	0.41
1:B:217:THR:HG23	1:B:242:THR:HA	2.02	0.41
1:B:161:SER:HB2	1:B:225:TYR:CE2	2.56	0.41
1:C:131:LEU:HD13	1:C:131:LEU:HA	1.78	0.41
1:C:199:ASP:OD1	1:C:201:SER:OG	2.29	0.41
1:A:107:LYS:N	3:A:404:HOH:O	2.49	0.41
1:B:209:MET:HE1	1:B:241:VAL:HG21	2.03	0.41
1:C:127:GLU:O	1:C:129:GLN:NE2	2.54	0.41
1:B:191:LYS:HE2	1:B:191:LYS:HB3	1.82	0.41
1:C:42:CYS:HB3	1:C:119:PRO:HD2	2.02	0.41
1:D:42:CYS:HB3	1:D:119:PRO:HD2	2.03	0.41
1:A:144:LEU:HD12	1:A:144:LEU:HA	1.96	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:169:LYS:HA	1:D:169:LYS:HD3	1.89	0.40
1:D:199:ASP:OD1	1:D:201:SER:OG	2.31	0.40
1:C:214:ALA:HA	1:C:217:THR:HG23	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:456:HOH:O	3:B:401:HOH:O[4_445]	2.13	0.07

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	238/252 (94%)	230 (97%)	7 (3%)	1 (0%)	30	15
1	B	243/252 (96%)	236 (97%)	7 (3%)	0	100	100
1	C	240/252 (95%)	228 (95%)	11 (5%)	1 (0%)	30	15
1	D	239/252 (95%)	233 (98%)	6 (2%)	0	100	100
All	All	960/1008 (95%)	927 (97%)	31 (3%)	2 (0%)	43	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	127	GLU
1	C	140	PRO

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	190/201 (94%)	189 (100%)	1 (0%)	81	72
1	B	190/201 (94%)	187 (98%)	3 (2%)	55	34
1	C	192/201 (96%)	188 (98%)	4 (2%)	47	24
1	D	192/201 (96%)	191 (100%)	1 (0%)	81	72
All	All	764/804 (95%)	755 (99%)	9 (1%)	63	45

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	96	LEU
1	B	89	GLN
1	B	213	ARG
1	B	231	GLU
1	C	131	LEU
1	C	145	ARG
1	C	243	VAL
1	C	244	SER
1	D	22	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	GLN
1	B	208	GLN
1	B	210	ASN
1	D	208	GLN
1	D	210	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

1 ligand 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	$\# Z > 2$
2	SO4	A	301	-	4,4,4	0.28	0	6,6,6	0.16	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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	SO4	2	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	242/252 (96%)	0.44	14 (5%) 29 32	15, 23, 34, 49	0
1	B	245/252 (97%)	0.53	11 (4%) 38 42	15, 24, 38, 52	0
1	C	244/252 (96%)	1.22	51 (20%) 2 3	17, 30, 57, 67	0
1	D	243/252 (96%)	0.46	7 (2%) 53 58	17, 25, 35, 40	0
All	All	974/1008 (96%)	0.66	83 (8%) 16 18	15, 25, 45, 67	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	140	PRO	7.5
1	C	243	VAL	6.6
1	A	123	GLY	6.3
1	C	138	VAL	5.6
1	C	194	PHE	5.5
1	C	141	GLY	5.1
1	C	214	ALA	5.0
1	C	213	ARG	4.9
1	A	124	SER	4.8
1	C	212	LEU	4.7
1	C	142	GLY	4.6
1	C	144	LEU	4.5
1	C	124	SER	4.4
1	A	245	SER	4.3
1	C	190	VAL	4.3
1	C	109	GLY	4.1
1	C	125	GLY	4.1
1	C	211	SER	4.0
1	B	109	GLY	4.0
1	C	242	THR	4.0
1	C	139	GLN	4.0

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Mol	Chain	Res	Type	RSRZ
1	C	244	SER	3.9
1	C	245	SER	3.9
1	A	126	GLY	3.9
1	B	126	GLY	3.8
1	C	210	ASN	3.6
1	A	1	ASP	3.6
1	C	143	SER	3.5
1	C	195	THR	3.5
1	B	127	GLU	3.4
1	C	137	LEU	3.4
1	A	106	ILE	3.3
1	C	134	GLY	3.2
1	A	127	GLU	3.2
1	B	125	GLY	3.2
1	B	106	ILE	3.1
1	C	166	ALA	3.1
1	C	136	GLY	3.0
1	D	191	LYS	3.0
1	C	188	ASP	2.9
1	C	128	VAL	2.9
1	C	217	THR	2.8
1	B	245	SER	2.8
1	C	146	LEU	2.8
1	C	108	ARG	2.8
1	C	209	MET	2.8
1	C	123	GLY	2.8
1	A	128	VAL	2.7
1	B	107	LYS	2.7
1	C	131	LEU	2.7
1	B	108	ARG	2.6
1	C	191	LYS	2.6
1	A	122	GLY	2.6
1	C	169	LYS	2.6
1	C	110	GLY	2.5
1	C	192	GLY	2.5
1	A	116	GLY	2.5
1	B	1	ASP	2.4
1	C	96	LEU	2.3
1	C	187	ALA	2.3
1	D	145	ARG	2.3
1	C	189	SER	2.3
1	D	127	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	148	CYS	2.3
1	C	168	GLY	2.3
1	A	131	LEU	2.2
1	D	243	VAL	2.2
1	D	139	GLN	2.2
1	C	1	ASP	2.1
1	C	216	ASP	2.1
1	C	133	SER	2.1
1	D	245	SER	2.1
1	A	19	VAL	2.1
1	C	127	GLU	2.1
1	A	234	PHE	2.1
1	C	218	ALA	2.0
1	C	145	ARG	2.0
1	A	11	LEU	2.0
1	C	215	GLU	2.0
1	B	123	GLY	2.0
1	C	220	TYR	2.0
1	B	128	VAL	2.0
1	D	107	LYS	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(Å ²)	Q<0.9
2	SO4	A	301	5/5	0.91	0.12	27,33,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.