



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:53 PM GMT

PDB ID : 3CD3  
Title : Crystal structure of phosphorylated human feline sarcoma viral oncogene homologue (v-FES) in complex with staurosporine and a consensus peptide  
Authors : Filippakopoulos, P.; Salah, E.; Cooper, C.; Picaud, S.S.; Elkins, J.M.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.M.; Weigelt, J.; Bountra, C.; Knapp, S.; Structural Genomics Consortium (SGC)  
Deposited on : 2008-02-26  
Resolution : 1.98 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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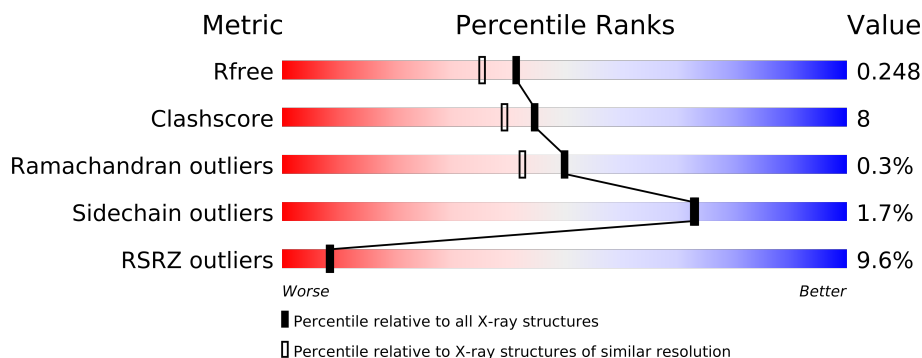
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.98 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))
$R_{free}$	66092	6577 (2.00-1.96)
Clashscore	79885	8091 (2.00-1.96)
Ramachandran outliers	78287	7989 (2.00-1.96)
Sidechain outliers	78261	7987 (2.00-1.96)
RSRZ outliers	66119	6578 (2.00-1.96)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	377	
2	B	6	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3232 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Proto-oncogene tyrosine-protein kinase Fes/Fps.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	353	Total	C	N	O	P	S	0	9	0
			2771	1777	477	502	1	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	446	SER	-	EXPRESSION TAG	UNP P07332
A	447	MET	-	EXPRESSION TAG	UNP P07332

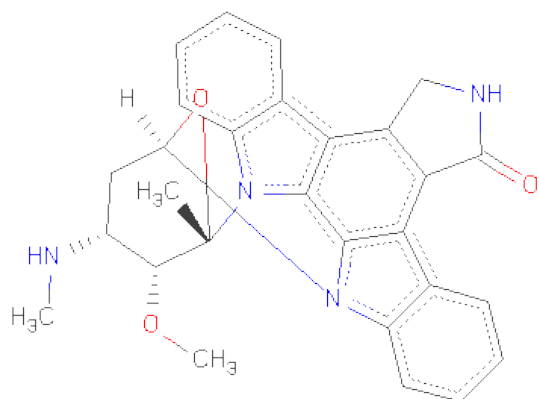
- Molecule 2 is a protein called Synthetic peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	0	0	0
			47	31	5	11			

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Cl	0	0
			2	2		

- Molecule 4 is STAUROSPORINE (three-letter code: STU) (formula: C<sub>28</sub>H<sub>26</sub>N<sub>4</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			35	28	4	3		
4	A	1	Total	C	N	O	0	0
			35	28	4	3		
4	A	1	Total	C	N	O	0	0
			35	28	4	3		

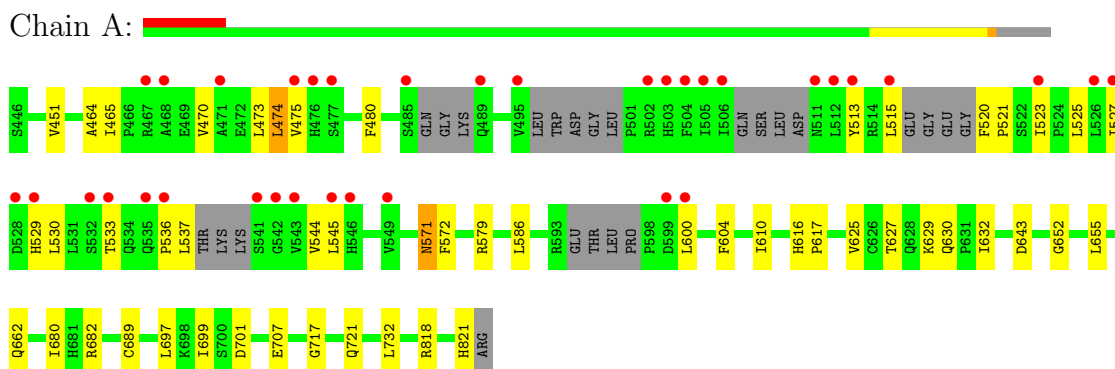
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	299	Total	O	0	0
			299	299		
5	B	8	Total	O	0	0
			8	8		

### 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: Proto-oncogene tyrosine-protein kinase Fes/Fps



- Molecule 2: Synthetic peptide

Chain B: 

There are no outlier residues recorded for this chain.

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	35.45Å 76.91Å 150.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.35 – 1.98 14.95 – 1.98	Depositor EDS
% Data completeness (in resolution range)	97.3 (15.35-1.98) 97.6 (14.95-1.98)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.95 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.4.0066	Depositor
R, $R_{free}$	0.185 , 0.247 0.190 , 0.248	Depositor DCC
$R_{free}$ test set	1447 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.7	Xtriage
Anisotropy	0.253	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.42 , 87.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 28827 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3232	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.88% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: STU, PTR, ACE, CL

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	0.68	0/2854	0.74	0/3863
2	B	0.81	0/45	0.89	0/59
All	All	0.68	0/2899	0.74	0/3922

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2771	0	2699	40	0
2	B	47	0	45	0	0
3	A	2	0	0	0	0
4	A	105	0	78	9	0
5	A	299	0	0	2	0
5	B	8	0	0	1	0
All	All	3232	0	2822	49	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (49) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:627:THR:HA	1:A:632:ILE:HD12	1.56	0.86
4:A:903:STU:H16	4:A:903:STU:H261	1.57	0.85
4:A:902:STU:H261	4:A:902:STU:H16	1.65	0.77
4:A:901:STU:H16	4:A:901:STU:H261	1.69	0.74
1:A:717:GLY:O	5:A:310:HOH:O	2.05	0.73
1:A:610:ILE:CD1	1:A:707[B]:GLU:HG3	2.29	0.63
1:A:625:VAL:HG12	1:A:627:THR:HG23	1.81	0.62
1:A:721:GLN:HG3	5:B:165:HOH:O	1.99	0.62
1:A:465:ILE:HD11	1:A:473:LEU:CD1	2.30	0.61
4:A:903:STU:H16	4:A:903:STU:C26	2.31	0.59
1:A:689:CYS:SG	1:A:699[A]:ILE:HD12	2.44	0.58
1:A:610:ILE:HD13	1:A:707[B]:GLU:HG3	1.86	0.57
1:A:465:ILE:HD11	1:A:473:LEU:HD12	1.87	0.56
1:A:530:LEU:HD13	1:A:537:LEU:CD2	2.35	0.56
1:A:610:ILE:HD13	1:A:707[B]:GLU:CG	2.36	0.55
1:A:610:ILE:HD13	1:A:707[B]:GLU:OE1	2.07	0.55
1:A:465:ILE:O	1:A:465:ILE:HG23	2.07	0.54
4:A:903:STU:C16	4:A:903:STU:H261	2.35	0.54
1:A:627:THR:HA	1:A:632:ILE:CD1	2.36	0.52
1:A:464:ALA:HB1	1:A:630:GLN:HE21	1.75	0.51
1:A:579:ARG:HG2	1:A:586:LEU:HD23	1.94	0.50
4:A:901:STU:H273	4:A:901:STU:C17	2.42	0.49
1:A:697:LEU:HD21	1:A:699[A]:ILE:CD1	2.42	0.49
1:A:652:GLY:HA2	1:A:655:LEU:HD12	1.95	0.48
1:A:662:GLN:HB2	1:A:818:ARG:NE	2.29	0.48
1:A:604:PHE:HE2	1:A:632:ILE:HG22	1.78	0.47
1:A:523:ILE:O	1:A:527:ILE:HG12	2.13	0.47
1:A:480:PHE:CG	1:A:545:LEU:HD22	2.49	0.47
4:A:903:STU:C16	4:A:903:STU:C26	2.91	0.47
1:A:465:ILE:HD11	1:A:473:LEU:HD11	1.96	0.47
1:A:697:LEU:HD23	1:A:697:LEU:C	2.36	0.46
1:A:470:VAL:HG12	1:A:474:LEU:HD22	1.96	0.46
1:A:529:HIS:CE1	1:A:533:THR:HG21	2.50	0.46
1:A:515:LEU:CD1	1:A:537:LEU:HD22	2.46	0.46
1:A:572:PHE:CE1	1:A:600:LEU:HG	2.51	0.46
1:A:629:LYS:HD2	5:A:332:HOH:O	2.17	0.45
1:A:536:PRO:HA	1:A:544:VAL:HG22	1.99	0.45
1:A:513:TYR:O	1:A:520:PHE:N	2.51	0.44
1:A:451:VAL:O	1:A:451:VAL:HG12	2.18	0.43
1:A:616:HIS:CG	1:A:617:PRO:HD2	2.53	0.43
1:A:521:PRO:HD2	1:A:525:LEU:HD13	2.01	0.43
1:A:465:ILE:O	1:A:465:ILE:CG2	2.67	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:732:LEU:HD23	1:A:732:LEU:C	2.41	0.41
4:A:902:STU:H261	4:A:902:STU:C16	2.42	0.41
1:A:464:ALA:HB3	1:A:632:ILE:HD11	2.03	0.41
4:A:901:STU:H16	4:A:901:STU:C26	2.43	0.40
1:A:680:ILE:HG22	1:A:682:ARG:HG3	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	347/377 (92%)	338 (97%)	8 (2%)	1 (0%)	50	42
2	B	4/6 (67%)	4 (100%)	0	0	100	100
All	All	351/383 (92%)	342 (97%)	8 (2%)	1 (0%)	50	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	701	ASP

### 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	293/326 (90%)	287 (98%)	6 (2%)	68	67
2	B	5/5 (100%)	5 (100%)	0	100	100
All	All	298/331 (90%)	292 (98%)	6 (2%)	73	67

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	474	LEU
1	A	475	VAL
1	A	571[A]	ASN
1	A	571[B]	ASN
1	A	643	ASP
1	A	821	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	529	HIS
1	A	630	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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	PTR	A	713	1	16,16,17	5.68	3 (18%)	20,22,24	1.53	2 (10%)

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	PTR	A	713	1	-	0/9/11/13	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	713	PTR	O-C	21.45	1.26	1.11
1	A	713	PTR	OH-CZ	-6.85	1.23	1.40
1	A	713	PTR	CA-C	2.28	1.52	1.48

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	713	PTR	C-CA-N	-5.49	108.35	113.83
1	A	713	PTR	OH-CZ-CE2	2.15	125.86	119.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	STU	A	901	-	40,42,42	1.92	9 (22%)	62,68,68	2.65	11 (17%)
4	STU	A	902	-	40,42,42	2.23	14 (35%)	62,68,68	2.48	15 (24%)
4	STU	A	903	-	40,42,42	2.32	10 (25%)	62,68,68	2.49	13 (20%)

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
4	STU	A	901	-	-	0/4/42/42	0/0/8/8
4	STU	A	902	-	-	0/4/42/42	0/0/8/8
4	STU	A	903	-	-	0/4/42/42	0/0/8/8

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	903	STU	C21-C22	6.56	1.61	1.53
4	A	902	STU	C20-N3	-5.83	1.31	1.39
4	A	901	STU	C21-C22	5.53	1.60	1.53
4	A	902	STU	C21-C22	5.39	1.60	1.53
4	A	901	STU	C7-C8	-5.20	1.40	1.49
4	A	902	STU	C7-C8	-4.64	1.41	1.49
4	A	903	STU	C7-C8	-4.52	1.41	1.49
4	A	903	STU	C24-C25	4.40	1.58	1.51
4	A	903	STU	C17-N2	-4.39	1.33	1.39
4	A	903	STU	C20-N3	-4.34	1.33	1.39
4	A	902	STU	C18-N2	-4.11	1.34	1.39
4	A	901	STU	O6-C22	3.83	1.50	1.42
4	A	901	STU	C24-C25	3.60	1.57	1.51
4	A	903	STU	O4-C25	3.59	1.49	1.43
4	A	903	STU	O6-C22	3.54	1.49	1.42
4	A	903	STU	C19-N3	-3.52	1.34	1.39
4	A	901	STU	C9-C10	3.42	1.55	1.50
4	A	903	STU	C11-C18	-3.27	1.38	1.42
4	A	902	STU	C9-N1	3.22	1.49	1.45
4	A	902	STU	C11-C18	-3.21	1.38	1.42
4	A	902	STU	O6-C22	3.18	1.49	1.42
4	A	902	STU	C26-C21	3.13	1.55	1.51
4	A	902	STU	C19-N3	-3.05	1.35	1.39
4	A	901	STU	C18-N2	-2.86	1.35	1.39
4	A	903	STU	C9-C10	2.71	1.54	1.50
4	A	902	STU	C17-N2	-2.45	1.36	1.39
4	A	902	STU	C24-C25	2.30	1.54	1.51
4	A	901	STU	C19-N3	-2.24	1.36	1.39
4	A	901	STU	C15-C16	2.23	1.41	1.36
4	A	902	STU	C8-N1	-2.14	1.31	1.34
4	A	901	STU	C6-C19	-2.04	1.40	1.42
4	A	902	STU	C6-C19	-2.02	1.40	1.42
4	A	902	STU	C15-C16	2.01	1.41	1.36

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	STU	C26-C21-N2	11.31	117.21	111.09
4	A	903	STU	C6-C7-C8	9.01	135.95	128.93
4	A	902	STU	C6-C7-C8	8.48	135.53	128.93
4	A	901	STU	C28-N4-C23	8.09	118.34	113.82
4	A	903	STU	C26-C21-N2	8.04	115.44	111.09
4	A	902	STU	C28-N4-C23	7.68	118.11	113.82
4	A	903	STU	C7-C8-N1	6.72	113.69	106.39
4	A	901	STU	C7-C8-N1	6.39	113.34	106.39
4	A	901	STU	C6-C7-C8	6.17	133.74	128.93
4	A	902	STU	C26-C21-N2	6.08	114.38	111.09
4	A	901	STU	C22-C21-N2	-5.91	104.11	112.78
4	A	902	STU	C22-C21-N2	-5.68	104.45	112.78
4	A	902	STU	C7-C8-N1	5.65	112.53	106.39
4	A	901	STU	O6-C22-C21	5.37	120.45	108.27
4	A	901	STU	C24-C25-N3	5.36	118.65	112.48
4	A	903	STU	C28-N4-C23	5.14	116.69	113.82
4	A	903	STU	C22-C21-N2	-5.11	105.28	112.78
4	A	902	STU	O5-C8-N1	-5.07	118.38	125.50
4	A	903	STU	O5-C8-N1	-5.03	118.43	125.50
4	A	903	STU	C24-C25-N3	4.14	117.25	112.48
4	A	902	STU	O6-C22-C21	4.13	117.65	108.27
4	A	902	STU	C25-C24-C23	-4.07	106.51	112.35
4	A	902	STU	C24-C25-N3	3.73	116.77	112.48
4	A	903	STU	O6-C22-C21	3.58	116.40	108.27
4	A	903	STU	O4-C25-C24	3.47	117.55	112.00
4	A	903	STU	C10-C7-C8	-3.38	105.33	108.89
4	A	903	STU	C25-C24-C23	-2.96	108.10	112.35
4	A	901	STU	C26-C21-C22	-2.68	106.48	112.87
4	A	902	STU	C9-N1-C8	-2.62	109.94	113.15
4	A	902	STU	O4-C25-C24	2.62	116.19	112.00
4	A	901	STU	O5-C8-N1	-2.58	121.87	125.50
4	A	903	STU	C9-N1-C8	-2.56	110.02	113.15
4	A	902	STU	C10-C7-C8	-2.52	106.23	108.89
4	A	903	STU	C22-C23-N4	2.45	114.91	110.54
4	A	901	STU	C9-N1-C8	-2.42	110.19	113.15
4	A	901	STU	C10-C7-C8	-2.37	106.39	108.89
4	A	902	STU	C10-C7-C6	-2.23	118.20	120.39
4	A	902	STU	C16-C17-C12	-2.21	117.54	120.73
4	A	902	STU	O4-C25-N3	-2.12	105.69	108.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	353/377 (93%)	0.37	35 (9%) <b>8</b> <b>8</b>	35, 47, 61, 74	0
2	B	6/6 (100%)	-0.43	0 <b>100</b> <b>100</b>	42, 43, 48, 49	0
All	All	359/383 (93%)	0.36	35 (9%) <b>8</b> <b>8</b>	35, 46, 61, 74	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	512	LEU	6.9
1	A	542	GLY	5.6
1	A	533	THR	5.2
1	A	468	ALA	4.7
1	A	475	VAL	4.4
1	A	476	HIS	4.3
1	A	600	LEU	4.3
1	A	532	SER	4.1
1	A	513	TYR	4.0
1	A	515	LEU	3.9
1	A	536	PRO	3.8
1	A	471	ALA	3.7
1	A	489	GLN	3.4
1	A	546	HIS	3.4
1	A	511	ASN	3.3
1	A	527	ILE	3.3
1	A	545	LEU	3.2
1	A	485	SER	3.1
1	A	541	SER	3.0
1	A	495	VAL	2.6
1	A	549	VAL	2.6
1	A	528	ASP	2.6
1	A	505	ILE	2.5
1	A	503	HIS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	599	ASP	2.5
1	A	506	ILE	2.5
1	A	543	VAL	2.4
1	A	529	HIS	2.4
1	A	535	GLN	2.4
1	A	502	ARG	2.4
1	A	477	SER	2.4
1	A	467	ARG	2.3
1	A	504	PHE	2.3
1	A	523	ILE	2.2
1	A	526	LEU	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, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	PTR	A	713	16/17	0.15	0.92	40,47,77,79	0

## 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, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	STU	A	903	35/35	0.14	0.43	36,45,63,73	0
3	CL	A	904	1/1	0.13	0.27	39,39,39,39	0
4	STU	A	902	35/35	0.10	-0.11	33,43,50,61	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	STU	A	901	35/35	0.08	-0.89	28,33,40,49	0
3	CL	A	905	1/1	0.05	-2.80	47,47,47,47	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.