



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:52 AM GMT

PDB ID : 2OIC
Title : Crystal structure of IRAK4 kinase domain complexed with staurosporine
Authors : Kuglstatter, A.; Villasenor, A.G.; Browner, M.F.
Deposited on : 2007-01-10
Resolution : 2.40 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
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

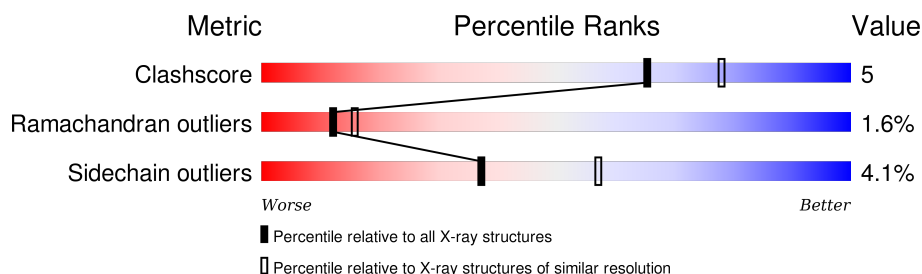
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	301	 79% 12% • 7%
1	B	301	 77% 13% • 8%
1	C	301	 80% 13% 7%
1	D	301	 86% 9% • 5%

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	STU	A	1	X	-	-	-
2	STU	B	2	X	-	-	-
2	STU	C	3	X	-	-	-
2	STU	D	4	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9302 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.

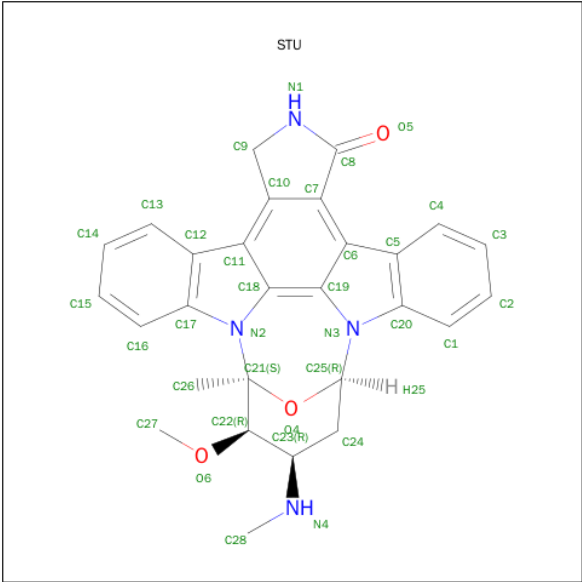
- Molecule 1 is a protein called Interleukin-1 receptor-associated kinase 4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	279	Total	C	N	O	P	S	0	0	0
			2210	1385	373	435	3	14			
1	B	278	Total	C	N	O	P	S	0	0	0
			2196	1378	372	430	2	14			
1	C	280	Total	C	N	O	P	S	0	0	0
			2221	1390	371	443	3	14			
1	D	287	Total	C	N	O	P	S	0	0	0
			2266	1417	382	450	3	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	342	TPO	THR	MODIFIED RESIDUE	UNP Q9NWZ3
A	345	TPO	THR	MODIFIED RESIDUE	UNP Q9NWZ3
A	346	SEP	SER	MODIFIED RESIDUE	UNP Q9NWZ3
B	342	TPO	THR	MODIFIED RESIDUE	UNP Q9NWZ3
B	345	TPO	THR	MODIFIED RESIDUE	UNP Q9NWZ3
B	346	SEP	SER	MODIFIED RESIDUE	UNP Q9NWZ3
C	342	TPO	THR	MODIFIED RESIDUE	UNP Q9NWZ3
C	345	TPO	THR	MODIFIED RESIDUE	UNP Q9NWZ3
C	346	SEP	SER	MODIFIED RESIDUE	UNP Q9NWZ3
D	342	TPO	THR	MODIFIED RESIDUE	UNP Q9NWZ3
D	345	TPO	THR	MODIFIED RESIDUE	UNP Q9NWZ3
D	346	SEP	SER	MODIFIED RESIDUE	UNP Q9NWZ3

- Molecule 2 is STAUROSPORINE (three-letter code: STU) (formula: $C_{28}H_{26}N_4O_3$).



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

- Molecule 3 is water.

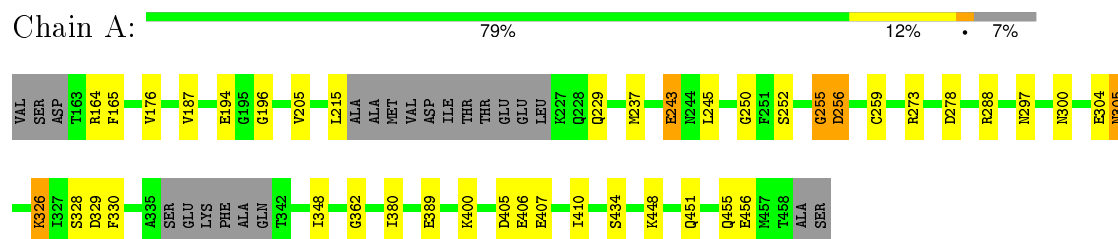
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	78	Total	O	0	0
			78	78		
3	B	58	Total	O	0	0
			58	58		
3	C	59	Total	O	0	0
			59	59		
3	D	74	Total	O	0	0
			74	74		

3 Residue-property plots [i](#)

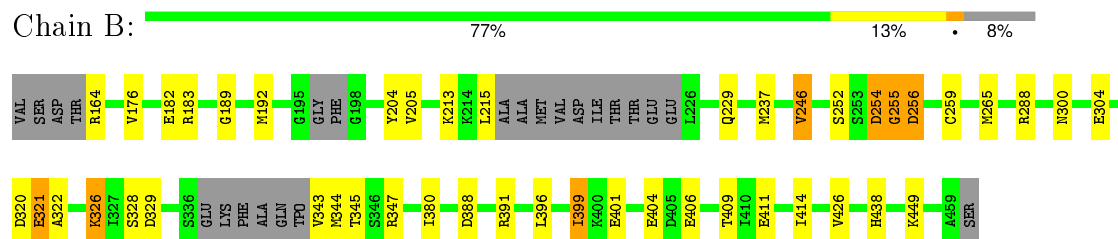
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

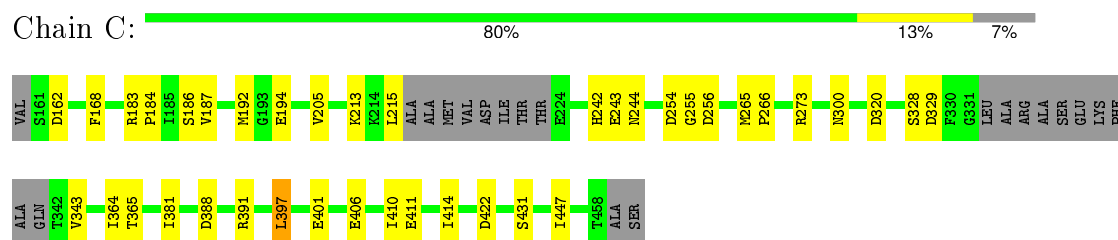
- Molecule 1: Interleukin-1 receptor-associated kinase 4



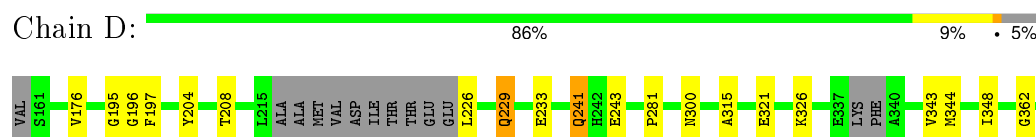
- Molecule 1: Interleukin-1 receptor-associated kinase 4



- Molecule 1: Interleukin-1 receptor-associated kinase 4



- Molecule 1: Interleukin-1 receptor-associated kinase 4





4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	153.89Å 141.32Å 88.26Å 90.00° 122.27° 90.00°	Depositor
Resolution (Å)	48.17 – 2.40	Depositor
% Data completeness (in resolution range)	99.9 (48.17-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	7.10	Depositor
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.223 , 0.278	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9302	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

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: TPO, STU, SEP

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	0.55	0/2214	0.62	0/2981
1	B	0.50	0/2209	0.61	0/2973
1	C	0.50	0/2225	0.61	1/2996 (0.0%)
1	D	0.52	0/2269	0.59	0/3054
All	All	0.52	0/8917	0.61	1/12004 (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	397	LEU	CA-CB-CG	5.74	128.50	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	2210	0	2175	22	0
1	B	2196	0	2168	26	0
1	C	2221	0	2173	17	0
1	D	2266	0	2222	16	0
2	A	35	0	26	6	0
2	B	35	0	26	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	35	0	26	2	0
2	D	35	0	26	2	0
3	A	78	0	0	2	0
3	B	58	0	0	1	0
3	C	59	0	0	2	0
3	D	74	0	0	0	0
All	All	9302	0	8842	90	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 5.

All (90) 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:265:MET:HE1	1:B:320:ASP:HB3	1.52	0.90
2:A:1:STU:H16	2:A:1:STU:H261	1.60	0.83
1:D:411:GLU:HA	1:D:414:ILE:HD12	1.62	0.81
1:D:176:VAL:HG21	1:D:204:TYR:O	1.83	0.79
1:C:242:HIS:HE1	3:C:2240:HOH:O	1.69	0.74
1:B:265:MET:CE	1:B:320:ASP:HB3	2.17	0.73
2:C:3:STU:H261	2:C:3:STU:H16	1.69	0.73
2:B:2:STU:H16	2:B:2:STU:H261	1.73	0.70
2:D:4:STU:H16	2:D:4:STU:H261	1.73	0.69
1:D:411:GLU:HA	1:D:414:ILE:CD1	2.23	0.68
2:B:2:STU:C17	2:B:2:STU:H273	2.24	0.68
1:B:237:MET:HE2	1:B:237:MET:HA	1.75	0.68
1:B:246:VAL:CG1	1:B:328:SER:HB3	2.25	0.67
1:B:246:VAL:HG11	1:B:328:SER:HB3	1.80	0.63
1:A:243:GLU:O	1:A:326:LYS:NZ	2.32	0.63
1:A:305:ASN:HD22	1:A:305:ASN:N	1.96	0.63
1:A:328:SER:OG	2:A:1:STU:H271	2.00	0.62
1:B:176:VAL:HG22	1:B:204:TYR:O	2.00	0.60
1:B:321:GLU:OE1	3:B:3020:HOH:O	2.16	0.59
1:B:388:ASP:HB3	1:B:391:ARG:HB3	1.83	0.59
1:A:297:ASN:OD1	1:A:451:GLN:NE2	2.33	0.59
1:B:254:ASP:O	1:B:255:GLY:O	2.21	0.59
1:A:252:SER:HB3	1:A:259:CYS:HB2	1.84	0.58
1:A:255:GLY:O	1:A:256:ASP:HB2	2.04	0.56
1:A:389:GLU:HG2	3:A:2031:HOH:O	2.06	0.56
1:C:388:ASP:HB3	1:C:391:ARG:HB3	1.87	0.56
1:B:396:LEU:O	1:B:399:ILE:HB	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:343:VAL:HG23	1:C:364:ILE:HB	1.87	0.54
1:C:183:ARG:HB3	1:C:187:VAL:HG11	1.89	0.54
1:C:411:GLU:HA	1:C:414:ILE:HD12	1.90	0.54
1:C:168:PHE:CE2	1:C:205:VAL:HG11	2.43	0.53
1:C:213:LYS:HE3	1:C:215:LEU:HD21	1.91	0.52
1:B:300:ASN:O	1:B:304:GLU:HB2	2.09	0.52
2:A:1:STU:O5	2:A:1:STU:H4	2.10	0.52
2:A:1:STU:H16	2:A:1:STU:C26	2.36	0.51
1:C:422:ASP:N	1:C:422:ASP:OD2	2.43	0.51
1:A:451:GLN:O	1:A:455:GLN:HG2	2.10	0.50
1:A:245:LEU:HG	1:A:330:PHE:HZ	1.76	0.50
1:D:229:GLN:O	1:D:233:GLU:HB2	2.11	0.50
1:D:343:VAL:HG22	1:D:344:MET:H	1.77	0.50
1:B:328:SER:OG	2:B:2:STU:H271	2.12	0.49
1:A:176:VAL:HG21	1:A:205:VAL:HG22	1.93	0.49
1:D:176:VAL:CG2	1:D:204:TYR:O	2.59	0.49
1:B:322:ALA:HB2	1:D:281:PRO:HB3	1.95	0.49
1:B:411:GLU:HA	1:B:414:ILE:HD12	1.94	0.49
1:B:414:ILE:HG12	1:B:426:VAL:HG11	1.96	0.48
1:B:213:LYS:HE3	1:B:215:LEU:HD21	1.94	0.47
1:A:305:ASN:ND2	1:A:305:ASN:N	2.60	0.47
1:A:278:ASP:O	1:D:419:ASN:HB3	2.15	0.47
1:D:243:GLU:O	1:D:326:LYS:HE3	2.15	0.47
1:D:416:LYS:HD3	1:D:416:LYS:HA	1.77	0.47
1:C:242:HIS:CD2	1:C:244:ASN:H	2.33	0.47
1:D:300:ASN:HA	1:D:447:ILE:HG21	1.97	0.47
1:B:255:GLY:O	1:B:256:ASP:CB	2.62	0.46
1:D:348:ILE:HG12	1:D:362:GLY:HA2	1.97	0.46
1:D:315:ALA:HB1	2:D:4:STU:H283	1.96	0.46
1:B:388:ASP:HB3	1:B:391:ARG:CB	2.45	0.46
1:A:348:ILE:HG12	1:A:362:GLY:HA2	1.98	0.45
1:A:215:LEU:O	3:A:2006:HOH:O	2.21	0.45
1:D:241:GLN:HE21	1:D:241:GLN:HB2	1.65	0.45
2:C:3:STU:O5	2:C:3:STU:H4	2.16	0.45
2:A:1:STU:C16	2:A:1:STU:H261	2.38	0.45
2:B:2:STU:C16	2:B:2:STU:H261	2.46	0.45
1:C:273:ARG:NE	1:C:273:ARG:HA	2.31	0.45
1:C:300:ASN:HA	1:C:447:ILE:HG21	1.97	0.44
1:B:265:MET:CE	1:B:326:LYS:HD3	2.47	0.44
1:B:252:SER:HB3	1:B:259:CYS:HB2	2.00	0.43
1:B:176:VAL:HG21	1:B:205:VAL:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:195:GLY:O	1:D:197:PHE:N	2.47	0.43
1:A:165:PHE:HB3	1:A:250:GLY:HA2	2.01	0.43
1:D:414:ILE:HG12	1:D:426:VAL:HG11	2.00	0.43
1:A:304:GLU:HB3	1:A:305:ASN:ND2	2.34	0.42
1:B:183:ARG:O	1:B:189:GLY:HA3	2.19	0.42
1:B:288:ARG:HB3	1:B:380:ILE:HG23	2.01	0.42
1:C:184:PRO:O	1:C:187:VAL:HG12	2.19	0.42
1:C:265:MET:HA	1:C:266:PRO:HD3	1.88	0.42
1:A:400:LYS:HE3	1:A:434:SER:O	2.19	0.42
1:A:405:ASP:O	1:A:407:GLU:N	2.53	0.42
1:B:246:VAL:HG13	1:B:328:SER:HB3	1.99	0.41
1:A:300:ASN:O	1:A:304:GLU:HB2	2.19	0.41
2:A:1:STU:C16	2:A:1:STU:C26	2.97	0.41
1:B:345:TPO:O1P	1:B:347:ARG:NH2	2.53	0.41
1:A:237:MET:HA	1:A:237:MET:HE2	2.01	0.41
1:A:288:ARG:HB3	1:A:380:ILE:HG23	2.02	0.41
1:C:242:HIS:CE1	3:C:2240:HOH:O	2.57	0.41
1:C:381:ILE:HG21	1:C:410:ILE:HD11	2.03	0.41
1:C:254:ASP:O	1:C:256:ASP:N	2.54	0.40
1:A:273:ARG:NE	1:A:273:ARG:HA	2.36	0.40
1:C:266:PRO:HD2	1:C:320:ASP:HA	2.02	0.40
1:B:288:ARG:HB3	1:B:380:ILE:CG2	2.52	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	271/301 (90%)	253 (93%)	12 (4%)	6 (2%)	8	9
1	B	268/301 (89%)	251 (94%)	12 (4%)	5 (2%)	10	12
1	C	272/301 (90%)	257 (94%)	11 (4%)	4 (2%)	13	17

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	278/301 (92%)	265 (95%)	11 (4%)	2 (1%)	26	38
All	All	1089/1204 (90%)	1026 (94%)	46 (4%)	17 (2%)	12	16

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	255	GLY
1	A	256	ASP
1	A	406	GLU
1	B	254	ASP
1	B	255	GLY
1	B	256	ASP
1	D	406	GLU
1	B	406	GLU
1	C	186	SER
1	C	255	GLY
1	B	329	ASP
1	A	164	ARG
1	C	406	GLU
1	C	329	ASP
1	D	196	GLY
1	A	196	GLY
1	A	329	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	241/259 (93%)	232 (96%)	9 (4%)	41	62
1	B	241/259 (93%)	226 (94%)	15 (6%)	23	35
1	C	244/259 (94%)	235 (96%)	9 (4%)	41	62
1	D	247/259 (95%)	240 (97%)	7 (3%)	51	72
All	All	973/1036 (94%)	933 (96%)	40 (4%)	37	57

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	187	VAL
1	A	194	GLU
1	A	229	GLN
1	A	243	GLU
1	A	305	ASN
1	A	326	LYS
1	A	410	ILE
1	A	448	LYS
1	A	456	GLU
1	B	164	ARG
1	B	182	GLU
1	B	192	MET
1	B	229	GLN
1	B	246	VAL
1	B	321	GLU
1	B	326	LYS
1	B	343	VAL
1	B	344	MET
1	B	399	ILE
1	B	401	GLU
1	B	404	GLU
1	B	409	THR
1	B	438	HIS
1	B	449	LYS
1	C	162	ASP
1	C	192	MET
1	C	194	GLU
1	C	243	GLU
1	C	328	SER
1	C	365	THR
1	C	397	LEU
1	C	401	GLU
1	C	431	SER
1	D	208	THR
1	D	226	LEU
1	D	229	GLN
1	D	241	GLN
1	D	321	GLU
1	D	407	GLU
1	D	441	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	HIS
1	A	241	GLN
1	A	293	GLN
1	A	305	ASN
1	A	394	GLN
1	B	166	HIS
1	B	207	ASN
1	B	229	GLN
1	B	455	GLN
1	C	179	ASN
1	C	242	HIS
1	D	206	ASN
1	D	241	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

11 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	TPO	A	342	1	8,10,11	0.68	0	7,14,16	1.15	1 (14%)
1	TPO	A	345	1	8,10,11	0.63	0	7,14,16	1.23	0
1	SEP	A	346	1	8,9,10	1.58	2 (25%)	8,12,14	1.67	1 (12%)
1	TPO	B	345	1	8,10,11	0.62	0	7,14,16	1.97	2 (28%)
1	SEP	B	346	1	8,9,10	1.54	1 (12%)	8,12,14	2.03	2 (25%)
1	TPO	C	342	1	8,10,11	0.64	0	7,14,16	1.35	0
1	TPO	C	345	1	8,10,11	0.69	0	7,14,16	1.44	1 (14%)
1	SEP	C	346	1	8,9,10	1.58	1 (12%)	8,12,14	2.00	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	TPO	D	342	1	8,10,11	0.60	0	7,14,16	1.42	1 (14%)
1	TPO	D	345	1	8,10,11	0.65	0	7,14,16	1.26	1 (14%)
1	SEP	D	346	1	8,9,10	1.54	1 (12%)	8,12,14	2.01	2 (25%)

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	TPO	A	342	1	-	0/8/11/13	0/0/0/0
1	TPO	A	345	1	-	1/8/11/13	0/0/0/0
1	SEP	A	346	1	-	0/6/8/10	0/0/0/0
1	TPO	B	345	1	-	1/8/11/13	0/0/0/0
1	SEP	B	346	1	-	0/6/8/10	0/0/0/0
1	TPO	C	342	1	-	0/8/11/13	0/0/0/0
1	TPO	C	345	1	-	1/8/11/13	0/0/0/0
1	SEP	C	346	1	-	0/6/8/10	0/0/0/0
1	TPO	D	342	1	-	0/8/11/13	0/0/0/0
1	TPO	D	345	1	-	1/8/11/13	0/0/0/0
1	SEP	D	346	1	-	0/6/8/10	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	346	SEP	P-O2P	2.01	1.61	1.54
1	B	346	SEP	P-O1P	3.10	1.61	1.51
1	D	346	SEP	P-O1P	3.16	1.61	1.51
1	C	346	SEP	P-O1P	3.18	1.61	1.51
1	A	346	SEP	P-O1P	3.24	1.61	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	346	SEP	O-C-CA	-2.04	120.19	125.49
1	C	346	SEP	O-C-CA	-2.03	120.21	125.49
1	D	346	SEP	O-C-CA	-2.01	120.25	125.49
1	A	342	TPO	O3P-P-O2P	2.01	115.02	107.38
1	D	342	TPO	C-CA-N	2.09	114.20	109.83
1	B	345	TPO	O3P-P-O2P	2.12	115.46	107.38
1	D	345	TPO	O3P-P-O2P	2.13	115.48	107.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	345	TPO	C-CA-N	2.39	114.82	109.83
1	B	345	TPO	C-CA-N	3.77	117.71	109.83
1	A	346	SEP	OG-CB-CA	3.95	111.64	108.27
1	C	346	SEP	OG-CB-CA	4.80	112.37	108.27
1	B	346	SEP	OG-CB-CA	4.97	112.52	108.27
1	D	346	SEP	OG-CB-CA	4.98	112.52	108.27

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	345	TPO	OG1-CB-CA-N
1	B	345	TPO	OG1-CB-CA-N
1	D	345	TPO	OG1-CB-CA-N
1	A	345	TPO	OG1-CB-CA-N

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	345	TPO	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	STU	A	1	-	27,42,42	1.41	2 (7%)	23,68,68	1.43	5 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	STU	B	2	-	27,42,42	1.45	5 (18%)	23,68,68	1.21	1 (4%)
2	STU	C	3	-	27,42,42	1.37	4 (14%)	23,68,68	1.32	2 (8%)
2	STU	D	4	-	27,42,42	1.41	4 (14%)	23,68,68	1.24	1 (4%)

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
2	STU	A	1	-	1/1/5/7	0/4/42/42	0/0/8/8
2	STU	B	2	-	1/1/5/7	0/4/42/42	0/0/8/8
2	STU	C	3	-	1/1/5/7	0/4/42/42	0/0/8/8
2	STU	D	4	-	1/1/5/7	0/4/42/42	0/0/8/8

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	STU	C11-C18	-2.93	1.38	1.42
2	D	4	STU	C11-C18	-2.61	1.39	1.42
2	C	3	STU	C6-C19	-2.58	1.39	1.42
2	C	3	STU	C7-C8	-2.52	1.39	1.49
2	B	2	STU	C7-C8	-2.51	1.39	1.49
2	C	3	STU	C11-C18	-2.44	1.39	1.42
2	A	1	STU	C6-C19	-2.28	1.39	1.42
2	D	4	STU	C7-C8	-2.11	1.41	1.49
2	B	2	STU	C6-C19	-2.09	1.40	1.42
2	C	3	STU	C3-C4	2.01	1.41	1.36
2	D	4	STU	C3-C4	2.13	1.41	1.36
2	B	2	STU	C2-C1	2.14	1.41	1.36
2	B	2	STU	C3-C4	2.24	1.41	1.36
2	D	4	STU	O5-C8	3.15	1.29	1.23
2	A	1	STU	O5-C8	3.49	1.30	1.23

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	STU	C1-C20-N3	-3.35	128.15	132.18
2	D	4	STU	C1-C20-N3	-3.10	128.46	132.18
2	A	1	STU	C16-C17-N2	-2.76	128.86	132.22
2	B	2	STU	C1-C20-N3	-2.74	128.89	132.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	STU	C1-C20-N3	-2.57	129.09	132.18
2	C	3	STU	O5-C8-C7	-2.40	124.61	128.62
2	A	1	STU	C14-C13-C12	-2.07	117.21	120.79
2	A	1	STU	C3-C4-C5	-2.03	117.28	120.79
2	A	1	STU	C4-C5-C20	2.22	122.20	119.39

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	3	STU	C23
2	A	1	STU	C23
2	D	4	STU	C23
2	B	2	STU	C23

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	STU	6	0
2	B	2	STU	4	0
2	C	3	STU	2	0
2	D	4	STU	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 [i](#)

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.