



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 02:51 am GMT

PDB ID : 2WTK
Title : STRUCTURE OF THE HETEROTRIMERIC LKB1-STRADALPHA-MO25
ALPHA COMPLEX
Authors : Zeqiraj, E.; Van Aalten, D.M.F.
Deposited on : 2009-09-16
Resolution : 2.65 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

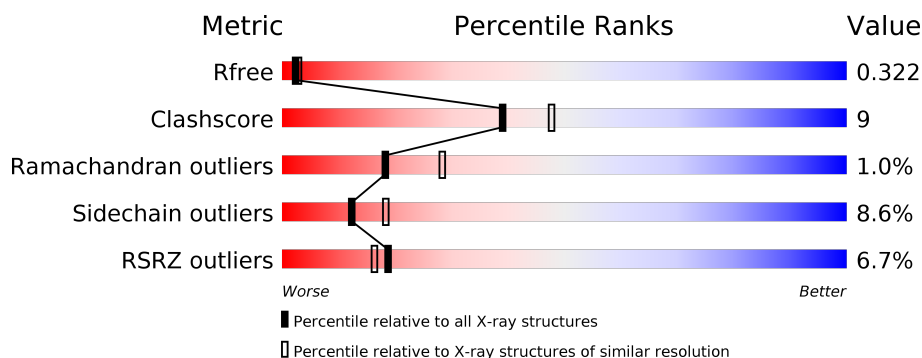
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	100719	3491 (2.70-2.62)
Clashscore	112137	1026 (2.68-2.64)
Ramachandran outliers	110173	1010 (2.68-2.64)
Sidechain outliers	110143	1010 (2.68-2.64)
RSRZ outliers	101464	3511 (2.70-2.62)

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	341	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>13%</div> <div>6%</div> </div> </div>
1	D	341	<div> <div>3%</div> <div> <div></div> <div>77%</div> <div>16%</div> <div>5%</div> </div> </div>
2	B	373	<div> <div>5%</div> <div> <div></div> <div>59%</div> <div>22%</div> <div>17%</div> </div> </div>
2	E	373	<div> <div>7%</div> <div> <div></div> <div>58%</div> <div>22%</div> <div>17%</div> </div> </div>
3	C	305	<div> <div>6%</div> <div> <div></div> <div>68%</div> <div>21%</div> <div>7%</div> </div> </div>
3	F	305	<div> <div>13%</div> <div> <div></div> <div>68%</div> <div>17%</div> <div>12%</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14850 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 CALCIUM-BINDING PROTEIN 39.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2655	1708	448	488	11			
1	D	325	Total	C	N	O	S	0	0	0
			2681	1721	450	499	11			

- Molecule 2 is a protein called STE20-RELATED KINASE ADAPTER PROTEIN ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	311	Total	C	N	O	S	0	0	0
			2473	1588	422	448	15			
2	E	308	Total	C	N	O	S	0	0	0
			2454	1576	419	444	15			

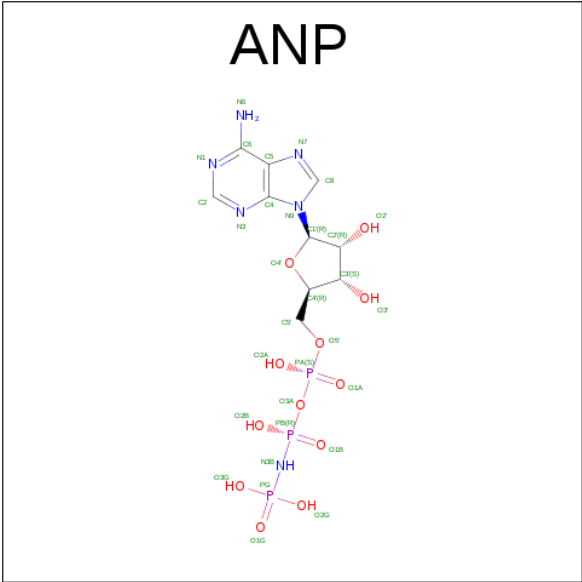
- Molecule 3 is a protein called SERINE/THREONINE-PROTEIN KINASE 11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	285	Total	C	N	O	S	0	0	0
			2254	1448	386	406	14			
3	F	268	Total	C	N	O	S	0	0	0
			2110	1360	363	374	13			

There are 2 discrepancies between the modelled and reference sequences:

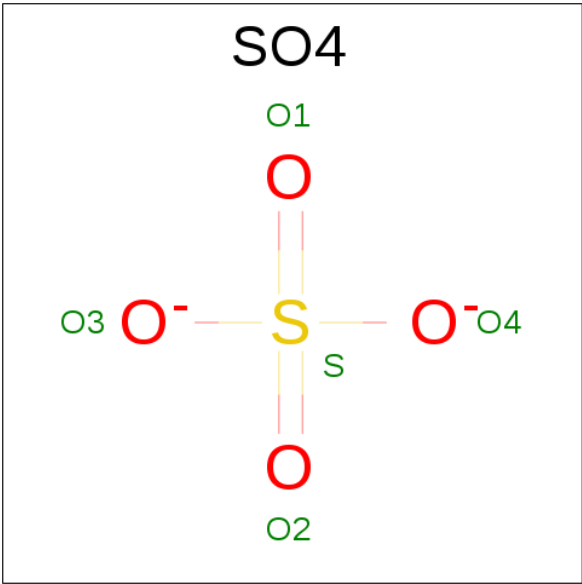
Chain	Residue	Modelled	Actual	Comment	Reference
C	194	ALA	ASP	ENGINEERED MUTATION	UNP Q15831
F	194	ALA	ASP	ENGINEERED MUTATION	UNP Q15831

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
4	C	1	Total	C	N	O	P	0	0
			27	10	6	9	2		
4	E	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
4	F	1	Total	C	N	O	P	0	0
			27	10	6	9	2		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

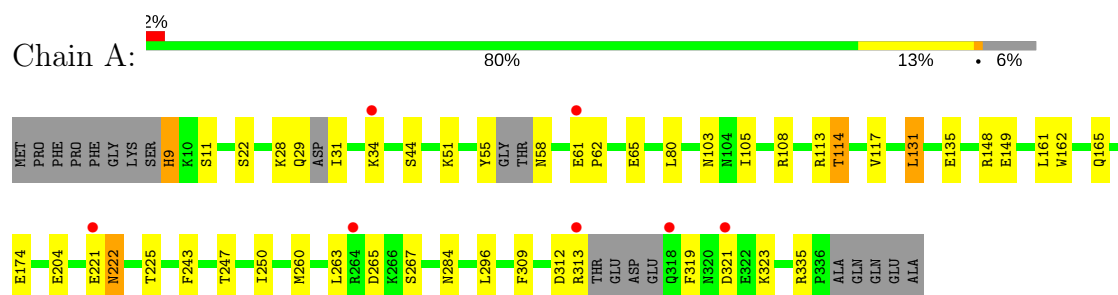
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	24	Total	O	0	0
			24	24		
6	B	8	Total	O	0	0
			8	8		
6	C	18	Total	O	0	0
			18	18		
6	D	26	Total	O	0	0
			26	26		
6	E	18	Total	O	0	0
			18	18		
6	F	8	Total	O	0	0
			8	8		

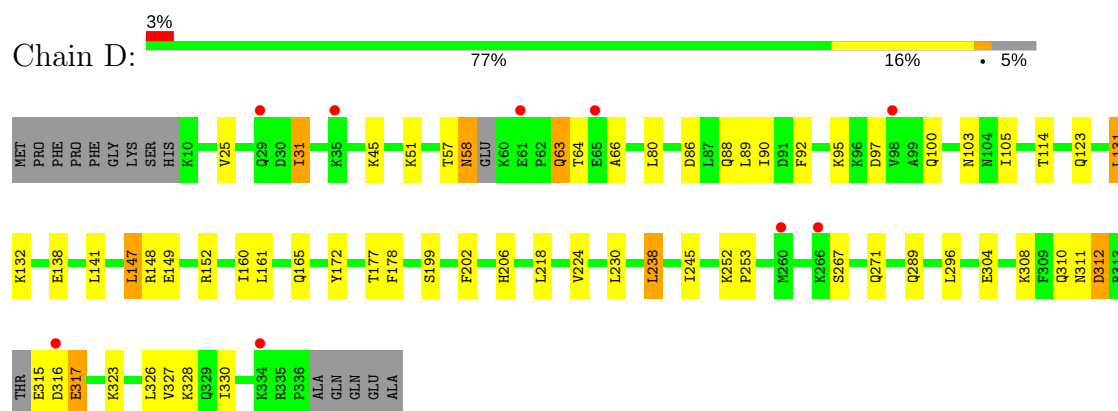
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

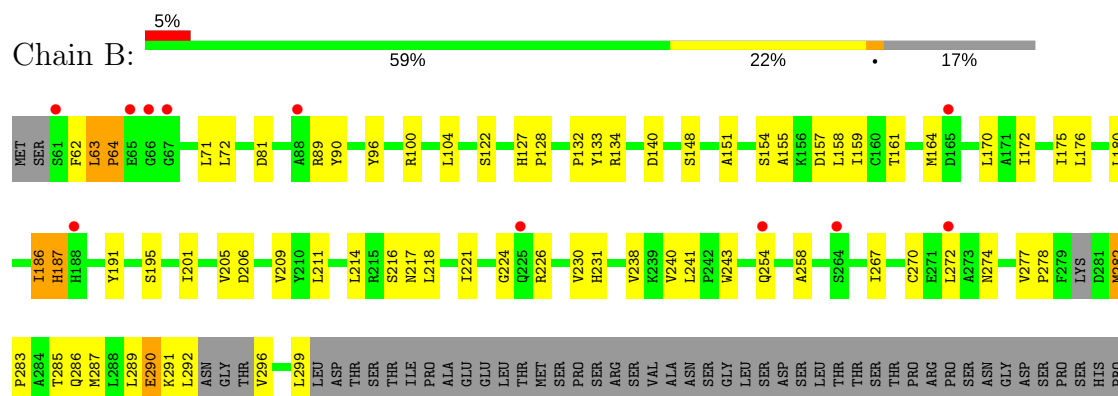
- Molecule 1: CALCIUM-BINDING PROTEIN 39

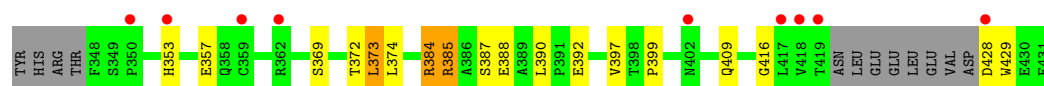


- Molecule 1: CALCIUM-BINDING PROTEIN 39

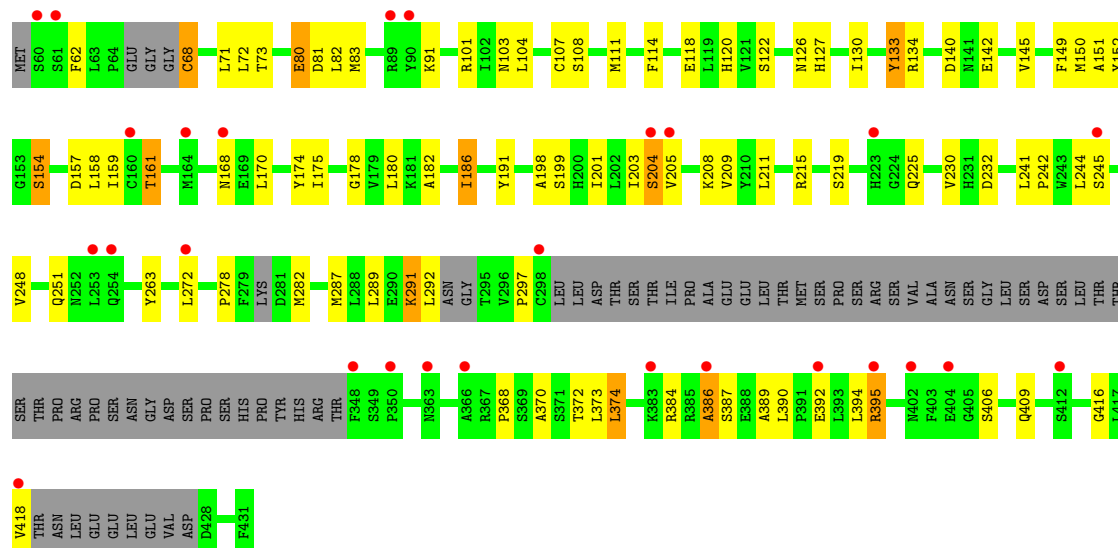


- Molecule 2: STE20-RELATED KINASE ADAPTER PROTEIN ALPHA

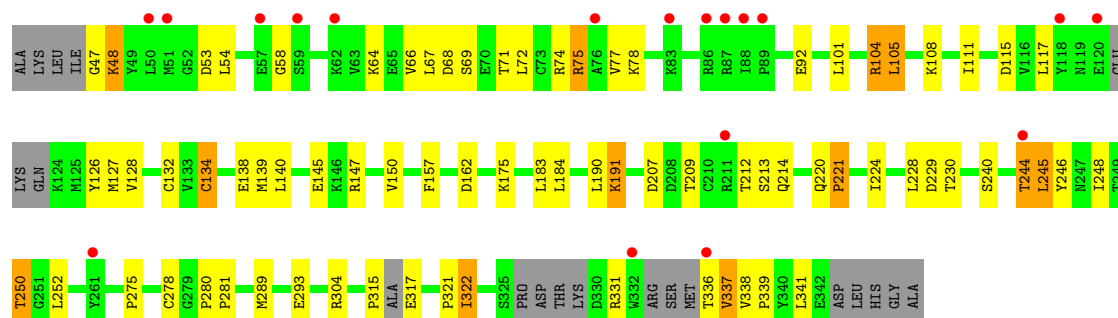




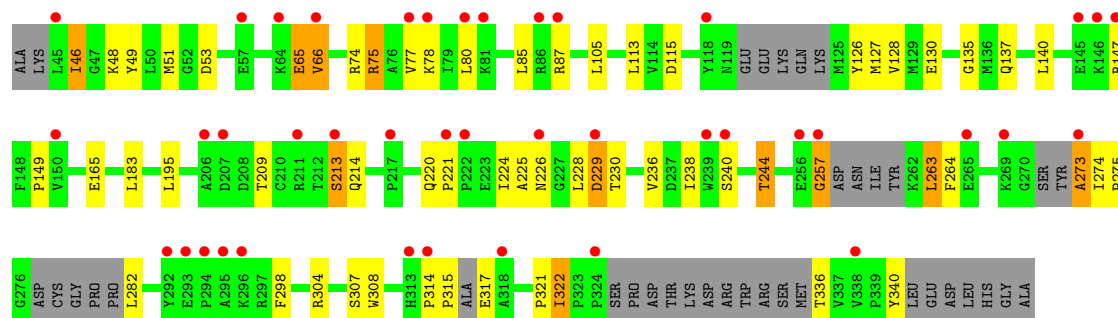
• Molecule 2: STE20-RELATED KINASE ADAPTER PROTEIN ALPHA



• Molecule 3: SERINE/THREONINE-PROTEIN KINASE 11



• Molecule 3: SERINE/THREONINE-PROTEIN KINASE 11



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.36Å 118.36Å 390.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.90 – 2.65 19.96 – 2.65	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.90-2.65) 99.2 (19.96-2.65)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 2.67Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.240 , 0.291 0.286 , 0.322	Depositor DCC
R_{free} test set	917 reflections (1.00%)	DCC
Wilson B-factor (Å ²)	60.4	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 60.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.015 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	14850	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% 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: ANP, 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	0.54	0/2700	0.62	0/3627
1	D	0.56	0/2726	0.62	0/3664
2	B	0.55	0/2534	0.69	0/3432
2	E	0.57	0/2514	0.67	0/3404
3	C	0.48	0/2305	0.62	1/3110 (0.0%)
3	F	0.68	5/2153 (0.2%)	0.62	0/2899
All	All	0.56	5/14932 (0.0%)	0.64	1/20136 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	273	ALA	CA-CB	15.55	1.85	1.52
3	F	273	ALA	N-CA	8.80	1.64	1.46
3	F	273	ALA	C-N	8.24	1.52	1.34
3	F	273	ALA	C-O	7.15	1.36	1.23
3	F	257	GLY	CA-C	6.12	1.61	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	C	105	LEU	CA-CB-CG	5.13	127.10	115.30

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	2655	0	2710	30	0
1	D	2681	0	2728	31	0
2	B	2473	0	2439	52	0
2	E	2454	0	2420	53	0
3	C	2254	0	2280	54	0
3	F	2110	0	2162	37	0
4	B	31	0	13	2	0
4	C	27	0	12	3	0
4	E	31	0	13	1	0
4	F	27	0	12	0	0
5	D	5	0	0	1	0
6	A	24	0	0	0	0
6	B	8	0	0	1	0
6	C	18	0	0	1	0
6	D	26	0	0	0	0
6	E	18	0	0	0	0
6	F	8	0	0	1	0
All	All	14850	0	14789	252	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 9.

The worst 5 of 252 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)
3:F:273:ALA:CB	3:F:273:ALA:CA	1.85	1.50
3:C:104:ARG:HH11	3:C:104:ARG:HG3	1.11	1.10
2:B:161:THR:HG21	2:B:416:GLY:HA3	1.07	1.05
2:B:384:ARG:HG2	2:B:384:ARG:HH11	1.26	1.01
2:B:161:THR:CG2	2:B:416:GLY:HA3	1.95	0.97

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	313/341 (92%)	300 (96%)	11 (4%)	2 (1%)	28	43
1	D	319/341 (94%)	301 (94%)	16 (5%)	2 (1%)	28	43
2	B	301/373 (81%)	277 (92%)	20 (7%)	4 (1%)	14	22
2	E	296/373 (79%)	270 (91%)	22 (7%)	4 (1%)	13	20
3	C	275/305 (90%)	258 (94%)	15 (6%)	2 (1%)	25	39
3	F	254/305 (83%)	235 (92%)	15 (6%)	4 (2%)	11	17
All	All	1758/2038 (86%)	1641 (93%)	99 (6%)	18 (1%)	18	28

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	ASN
2	B	62	PHE
1	D	317	GLU
2	E	387	SER
3	F	214	GLN

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	296/312 (95%)	278 (94%)	18 (6%)	22	35
1	D	299/312 (96%)	277 (93%)	22 (7%)	16	26
2	B	274/331 (83%)	245 (89%)	29 (11%)	8	12
2	E	273/331 (82%)	243 (89%)	30 (11%)	7	11
3	C	246/262 (94%)	225 (92%)	21 (8%)	12	19
3	F	230/262 (88%)	211 (92%)	19 (8%)	13	20
All	All	1618/1810 (89%)	1479 (91%)	139 (9%)	12	18

5 of 139 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	245	LEU
1	D	132	LYS
3	F	115	ASP
3	C	278	CYS
1	D	58	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	63	GLN
1	D	228	GLN
1	D	100	GLN
2	B	254	GLN
1	D	205	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

5 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
4	ANP	B	432	-	29,33,33	2.00	6 (20%)	28,52,52	2.02	8 (28%)
4	ANP	C	2	-	24,29,33	1.56	3 (12%)	22,45,52	1.81	4 (18%)
5	SO4	D	342	-	4,4,4	0.17	0	6,6,6	0.12	0
4	ANP	E	432	-	29,33,33	2.07	7 (24%)	28,52,52	2.04	7 (25%)
4	ANP	F	4	-	24,29,33	1.61	4 (16%)	22,45,52	1.68	3 (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
4	ANP	B	432	-	-	0/13/38/38	0/3/3/3
4	ANP	C	2	-	-	0/9/32/38	0/3/3/3
5	SO4	D	342	-	-	0/0/0/0	0/0/0/0
4	ANP	E	432	-	-	1/13/38/38	0/3/3/3
4	ANP	F	4	-	-	0/9/32/38	0/3/3/3

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	432	ANP	PB-O2B	-2.21	1.50	1.56
4	E	432	ANP	PB-O2B	-2.14	1.50	1.56
4	F	4	ANP	C2-N3	2.09	1.35	1.32
4	E	432	ANP	PB-O3A	2.32	1.62	1.59
4	C	2	ANP	PB-O3A	2.35	1.62	1.59

The worst 5 of 22 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	432	ANP	N3-C2-N1	-7.04	122.73	128.86
4	B	432	ANP	N3-C2-N1	-6.57	123.14	128.86
4	C	2	ANP	N3-C2-N1	-5.91	123.71	128.86
4	F	4	ANP	N3-C2-N1	-5.68	123.91	128.86
4	E	432	ANP	O1G-PG-N3B	-4.35	105.29	111.79

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	432	ANP	O1B-PB-N3B-PG

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	432	ANP	2	0
4	C	2	ANP	3	0
5	D	342	SO4	1	0
4	E	432	ANP	1	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 [i](#)

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

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	321/341 (94%)	0.09	7 (2%) 62 61	28, 57, 107, 145	0
1	D	325/341 (95%)	0.10	9 (2%) 53 52	30, 61, 105, 155	0
2	B	311/373 (83%)	0.66	20 (6%) 20 18	51, 72, 106, 139	0
2	E	308/373 (82%)	0.57	27 (8%) 11 9	51, 74, 114, 146	0
3	C	285/305 (93%)	0.42	18 (6%) 21 19	39, 75, 128, 194	0
3	F	268/305 (87%)	0.88	41 (15%) 2 1	47, 95, 137, 171	0
All	All	1818/2038 (89%)	0.44	122 (6%) 19 16	28, 71, 121, 194	0

The worst 5 of 122 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	206	ALA	5.5
2	E	348	PHE	5.4
3	C	332	TRP	4.8
2	E	253	LEU	4.8
2	E	350	PRO	4.7

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 carbohydrates 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. 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(\AA^2)	Q<0.9
4	ANP	F	4	27/31	0.85	0.24	-0.16	89,98,102,103	0
4	ANP	B	432	31/31	0.90	0.20	-0.40	44,56,69,73	0
4	ANP	C	2	27/31	0.85	0.17	-0.65	82,102,121,123	0
4	ANP	E	432	31/31	0.93	0.17	-0.73	48,59,81,83	0
5	SO4	D	342	5/5	0.95	0.28	-	69,75,78,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.