



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 12:57 PM GMT

PDB ID : 3E8C  
Title : Crystal structures of the kinase domain of PKA in complex with ATP-competitive inhibitors  
Authors : Concha, N.O.; Elkins, P.A.; Smallwood, A.; Ward, P.  
Deposited on : 2008-08-19  
Resolution : 2.20 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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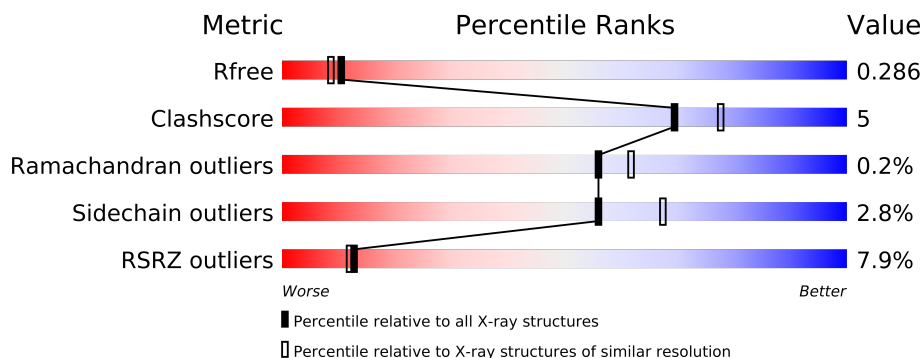
The following versions of software and data (see [references](#)) were used in the production of this report:

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 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))
$R_{free}$	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (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 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	350	
1	B	350	
1	C	350	
1	D	350	
1	E	350	
1	F	350	
2	G	20	
2	H	20	
2	I	20	
2	J	20	
2	K	20	
2	L	20	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18893 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.

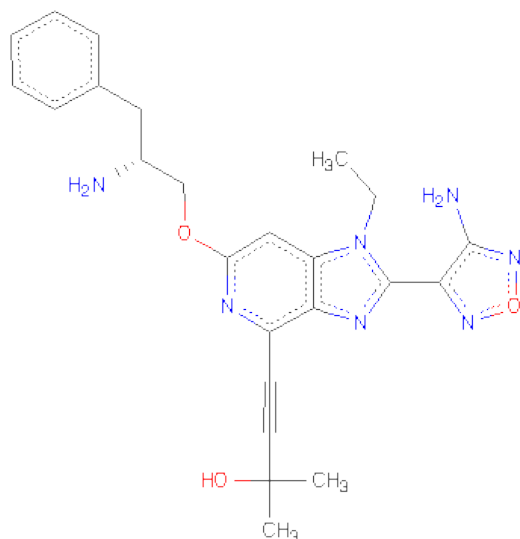
- Molecule 1 is a protein called cAMP-dependent protein kinase catalytic subunit alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	P	S	0	0	0
			2845	1838	477	518	3	9			
1	B	347	Total	C	N	O	P	S	0	0	0
			2860	1847	480	521	3	9			
1	C	344	Total	C	N	O	P	S	0	1	0
			2844	1838	477	517	3	9			
1	D	347	Total	C	N	O	P	S	0	0	0
			2842	1838	475	517	3	9			
1	E	338	Total	C	N	O	P	S	0	0	0
			2785	1804	467	503	2	9			
1	F	344	Total	C	N	O	P	S	0	0	0
			2845	1838	477	518	3	9			

- Molecule 2 is a protein called cAMP-dependent protein kinase inhibitor peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	18	Total	C	N	O	0	0	0
			138	84	28	26			
2	H	20	Total	C	N	O	0	0	0
			157	94	32	31			
2	I	20	Total	C	N	O	0	0	0
			157	94	32	31			
2	J	19	Total	C	N	O	0	0	0
			148	90	31	27			
2	K	20	Total	C	N	O	0	0	0
			157	94	32	31			
2	L	19	Total	C	N	O	0	0	0
			149	90	30	29			

- Molecule 3 is 4-[2-(4-AMINO-1,2,5-OXADIAZOL-3-YL)-6-{[(2R)-2-AMINO-3-PHENYLPROPYL]OXY}-1-ETHYL-1H-IMIDAZO[4,5-C]PYRIDIN-4-YL]-2-METHYLBUT-3-YN-2-OL (three-letter code: G96) (formula: C<sub>24</sub>H<sub>27</sub>N<sub>7</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			34	24	7	3		
3	B	1	Total	C	N	O	0	0
			34	24	7	3		
3	C	1	Total	C	N	O	0	0
			34	24	7	3		
3	D	1	Total	C	N	O	0	0
			34	24	7	3		
3	E	1	Total	C	N	O	0	0
			34	24	7	3		
3	F	1	Total	C	N	O	0	0
			34	24	7	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		
4	B	160	Total	O	0	0
			160	160		
4	C	122	Total	O	0	0
			122	122		
4	D	106	Total	O	0	0
			106	106		
4	E	95	Total	O	0	0
			95	95		
4	F	103	Total	O	0	0
			103	103		

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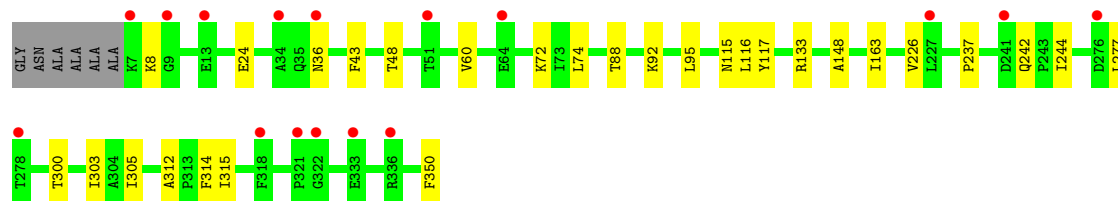
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	4	Total 4	O 4	0	0
4	H	12	Total 12	O 12	0	0
4	I	12	Total 12	O 12	0	0
4	J	6	Total 6	O 6	0	0
4	K	8	Total 8	O 8	0	0
4	L	4	Total 4	O 4	0	0

### 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.

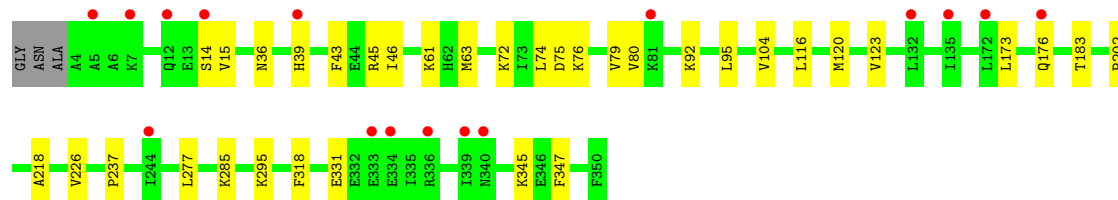
- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

Chain A: 



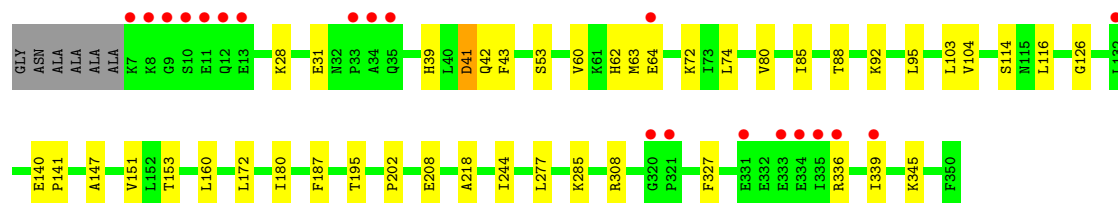
- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

Chain B: 



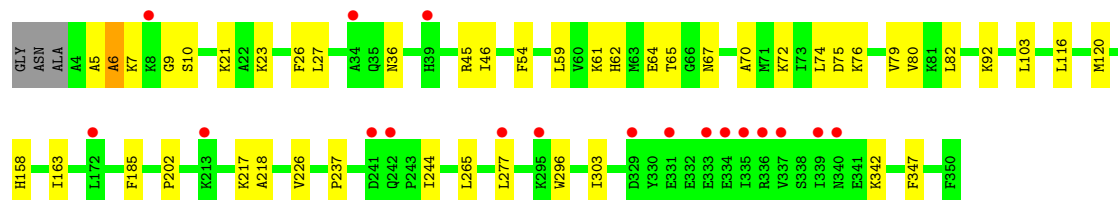
- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

Chain C: 

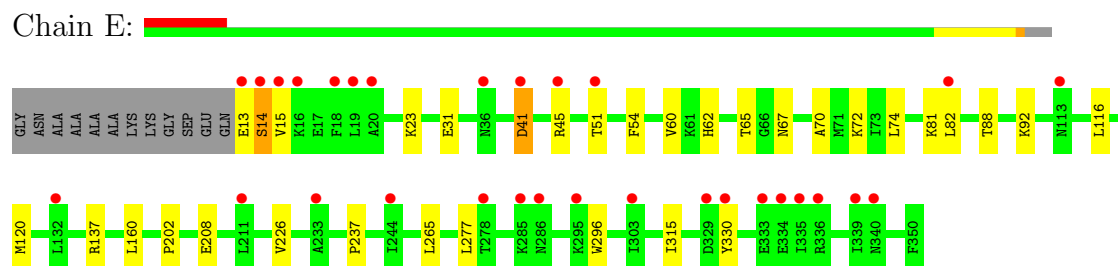


- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

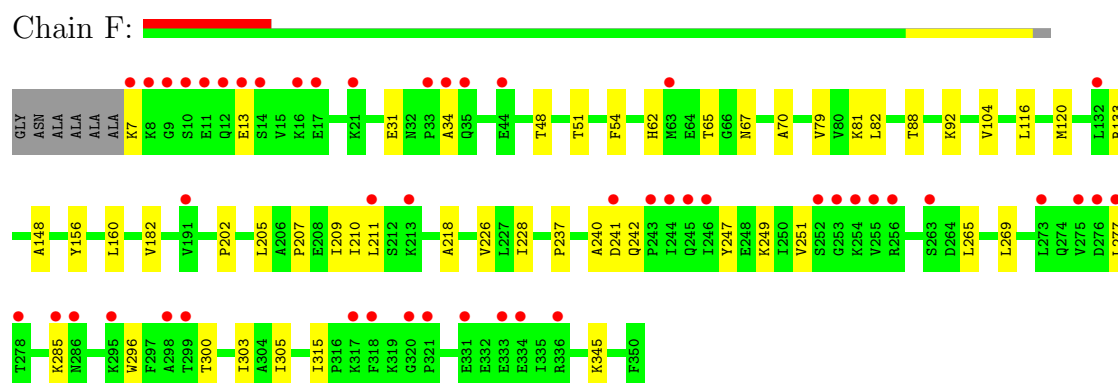
Chain D: 



- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha



- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha



- Molecule 2: cAMP-dependent protein kinase inhibitor peptide



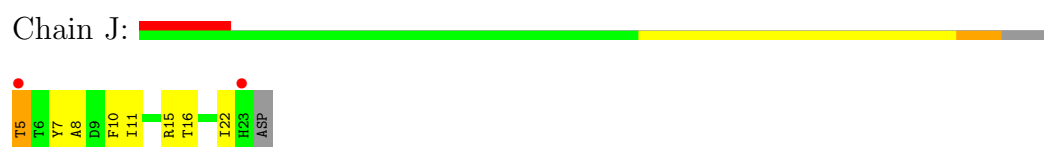
- Molecule 2: cAMP-dependent protein kinase inhibitor peptide



- Molecule 2: cAMP-dependent protein kinase inhibitor peptide

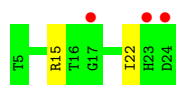


- Molecule 2: cAMP-dependent protein kinase inhibitor peptide



- Molecule 2: cAMP-dependent protein kinase inhibitor peptide

Chain K: 



- Molecule 2: cAMP-dependent protein kinase inhibitor peptide

Chain L: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.46Å 96.39Å 180.04Å 90.00° 102.66° 90.00°	Depositor
Resolution (Å)	46.93 – 2.20 43.06 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.8 (46.93-2.20) 95.8 (43.06-2.20)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.02 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.264 , 0.290 0.263 , 0.286	Depositor DCC
$R_{free}$ test set	6759 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.3	Xtriage
Anisotropy	0.772	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 32.0	EDS
Estimated twinning fraction	0.001 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	5 of 134346 reflections (0.004%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	18893	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.06 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.8220e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: TPO, G96, 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	0.38	0/2884	0.57	0/3880
1	B	0.39	0/2899	0.58	0/3901
1	C	0.38	0/2886	0.57	1/3884 (0.0%)
1	D	0.37	0/2881	0.53	0/3880
1	E	0.35	0/2835	0.53	0/3820
1	F	0.36	0/2884	0.52	0/3880
2	G	0.61	0/139	0.69	0/186
2	H	0.44	0/159	0.68	0/212
2	I	0.43	0/159	0.66	0/212
2	J	0.40	0/150	0.64	0/201
2	K	0.40	0/159	0.63	0/212
2	L	0.42	0/150	0.59	0/197
All	All	0.38	0/18185	0.55	1/24465 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	345	LYS	N-CA-CB	-5.03	101.55	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	15	VAL	Peptide

## 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	2845	0	2819	17	0
1	B	2860	0	2834	15	0
1	C	2844	0	2813	26	0
1	D	2842	0	2800	34	0
1	E	2785	0	2750	21	0
1	F	2845	0	2819	39	0
2	G	138	0	135	1	0
2	H	157	0	146	3	0
2	I	157	0	146	5	0
2	J	148	0	142	5	0
2	K	157	0	146	1	0
2	L	149	0	137	6	0
3	A	34	0	27	2	0
3	B	34	0	27	4	0
3	C	34	0	27	2	0
3	D	34	0	27	2	0
3	E	34	0	27	0	0
3	F	34	0	27	0	0
4	A	130	0	0	3	0
4	B	160	0	0	3	0
4	C	122	0	0	3	0
4	D	106	0	0	1	0
4	E	95	0	0	0	0
4	F	103	0	0	6	0
4	G	4	0	0	0	0
4	H	12	0	0	1	0
4	I	12	0	0	0	0
4	J	6	0	0	0	0
4	K	8	0	0	0	0
4	L	4	0	0	0	0
All	All	18893	0	17849	168	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 5.

All (168) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:208:GLU:HG2	1:E:277:LEU:HD21	1.58	0.84
1:A:48:THR:HG23	4:A:408:HOH:O	1.83	0.79
1:C:180:ILE:O	4:C:467:HOH:O	2.00	0.78
2:J:5:THR:HG23	2:J:8:ALA:HB3	1.68	0.75
1:F:210:ILE:CD1	1:F:251:VAL:CG1	2.67	0.72
1:F:210:ILE:HD11	1:F:251:VAL:HG12	1.69	0.72
1:D:65:THR:HG23	1:D:67:ASN:H	1.54	0.71
1:F:65:THR:HG23	1:F:67:ASN:H	1.56	0.70
2:J:5:THR:HG23	2:J:8:ALA:CB	2.22	0.70
1:D:62:HIS:ND1	1:D:65:THR:HG22	2.06	0.69
1:F:62:HIS:CD2	1:F:65:THR:HG22	2.27	0.69
1:F:210:ILE:HD11	1:F:251:VAL:CG1	2.23	0.69
1:D:5:ALA:CA	1:D:6:ALA:HB3	2.23	0.68
1:F:218:ALA:HB3	1:F:277:LEU:HD11	1.74	0.67
1:A:95:LEU:HD11	3:A:351:G96:H19	1.75	0.67
1:D:5:ALA:N	1:D:6:ALA:HB3	2.11	0.66
1:E:65:THR:HG23	1:E:67:ASN:H	1.60	0.65
1:F:218:ALA:CB	1:F:277:LEU:HD11	2.26	0.65
1:E:62:HIS:CD2	1:E:65:THR:HG22	2.32	0.64
2:H:22:ILE:O	4:H:448:HOH:O	2.15	0.64
1:E:72:LYS:HE2	1:E:74:LEU:HD11	1.79	0.63
1:F:62:HIS:HD2	1:F:65:THR:HG22	1.64	0.63
1:C:218:ALA:HB3	1:C:277:LEU:HD11	1.79	0.63
1:A:133:ARG:HD2	2:G:16:THR:O	1.99	0.62
1:D:5:ALA:HA	1:D:6:ALA:HB3	1.81	0.62
1:F:210:ILE:HD13	1:F:251:VAL:CG1	2.29	0.62
1:E:265:LEU:HD13	1:E:296:TRP:CE2	2.35	0.60
1:B:218:ALA:HB3	1:B:277:LEU:HD11	1.84	0.59
1:F:202:PRO:HD3	2:L:22:ILE:HG23	1.84	0.59
2:L:6:THR:C	2:L:7:TYR:CA	2.71	0.59
1:D:218:ALA:HB3	1:D:277:LEU:HD11	1.85	0.58
1:D:72:LYS:HE2	1:D:74:LEU:HD11	1.86	0.57
1:D:226:VAL:HG13	1:D:237:PRO:HD2	1.85	0.57
1:A:88:THR:HG21	1:A:116:LEU:CD1	2.34	0.57
1:B:202:PRO:HD3	2:H:22:ILE:HG23	1.89	0.55
2:J:10:PHE:CE2	2:J:16:THR:HG22	2.42	0.54
2:J:7:TYR:CE2	2:J:11:ILE:HD11	2.42	0.54
1:E:54:PHE:HB3	1:E:82:LEU:HD12	1.89	0.54
1:E:226:VAL:HG13	1:E:237:PRO:HD2	1.89	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:62:HIS:CE1	1:D:64:GLU:HG2	2.42	0.54
1:F:207:PRO:O	1:F:210:ILE:HD12	2.08	0.54
1:B:218:ALA:CB	1:B:277:LEU:HD11	2.38	0.54
1:D:244:ILE:H	1:D:244:ILE:HD12	1.73	0.54
1:F:205:LEU:HD22	1:F:209:ILE:HG21	1.90	0.53
1:A:244:ILE:HG23	4:A:445:HOH:O	2.08	0.53
1:C:88:THR:HG21	1:C:116:LEU:CD1	2.39	0.53
1:B:72:LYS:HE2	1:B:74:LEU:HD21	1.90	0.53
1:F:88:THR:HG21	1:F:116:LEU:CD1	2.38	0.53
1:D:23:LYS:O	1:D:27:LEU:HD23	2.10	0.52
1:F:48:THR:HG21	4:F:448:HOH:O	2.08	0.52
3:B:351:G96:H21	3:B:351:G96:N1	2.24	0.52
1:F:148:ALA:HB3	4:F:457:HOH:O	2.09	0.52
1:D:54:PHE:HB3	1:D:82:LEU:HD12	1.91	0.52
1:E:208:GLU:CG	1:E:277:LEU:HD11	2.39	0.52
1:A:24:GLU:OE2	1:C:244:ILE:HD12	2.11	0.51
1:F:300:THR:HG23	1:F:305:ILE:HD11	1.93	0.51
1:F:210:ILE:HD13	1:F:251:VAL:HG11	1.92	0.51
1:D:202:PRO:HD3	2:J:22:ILE:HG23	1.92	0.51
1:F:54:PHE:HB3	1:F:82:LEU:HD12	1.93	0.51
1:D:26:PHE:HD2	1:D:27:LEU:HD22	1.75	0.50
1:B:226:VAL:HG13	1:B:237:PRO:HD2	1.94	0.50
1:C:95:LEU:CD1	3:C:351:G96:H20	2.42	0.50
3:B:351:G96:H19A	4:B:752:HOH:O	2.11	0.50
1:F:7:LYS:NZ	1:F:303:ILE:HD13	2.27	0.49
1:D:74:LEU:HD22	1:D:74:LEU:N	2.27	0.49
1:D:120:MET:SD	3:D:351:G96:H20B	2.52	0.49
3:D:351:G96:H21	3:D:351:G96:N1	2.28	0.49
1:E:51:THR:HG22	1:E:330:TYR:CE2	2.47	0.49
1:C:103:LEU:HD21	1:C:153:THR:CG2	2.43	0.49
1:E:208:GLU:HG3	1:E:277:LEU:HD11	1.93	0.49
1:F:228:ILE:CG2	1:F:269:LEU:HD11	2.43	0.48
2:I:5:THR:O	2:I:9:ASP:OD2	2.31	0.48
4:B:516:HOH:O	2:H:21:ALA:HB2	2.14	0.48
1:D:103:LEU:HD22	1:D:185:PHE:HZ	1.77	0.48
1:C:187:PHE:CE2	2:I:21:ALA:HB1	2.49	0.48
3:A:351:G96:N1	3:A:351:G96:H21	2.29	0.48
1:F:300:THR:HG23	1:F:305:ILE:CD1	2.44	0.48
1:B:95:LEU:CD1	3:B:351:G96:H20	2.43	0.48
1:F:247:TYR:O	1:F:251:VAL:HG22	2.14	0.47
1:A:88:THR:HG21	1:A:116:LEU:HD13	1.96	0.47
1:A:300:THR:HG23	1:A:305:ILE:CD1	2.43	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:147:ALA:O	1:C:151:VAL:HG23	2.13	0.47
1:B:46:ILE:HD11	1:B:61:LYS:HB2	1.94	0.47
1:D:5:ALA:N	1:D:6:ALA:CB	2.77	0.47
1:A:303:ILE:HD13	4:A:451:HOH:O	2.14	0.47
1:F:226:VAL:HG13	1:F:237:PRO:HD2	1.96	0.47
1:B:75:ASP:O	1:B:79:VAL:HG23	2.15	0.47
1:F:265:LEU:HD13	1:F:296:TRP:CE2	2.49	0.47
1:B:116:LEU:HD11	1:B:347:PHE:CD1	2.50	0.47
1:D:62:HIS:ND1	1:D:65:THR:CG2	2.76	0.47
1:F:305:ILE:HG23	4:F:453:HOH:O	2.14	0.47
1:F:133:ARG:HD2	2:L:16:THR:O	2.15	0.47
1:C:187:PHE:HE2	2:I:21:ALA:HB1	1.79	0.47
1:A:72:LYS:HE2	1:A:74:LEU:HD11	1.97	0.47
1:B:43:PHE:C	1:B:63:MET:HE2	2.35	0.47
1:C:39:HIS:CE1	1:C:42:GLN:HG3	2.50	0.47
2:I:23:HIS:CG	2:I:24:ASP:H	2.33	0.46
1:D:218:ALA:CB	1:D:277:LEU:HD11	2.44	0.46
1:A:226:VAL:HG13	1:A:237:PRO:HD2	1.98	0.46
1:C:88:THR:HG21	1:C:116:LEU:HD13	1.97	0.46
1:D:265:LEU:HD13	1:D:296:TRP:CE2	2.50	0.46
1:A:312:ALA:HB3	1:A:315:ILE:HD11	1.98	0.46
1:F:305:ILE:HD13	4:F:457:HOH:O	2.15	0.46
1:C:103:LEU:HD21	1:C:153:THR:HG23	1.97	0.46
1:C:339:ILE:H	1:C:339:ILE:HD12	1.81	0.46
2:L:6:THR:O	2:L:6:THR:HG22	2.15	0.45
1:F:34:ALA:HB3	4:F:455:HOH:O	2.15	0.45
1:D:46:ILE:HD11	1:D:61:LYS:HB2	1.97	0.45
1:C:172:LEU:HB3	1:C:180:ILE:HD12	1.98	0.45
1:D:103:LEU:HD22	1:D:185:PHE:CZ	2.51	0.45
1:A:92:LYS:HE2	1:A:350:PHE:HA	1.98	0.45
1:C:43:PHE:C	1:C:63:MET:HE2	2.37	0.45
1:E:265:LEU:HD13	1:E:296:TRP:CZ2	2.51	0.45
1:C:202:PRO:HD3	2:I:22:ILE:HG23	1.99	0.45
1:E:202:PRO:HD3	2:K:22:ILE:HG23	1.98	0.44
1:C:80:VAL:HG22	1:C:85:ILE:HD11	2.00	0.44
1:A:43:PHE:HB3	1:A:60:VAL:HG13	1.99	0.44
1:D:75:ASP:O	1:D:79:VAL:HG23	2.17	0.44
2:L:10:PHE:CE2	2:L:16:THR:HG22	2.52	0.44
1:C:41:ASP:O	1:C:63:MET:HE3	2.18	0.44
1:D:5:ALA:CA	1:D:6:ALA:CB	2.91	0.44
1:F:207:PRO:O	1:F:210:ILE:CD1	2.65	0.44
1:E:74:LEU:N	1:E:74:LEU:HD22	2.33	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:46:ILE:HD12	1:D:59:LEU:HG	2.00	0.43
1:B:95:LEU:HD11	3:B:351:G96:H20	1.99	0.43
1:F:156:TYR:O	1:F:160:LEU:HD13	2.18	0.43
1:F:104:VAL:HG22	1:F:182:VAL:O	2.18	0.43
1:E:88:THR:HG21	1:E:116:LEU:CD1	2.48	0.43
1:B:123:VAL:HB	1:B:173:LEU:HD13	2.00	0.43
1:C:308:ARG:NH2	4:C:467:HOH:O	2.51	0.43
1:F:210:ILE:HD12	1:F:211:LEU:HG	2.01	0.43
1:F:34:ALA:HB2	4:F:378:HOH:O	2.18	0.43
1:D:5:ALA:HA	1:D:7:LYS:H	1.83	0.43
1:C:218:ALA:CB	1:C:277:LEU:HD11	2.46	0.42
1:E:208:GLU:CG	1:E:277:LEU:HD21	2.40	0.42
1:D:54:PHE:CD2	1:D:79:VAL:HG22	2.54	0.42
1:C:72:LYS:HE2	1:C:74:LEU:HD11	2.01	0.42
1:C:104:VAL:HG22	4:C:479:HOH:O	2.18	0.42
1:E:13:GLU:O	1:E:14:SER:HB2	2.19	0.42
1:E:13:GLU:O	1:E:14:SER:CB	2.68	0.42
1:F:207:PRO:HA	1:F:210:ILE:HD11	2.02	0.42
1:D:70:ALA:HB3	1:D:120:MET:HG3	2.01	0.42
1:A:148:ALA:CB	1:A:305:ILE:HD13	2.50	0.42
1:C:62:HIS:CD2	1:C:64:GLU:HG2	2.55	0.42
1:C:140:GLU:N	1:C:141:PRO:CD	2.83	0.42
1:E:315:ILE:HD12	1:E:315:ILE:N	2.35	0.42
1:B:176:GLN:NE2	1:B:318:PHE:CE2	2.87	0.42
1:E:70:ALA:HB3	1:E:120:MET:HG2	2.02	0.42
1:F:70:ALA:HB3	1:F:120:MET:HG2	2.01	0.41
1:A:314:PHE:C	1:A:315:ILE:HD12	2.41	0.41
1:D:116:LEU:HD11	1:D:347:PHE:CD1	2.55	0.41
1:F:240:ALA:HB2	1:F:249:LYS:NZ	2.35	0.41
1:D:9:GLY:HA2	1:D:10:SEP:HB3	2.01	0.41
1:D:303:ILE:HD11	4:D:438:HOH:O	2.21	0.41
2:L:11:ILE:HD13	2:L:11:ILE:N	2.36	0.41
1:F:265:LEU:HD13	1:F:296:TRP:CZ2	2.56	0.41
1:E:23:LYS:HA	1:E:160:LEU:HD21	2.03	0.41
1:E:41:ASP:OD1	1:E:41:ASP:N	2.44	0.41
3:C:351:G96:H21	3:C:351:G96:N1	2.36	0.40
1:C:126:GLY:HA2	1:C:327:PHE:CZ	2.56	0.40
1:D:76:LYS:O	1:D:80:VAL:HG23	2.21	0.40
1:D:158:HIS:ND1	1:D:217:LYS:HD2	2.36	0.40
1:F:54:PHE:CD2	1:F:79:VAL:HG22	2.56	0.40
1:F:315:ILE:HD12	1:F:315:ILE:N	2.37	0.40
1:B:76:LYS:O	1:B:80:VAL:HG23	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:208:GLU:OE1	1:C:277:LEU:HD13	2.22	0.40
1:B:104:VAL:HG21	4:B:752:HOH:O	2.21	0.40
1:A:115:ASN:HB2	1:A:117:TYR:CE1	2.57	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	339/350 (97%)	330 (97%)	9 (3%)	0	100	100
1	B	342/350 (98%)	330 (96%)	10 (3%)	2 (1%)	33	32
1	C	340/350 (97%)	331 (97%)	9 (3%)	0	100	100
1	D	342/350 (98%)	330 (96%)	11 (3%)	1 (0%)	50	53
1	E	334/350 (95%)	324 (97%)	9 (3%)	1 (0%)	50	53
1	F	339/350 (97%)	331 (98%)	8 (2%)	0	100	100
2	G	16/20 (80%)	15 (94%)	1 (6%)	0	100	100
2	H	18/20 (90%)	16 (89%)	2 (11%)	0	100	100
2	I	18/20 (90%)	15 (83%)	3 (17%)	0	100	100
2	J	17/20 (85%)	16 (94%)	1 (6%)	0	100	100
2	K	18/20 (90%)	16 (89%)	2 (11%)	0	100	100
2	L	16/20 (80%)	14 (88%)	2 (12%)	0	100	100
All	All	2139/2220 (96%)	2068 (97%)	67 (3%)	4 (0%)	56	62

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	14	SER
1	B	15	VAL
1	D	6	ALA
1	B	14	SER



### 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	301/302 (100%)	296 (98%)	5 (2%)	73	84
1	B	301/302 (100%)	291 (97%)	10 (3%)	50	60
1	C	300/302 (99%)	289 (96%)	11 (4%)	45	54
1	D	296/302 (98%)	290 (98%)	6 (2%)	68	79
1	E	294/302 (97%)	287 (98%)	7 (2%)	61	73
1	F	301/302 (100%)	292 (97%)	9 (3%)	53	64
2	G	13/15 (87%)	13 (100%)	0	100	100
2	H	15/15 (100%)	15 (100%)	0	100	100
2	I	15/15 (100%)	14 (93%)	1 (7%)	23	24
2	J	14/15 (93%)	12 (86%)	2 (14%)	5	4
2	K	15/15 (100%)	14 (93%)	1 (7%)	23	24
2	L	13/15 (87%)	12 (92%)	1 (8%)	18	18
All	All	1878/1902 (99%)	1825 (97%)	53 (3%)	56	67

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	36	ASN
1	A	163	ILE
1	A	242	GLN
1	A	277	LEU
1	B	36	ASN
1	B	39	HIS
1	B	45	ARG
1	B	92	LYS
1	B	120	MET
1	B	183	THR
1	B	285	LYS
1	B	295	LYS
1	B	331	GLU
1	B	345	LYS

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Mol	Chain	Res	Type
1	C	28	LYS
1	C	31	GLU
1	C	41	ASP
1	C	53	SER
1	C	60	VAL
1	C	92	LYS
1	C	114	SER
1	C	160	LEU
1	C	195	THR
1	C	285	LYS
1	C	336	ARG
1	D	21	LYS
1	D	36	ASN
1	D	45	ARG
1	D	92	LYS
1	D	163	ILE
1	D	342	LYS
1	E	31	GLU
1	E	41	ASP
1	E	45	ARG
1	E	60	VAL
1	E	81	LYS
1	E	92	LYS
1	E	137	ARG
1	F	13	GLU
1	F	31	GLU
1	F	51	THR
1	F	81	LYS
1	F	92	LYS
1	F	241	ASP
1	F	242	GLN
1	F	285	LYS
1	F	345	LYS
2	I	15	ARG
2	J	5	THR
2	J	15	ARG
2	K	15	ARG
2	L	24	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	39	HIS
1	B	293	ASN
1	C	286	ASN
1	E	32	ASN
1	E	62	HIS
1	E	271	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

17 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	SEP	A	10	1	9,9,10	6.25	3 (33%)	10,12,14	1.87	3 (30%)
1	TPO	A	197	1	10,10,11	5.62	1 (10%)	12,14,16	1.13	1 (8%)
1	SEP	A	338	1	9,9,10	6.21	3 (33%)	10,12,14	1.36	3 (30%)
1	SEP	B	10	1	9,9,10	6.11	3 (33%)	10,12,14	1.15	1 (10%)
1	TPO	B	197	1	10,10,11	5.52	1 (10%)	12,14,16	1.25	1 (8%)
1	SEP	B	338	1	9,9,10	6.12	3 (33%)	10,12,14	1.17	1 (10%)
1	SEP	C	10	1	9,9,10	6.16	3 (33%)	10,12,14	1.32	1 (10%)
1	TPO	C	197	1	10,10,11	5.96	1 (10%)	12,14,16	1.12	0
1	SEP	C	338	1	9,9,10	6.21	3 (33%)	10,12,14	1.21	2 (20%)
1	SEP	D	10	1	9,9,10	6.26	3 (33%)	10,12,14	1.07	0
1	TPO	D	197	1	10,10,11	5.60	1 (10%)	12,14,16	1.28	1 (8%)
1	SEP	D	338	1	9,9,10	6.07	3 (33%)	10,12,14	1.24	2 (20%)
1	TPO	E	197	1	10,10,11	5.64	1 (10%)	12,14,16	1.10	1 (8%)
1	SEP	E	338	1	9,9,10	6.13	3 (33%)	10,12,14	1.20	1 (10%)
1	SEP	F	10	1	9,9,10	6.27	3 (33%)	10,12,14	1.25	1 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	TPO	F	197	1	10,10,11	5.68	1 (10%)	12,14,16	1.14	1 (8%)
1	SEP	F	338	1	9,9,10	6.15	3 (33%)	10,12,14	1.38	2 (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
1	SEP	A	10	1	-	0/6/8/10	0/0/0/0
1	TPO	A	197	1	-	0/9/11/13	0/0/0/0
1	SEP	A	338	1	-	0/6/8/10	0/0/0/0
1	SEP	B	10	1	-	0/6/8/10	0/0/0/0
1	TPO	B	197	1	-	0/9/11/13	0/0/0/0
1	SEP	B	338	1	-	0/6/8/10	0/0/0/0
1	SEP	C	10	1	-	0/6/8/10	0/0/0/0
1	TPO	C	197	1	-	0/9/11/13	0/0/0/0
1	SEP	C	338	1	-	0/6/8/10	0/0/0/0
1	SEP	D	10	1	-	0/6/8/10	0/0/0/0
1	TPO	D	197	1	-	0/9/11/13	0/0/0/0
1	SEP	D	338	1	-	0/6/8/10	0/0/0/0
1	TPO	E	197	1	-	0/9/11/13	0/0/0/0
1	SEP	E	338	1	-	0/6/8/10	0/0/0/0
1	SEP	F	10	1	-	0/6/8/10	0/0/0/0
1	TPO	F	197	1	-	0/9/11/13	0/0/0/0
1	SEP	F	338	1	-	0/6/8/10	0/0/0/0

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	197	TPO	O-C	18.67	1.24	1.11
1	F	10	SEP	O-C	18.09	1.23	1.11
1	A	10	SEP	O-C	18.04	1.23	1.11
1	D	10	SEP	O-C	18.04	1.23	1.11
1	C	338	SEP	O-C	18.00	1.23	1.11
1	A	338	SEP	O-C	17.97	1.23	1.11
1	F	338	SEP	O-C	17.78	1.23	1.11
1	F	197	TPO	O-C	17.78	1.23	1.11
1	C	10	SEP	O-C	17.77	1.23	1.11
1	E	338	SEP	O-C	17.72	1.23	1.11
1	B	338	SEP	O-C	17.69	1.23	1.11
1	E	197	TPO	O-C	17.64	1.23	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	197	TPO	O-C	17.61	1.23	1.11
1	B	10	SEP	O-C	17.60	1.23	1.11
1	D	338	SEP	O-C	17.55	1.23	1.11
1	D	197	TPO	O-C	17.49	1.23	1.11
1	B	197	TPO	O-C	17.27	1.23	1.11
1	C	10	SEP	P-O1P	3.19	1.61	1.51
1	B	10	SEP	P-O1P	3.18	1.61	1.51
1	B	338	SEP	P-O1P	3.16	1.61	1.51
1	D	10	SEP	CA-C	3.15	1.54	1.48
1	A	10	SEP	P-O1P	3.13	1.61	1.51
1	F	10	SEP	P-O1P	3.12	1.61	1.51
1	F	338	SEP	P-O1P	3.12	1.61	1.51
1	D	10	SEP	P-O1P	3.06	1.61	1.51
1	D	338	SEP	P-O1P	3.06	1.61	1.51
1	E	338	SEP	P-O1P	3.03	1.61	1.51
1	A	338	SEP	CA-C	2.96	1.53	1.48
1	F	10	SEP	CA-C	2.93	1.53	1.48
1	C	338	SEP	P-O1P	2.90	1.60	1.51
1	B	10	SEP	CA-C	2.88	1.53	1.48
1	A	10	SEP	CA-C	2.83	1.53	1.48
1	A	338	SEP	P-O1P	2.83	1.60	1.51
1	C	338	SEP	CA-C	2.77	1.53	1.48
1	C	10	SEP	CA-C	2.76	1.53	1.48
1	F	338	SEP	CA-C	2.70	1.53	1.48
1	E	338	SEP	CA-C	2.65	1.53	1.48
1	D	338	SEP	CA-C	2.53	1.53	1.48
1	B	338	SEP	CA-C	2.48	1.52	1.48

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	SEP	C-CA-N	-4.57	109.27	113.83
1	D	197	TPO	CB-CA-N	2.99	113.78	109.60
1	F	338	SEP	OG-CB-CA	2.86	112.73	108.69
1	C	10	SEP	OG-CB-CA	2.69	112.50	108.69
1	A	338	SEP	OG-P-O1P	2.48	113.96	106.71
1	E	338	SEP	OG-CB-CA	2.39	112.07	108.69
1	F	10	SEP	P-OG-CB	-2.32	111.48	118.19
1	B	10	SEP	P-OG-CB	-2.31	111.50	118.19
1	D	338	SEP	OG-CB-CA	2.29	111.93	108.69
1	A	10	SEP	OG-CB-CA	2.25	111.88	108.69
1	A	197	TPO	CB-CA-N	2.22	112.70	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	338	SEP	OG-CB-CA	2.20	111.81	108.69
1	F	197	TPO	O3P-P-O2P	2.20	116.17	107.61
1	A	10	SEP	P-OG-CB	-2.18	111.89	118.19
1	C	338	SEP	C-CA-N	-2.17	111.67	113.83
1	B	338	SEP	OG-CB-CA	2.16	111.75	108.69
1	B	197	TPO	CB-CA-N	2.14	112.58	109.60
1	A	338	SEP	C-CA-N	-2.06	111.78	113.83
1	D	338	SEP	C-CA-N	-2.05	111.78	113.83
1	C	338	SEP	OG-P-O1P	2.04	112.70	106.71
1	E	197	TPO	O3P-P-O2P	2.04	115.56	107.61
1	F	338	SEP	C-CA-N	-2.01	111.82	113.83

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 ⓘ

6 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$
3	G96	A	351	-	35,37,37	2.04	4 (11%)	48,53,53	3.83	11 (22%)
3	G96	B	351	-	35,37,37	2.13	4 (11%)	48,53,53	3.01	11 (22%)
3	G96	C	351	-	35,37,37	2.06	4 (11%)	48,53,53	3.18	10 (20%)
3	G96	D	351	-	35,37,37	2.09	4 (11%)	48,53,53	3.12	9 (18%)
3	G96	E	351	-	35,37,37	2.16	4 (11%)	48,53,53	3.49	12 (25%)
3	G96	F	351	-	35,37,37	2.00	4 (11%)	48,53,53	3.57	11 (22%)

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	G96	A	351	-	-	0/13/21/21	0/1/4/4
3	G96	B	351	-	-	0/13/21/21	0/1/4/4
3	G96	C	351	-	-	0/13/21/21	0/1/4/4
3	G96	D	351	-	-	0/13/21/21	0/1/4/4
3	G96	E	351	-	-	0/13/21/21	0/1/4/4
3	G96	F	351	-	-	0/13/21/21	0/1/4/4

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	351	G96	C12-C13	-10.77	1.28	1.43
3	B	351	G96	C12-C13	-10.71	1.28	1.43
3	D	351	G96	C12-C13	-10.44	1.28	1.43
3	C	351	G96	C12-C13	-10.10	1.29	1.43
3	A	351	G96	C12-C13	-9.98	1.29	1.43
3	F	351	G96	C12-C13	-9.79	1.29	1.43
3	A	351	G96	C15-N5	-3.78	1.34	1.38
3	C	351	G96	C15-N5	-3.64	1.34	1.38
3	B	351	G96	C15-N5	-3.44	1.34	1.38
3	D	351	G96	C15-N5	-3.35	1.34	1.38
3	E	351	G96	C15-N5	-3.33	1.34	1.38
3	F	351	G96	C15-N5	-3.21	1.34	1.38
3	F	351	G96	C9-C16	-3.15	1.40	1.48
3	A	351	G96	C9-C16	-3.11	1.40	1.48
3	C	351	G96	C9-C16	-3.06	1.40	1.48
3	E	351	G96	C9-C16	-2.99	1.41	1.48
3	D	351	G96	C9-C16	-2.92	1.41	1.48
3	B	351	G96	C9-C16	-2.78	1.41	1.48
3	A	351	G96	C9-C8	-2.69	1.39	1.44
3	C	351	G96	C9-C8	-2.63	1.39	1.44
3	F	351	G96	C9-C8	-2.57	1.39	1.44
3	E	351	G96	C9-C8	-2.53	1.39	1.44
3	B	351	G96	C9-C8	-2.46	1.39	1.44
3	D	351	G96	C9-C8	-2.36	1.40	1.44

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	351	G96	N3-C8-N2	21.69	122.49	119.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	351	G96	N3-C8-N2	20.37	122.31	119.50
3	E	351	G96	N3-C8-N2	19.82	122.23	119.50
3	C	351	G96	N3-C8-N2	17.12	121.86	119.50
3	D	351	G96	N3-C8-N2	15.28	121.61	119.50
3	B	351	G96	N3-C8-N2	14.74	121.53	119.50
3	D	351	G96	C12-C14-N6	8.77	133.66	130.78
3	A	351	G96	C12-C14-N6	8.33	133.51	130.78
3	B	351	G96	C12-C14-N6	8.25	133.49	130.78
3	F	351	G96	C12-N4-C10	7.61	121.78	115.81
3	A	351	G96	C12-N4-C10	7.51	121.70	115.81
3	D	351	G96	C12-N4-C10	6.98	121.28	115.81
3	C	351	G96	C12-N4-C10	6.97	121.28	115.81
3	C	351	G96	C12-C14-N6	6.89	133.04	130.78
3	B	351	G96	C12-N4-C10	6.52	120.93	115.81
3	E	351	G96	C12-N4-C10	6.31	120.76	115.81
3	E	351	G96	C12-C14-N6	6.15	132.80	130.78
3	F	351	G96	C12-C14-N6	6.11	132.79	130.78
3	F	351	G96	C8-C9-N1	5.08	111.49	107.56
3	E	351	G96	C7-O1-C10	5.05	122.91	116.83
3	E	351	G96	C8-C9-N1	4.89	111.34	107.56
3	A	351	G96	C8-C9-N1	4.70	111.19	107.56
3	D	351	G96	C8-C9-N1	4.62	111.14	107.56
3	C	351	G96	C8-C9-N1	4.57	111.10	107.56
3	B	351	G96	C7-O1-C10	4.29	121.99	116.83
3	D	351	G96	C7-O1-C10	4.19	121.87	116.83
3	B	351	G96	C8-C9-N1	4.02	110.67	107.56
3	D	351	G96	C21-N5-C16	3.96	131.46	126.69
3	B	351	G96	C21-N5-C16	3.69	131.13	126.69
3	F	351	G96	C7-O1-C10	3.34	120.85	116.83
3	A	351	G96	C21-N5-C16	3.32	130.69	126.69
3	D	351	G96	C11-C10-N4	-3.24	120.13	124.03
3	F	351	G96	C11-C10-N4	-3.20	120.17	124.03
3	B	351	G96	C11-C10-N4	-3.17	120.20	124.03
3	C	351	G96	C21-N5-C16	3.05	130.36	126.69
3	E	351	G96	C11-C10-N4	-3.03	120.37	124.03
3	C	351	G96	C11-C10-N4	-2.93	120.50	124.03
3	A	351	G96	C11-C10-N4	-2.87	120.56	124.03
3	A	351	G96	C7-O1-C10	2.80	120.20	116.83
3	B	351	G96	C9-C8-N2	2.77	113.46	111.43
3	C	351	G96	C7-O1-C10	2.76	120.15	116.83
3	F	351	G96	C21-N5-C16	2.75	130.00	126.69
3	E	351	G96	C21-N5-C16	2.70	129.94	126.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	351	G96	C14-C12-C13	2.66	123.73	119.22
3	E	351	G96	C1-C24-C23	-2.65	108.39	113.28
3	F	351	G96	C16-N5-C15	-2.64	104.40	107.80
3	A	351	G96	C16-N5-C15	-2.59	104.47	107.80
3	C	351	G96	C9-C8-N2	2.59	113.33	111.43
3	C	351	G96	C16-N5-C15	-2.55	104.52	107.80
3	A	351	G96	C9-C8-N2	2.54	113.29	111.43
3	E	351	G96	C16-N5-C15	-2.53	104.55	107.80
3	D	351	G96	C16-N5-C15	-2.52	104.56	107.80
3	F	351	G96	C9-C8-N2	2.44	113.22	111.43
3	B	351	G96	C1-C24-C23	-2.41	108.84	113.28
3	E	351	G96	C9-C16-N6	-2.41	121.09	125.47
3	A	351	G96	C1-C24-C23	-2.40	108.86	113.28
3	D	351	G96	C9-C8-N2	2.38	113.18	111.43
3	E	351	G96	C9-C8-N2	2.35	113.16	111.43
3	F	351	G96	C9-C16-N6	-2.35	121.19	125.47
3	B	351	G96	C9-C16-N6	-2.21	121.44	125.47
3	F	351	G96	C1-C24-C23	-2.18	109.25	113.28
3	B	351	G96	C16-N5-C15	-2.16	105.02	107.80
3	E	351	G96	C11-C15-C14	-2.12	118.55	120.55
3	C	351	G96	C1-C24-C23	-2.08	109.44	113.28

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(Å²)	Q<0.9
1	A	344/350 (98%)	0.52	16 (4%)	30	30	14, 16, 26, 48	0
1	B	347/350 (99%)	0.37	16 (4%)	31	31	14, 16, 27, 35	0
1	C	344/350 (98%)	0.66	20 (5%)	22	22	14, 16, 26, 53	0
1	D	347/350 (99%)	0.52	18 (5%)	26	26	14, 16, 24, 31	0
1	E	338/350 (96%)	0.68	30 (8%)	10	9	14, 16, 22, 31	0
1	F	344/350 (98%)	0.88	49 (14%)	3	3	14, 16, 26, 39	0
2	G	18/20 (90%)	0.89	2 (11%)	6	5	20, 21, 29, 32	0
2	H	20/20 (100%)	1.04	2 (10%)	8	7	20, 22, 37, 38	0
2	I	20/20 (100%)	1.70	4 (20%)	2	1	20, 22, 37, 38	0
2	J	19/20 (95%)	1.09	2 (10%)	7	6	20, 22, 32, 35	0
2	K	20/20 (100%)	1.42	3 (15%)	3	3	20, 22, 35, 36	0
2	L	19/20 (95%)	3.42	11 (57%)	0	0	20, 22, 35, 36	0
All	All	2180/2220 (98%)	0.66	173 (7%)	13	12	14, 16, 27, 53	0

All (173) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	24	ASP	10.7
1	F	9	GLY	10.4
1	C	9	GLY	9.2
2	L	23	HIS	9.1
2	I	24	ASP	8.9
2	K	23	HIS	7.8
2	I	23	HIS	7.7
2	H	24	ASP	7.3
1	F	13	GLU	7.3
2	K	24	ASP	7.0
2	L	7	TYR	6.1

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Mol	Chain	Res	Type	RSRZ
2	I	22	ILE	5.9
2	L	22	ILE	5.9
2	H	23	HIS	5.5
2	G	14	GLY	5.3
1	C	11	GLU	5.3
1	C	7	LYS	5.2
1	E	16	LYS	5.2
1	E	339	ILE	5.1
2	J	23	HIS	5.1
2	L	6	THR	5.0
1	E	19	LEU	4.9
1	E	15	VAL	4.9
1	F	321	PRO	4.8
1	F	11	GLU	4.8
1	E	18	PHE	4.7
1	A	9	GLY	4.6
1	E	336	ARG	4.5
1	C	8	LYS	4.5
1	F	34	ALA	4.5
2	L	13	SER	4.3
1	E	335	ILE	4.2
1	B	339	ILE	4.1
1	A	321	PRO	4.0
2	L	21	ALA	4.0
1	A	336	ARG	4.0
1	C	34	ALA	3.9
1	F	253	GLY	3.9
1	E	334	GLU	3.9
1	A	34	ALA	3.9
1	C	35	GLN	3.8
1	F	336	ARG	3.8
1	C	336	ARG	3.8
1	C	33	PRO	3.8
1	E	41	ASP	3.7
2	L	14	GLY	3.7
1	F	277	LEU	3.7
1	F	246	ILE	3.6
1	B	14	SER	3.6
1	C	13	GLU	3.4
1	C	64	GLU	3.4
1	F	318	PHE	3.4
1	F	63	MET	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	14	SER	3.4
2	J	5	THR	3.4
1	D	213	LYS	3.3
1	B	340	ASN	3.3
1	F	7	LYS	3.3
1	D	8	LYS	3.3
1	F	273	LEU	3.3
2	L	12	ALA	3.3
1	A	318	PHE	3.2
1	F	255	VAL	3.2
1	C	10	SEP	3.1
2	G	12	ALA	3.1
1	B	7	LYS	3.1
1	E	340	ASN	3.0
1	C	12	GLN	3.0
1	F	263	SER	3.0
1	E	333	GLU	3.0
1	F	285	LYS	3.0
1	B	39	HIS	3.0
1	F	298	ALA	2.9
1	F	12	GLN	2.9
1	D	336	ARG	2.9
1	F	211	LEU	2.9
1	D	335	ILE	2.9
1	F	317	LYS	2.9
1	A	64	GLU	2.9
1	F	243	PRO	2.9
1	F	295	LYS	2.9
1	C	334	GLU	2.9
1	B	336	ARG	2.9
1	F	132	LEU	2.8
1	D	241	ASP	2.8
1	B	244	ILE	2.8
1	A	13	GLU	2.8
1	F	213	LYS	2.8
1	D	333	GLU	2.8
1	C	331	GLU	2.8
1	F	17	GLU	2.8
2	L	10	PHE	2.8
1	F	14	SER	2.8
1	D	329	ASP	2.7
1	E	36	ASN	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	241	ASP	2.7
1	D	34	ALA	2.7
1	F	256	ARG	2.7
2	I	20	ASN	2.7
1	E	244	ILE	2.7
1	E	286	ASN	2.7
1	B	334	GLU	2.6
1	E	82	LEU	2.6
1	D	339	ILE	2.6
1	F	254	LYS	2.6
2	K	17	GLY	2.6
1	B	12	GLN	2.6
1	A	322	GLY	2.6
1	F	320	GLY	2.6
1	F	16	LYS	2.6
2	L	20	ASN	2.6
1	D	295	LYS	2.5
1	C	321	PRO	2.5
1	B	132	LEU	2.5
1	D	340	ASN	2.5
1	F	35	GLN	2.4
1	F	278	THR	2.4
1	F	299	THR	2.4
1	F	245	GLN	2.4
1	F	334	GLU	2.4
1	E	295	LYS	2.4
1	C	320	GLY	2.4
1	F	21	LYS	2.4
1	C	132	LEU	2.3
1	E	51	THR	2.3
1	C	333	GLU	2.3
1	C	339	ILE	2.3
1	E	45	ARG	2.3
1	E	211	LEU	2.3
1	E	233	ALA	2.3
1	F	252	SER	2.3
1	B	176	GLN	2.3
1	B	135	ILE	2.3
1	E	285	LYS	2.3
1	E	303	ILE	2.3
1	A	278	THR	2.3
1	E	20	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	33	PRO	2.3
1	A	333	GLU	2.2
1	D	334	GLU	2.2
1	A	241	ASP	2.2
1	A	7	LYS	2.2
1	C	335	ILE	2.2
1	A	227	LEU	2.2
1	D	172	LEU	2.2
1	F	286	ASN	2.2
1	B	81	LYS	2.2
1	E	113	ASN	2.2
1	B	333	GLU	2.2
1	D	331	GLU	2.2
1	B	172	LEU	2.2
1	F	276	ASP	2.1
1	D	39	HIS	2.1
1	F	191	VAL	2.1
1	A	276	ASP	2.1
1	E	329	ASP	2.1
1	F	333	GLU	2.1
1	D	337	VAL	2.1
1	F	44	GLU	2.1
1	F	331	GLU	2.1
1	E	278	THR	2.1
1	A	36	ASN	2.1
1	E	330	TYR	2.1
1	F	244	ILE	2.1
1	B	5	ALA	2.1
1	D	277	LEU	2.0
1	A	51	THR	2.0
1	E	13	GLU	2.0
1	F	10	SEP	2.0
1	D	242	GLN	2.0
1	F	8	LYS	2.0
1	E	132	LEU	2.0
1	F	275	VAL	2.0

## 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	SEP	D	10	10/11	0.18	0.00	22,23,23,24	0
1	SEP	C	10	10/11	0.46	-0.04	49,50,52,52	0
1	SEP	A	10	10/11	0.27	-0.21	45,46,46,46	0
1	SEP	F	10	10/11	0.36	-0.27	37,38,38,38	0
1	SEP	C	338	10/11	0.16	-0.68	25,27,27,27	0
1	SEP	B	10	10/11	0.15	-0.69	30,31,32,32	0
1	TPO	A	197	11/12	0.13	-0.69	11,13,15,15	0
1	SEP	F	338	10/11	0.10	-0.86	25,26,27,27	0
1	TPO	F	197	11/12	0.14	-0.87	12,14,15,15	0
1	SEP	A	338	10/11	0.13	-0.88	25,26,27,27	0
1	SEP	B	338	10/11	0.17	-0.95	26,27,27,27	0
1	TPO	D	197	11/12	0.13	-1.24	11,13,14,15	0
1	TPO	C	197	11/12	0.13	-1.24	11,13,15,15	0
1	TPO	B	197	11/12	0.11	-1.29	10,13,14,15	0
1	SEP	D	338	10/11	0.16	-1.30	26,26,27,27	0
1	TPO	E	197	11/12	0.10	-1.74	11,13,15,15	0
1	SEP	E	338	10/11	0.14	-2.05	26,26,27,27	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
3	G96	A	351	34/34	0.16	-0.22	4,5,7,7	0
3	G96	F	351	34/34	0.14	-0.38	13,14,17,17	0
3	G96	C	351	34/34	0.15	-0.43	10,11,12,13	0
3	G96	E	351	34/34	0.14	-0.56	13,14,19,20	0
3	G96	B	351	34/34	0.14	-0.62	5,6,12,12	0
3	G96	D	351	34/34	0.13	-0.63	8,10,15,15	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.