



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:48 PM GMT

PDB ID : 1CM8
Title : PHOSPHORYLATED MAP KINASE P38-GAMMA
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Deposited on : 1999-05-17
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

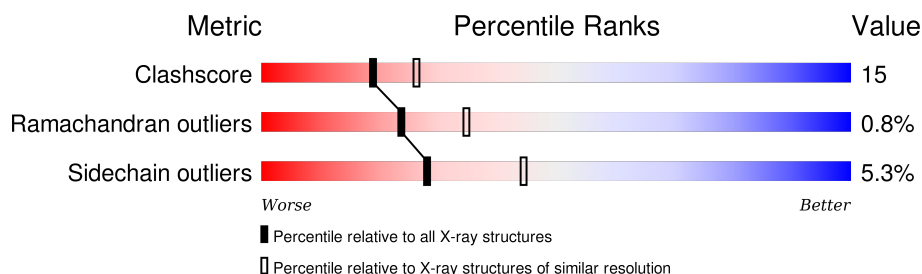
1 Overall quality at a glance

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	367	
1	B	367	

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
1	TPO	A	183	X	-	-	-
1	TPO	B	1183	X	-	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5442 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 PHOSPHORYLATED MAP KINASE P38-GAMMA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	P	S	0	0	0
			2595	1661	432	485	2	15			
1	B	329	Total	C	N	O	P	S	0	0	0
			2595	1661	432	485	2	15			

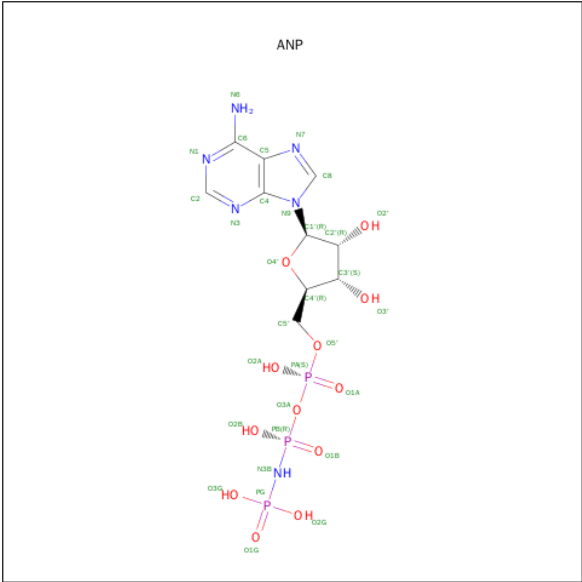
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	183	TPO	THR	MODIFIED RESIDUE	UNP P53778
B	1183	TPO	THR	MODIFIED RESIDUE	UNP P53778
A	185	PTR	TYR	MODIFIED RESIDUE	UNP P53778
B	1185	PTR	TYR	MODIFIED RESIDUE	UNP P53778

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mg	0	0
			2	2		
2	A	2	Total	Mg	0	0
			2	2		

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
3	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

- Molecule 4 is water.

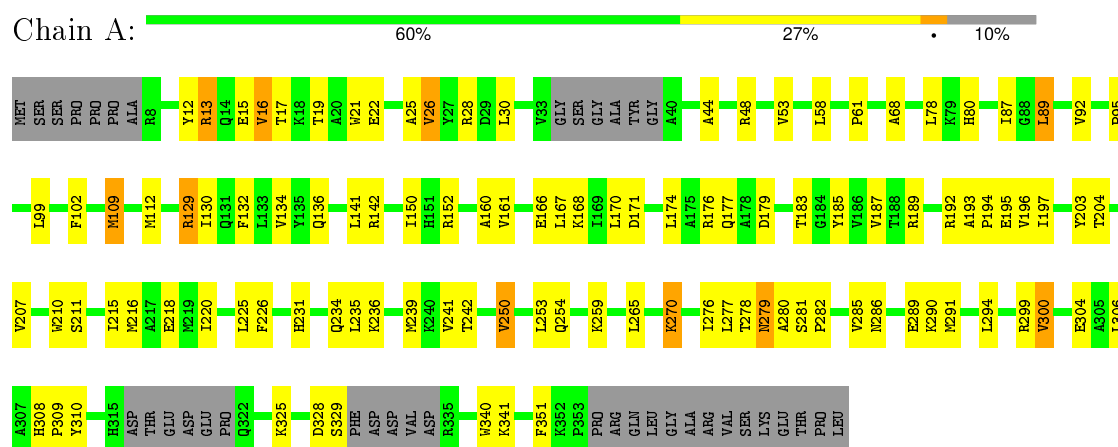
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	75	Total	O	0	0
			75	75		
4	B	111	Total	O	0	0
			111	111		

3 Residue-property plots

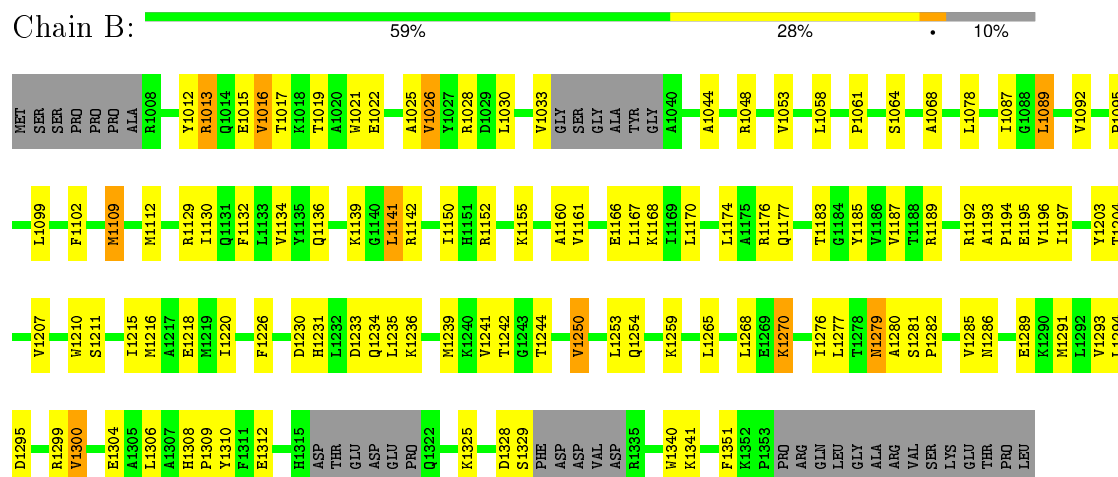
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

• Molecule 1: PHOSPHORYLATED MAP KINASE P38-GAMMA



• Molecule 1: PHOSPHORYLATED MAP KINASE P38-GAMMA



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.50 Å 66.82 Å 206.02 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40	Depositor
% Data completeness (in resolution range)	79.3 (50.00-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	6.70	Depositor
Refinement program	CNS 0.2	Depositor
R, R_{free}	0.232 , 0.283	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5442	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TPO, MG, ANP, PTR

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	0.49	0/2618	0.96	6/3542 (0.2%)
1	B	0.48	0/2618	0.97	6/3542 (0.2%)
All	All	0.49	0/5236	0.97	12/7084 (0.2%)

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	A	1	0
1	B	1	0
All	All	2	0

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1013	ARG	NE-CZ-NH2	-20.06	110.27	120.30
1	A	13	ARG	NE-CZ-NH2	-19.65	110.48	120.30
1	B	1129	ARG	NE-CZ-NH2	-18.95	110.82	120.30
1	B	1013	ARG	NE-CZ-NH1	18.64	129.62	120.30
1	A	129	ARG	NE-CZ-NH1	18.61	129.60	120.30
1	B	1129	ARG	NE-CZ-NH1	18.53	129.56	120.30
1	A	129	ARG	NE-CZ-NH2	-18.36	111.12	120.30
1	A	13	ARG	NE-CZ-NH1	17.22	128.91	120.30
1	B	1013	ARG	CD-NE-CZ	11.30	139.42	123.60
1	A	13	ARG	CD-NE-CZ	11.26	139.37	123.60
1	A	129	ARG	CD-NE-CZ	11.05	139.08	123.60
1	B	1129	ARG	CD-NE-CZ	10.95	138.93	123.60

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	183	TPO	CB
1	B	1183	TPO	CB

There are no planarity outliers.

5.2 Too-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 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	2595	0	2513	79	0
1	B	2595	0	2513	82	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	31	0	13	0	0
3	B	31	0	13	0	0
4	A	75	0	0	6	0
4	B	111	0	0	9	0
All	All	5442	0	5052	158	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 15.

All (158) 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:1279:ASN:H	1:B:1279:ASN:HD22	1.31	0.79
1:A:279:ASN:HD22	1:A:279:ASN:H	1.32	0.78
1:B:1134:VAL:HG21	1:B:1216:MET:HG3	1.69	0.75
1:A:134:VAL:HG21	1:A:216:MET:HG3	1.69	0.74
1:B:1152:ARG:HH22	1:B:1183:TPO:P	2.10	0.73
1:A:13:ARG:HG2	1:A:22:GLU:HG3	1.71	0.73
1:B:1078:LEU:HB3	1:B:1089:LEU:HD22	1.72	0.71
1:A:152:ARG:HH22	1:A:183:TPO:P	2.14	0.69
1:A:30:LEU:HD23	1:A:44:ALA:HB2	1.76	0.68
1:A:78:LEU:HB3	1:A:89:LEU:HD22	1.73	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:ASP:HB3	4:A:2104:HOH:O	1.94	0.67
1:B:1291:MET:O	1:B:1299:ARG:HD2	1.94	0.67
1:B:1030:LEU:HD23	1:B:1044:ALA:HB2	1.77	0.67
1:A:112:MET:SD	1:A:168:LYS:HD2	2.34	0.67
1:B:1235:LEU:O	1:B:1239:MET:HG2	1.95	0.67
1:B:1239:MET:SD	4:B:2170:HOH:O	2.52	0.66
1:B:1277:LEU:HD12	1:B:1285:VAL:HG22	1.77	0.66
1:A:160:ALA:HB2	1:A:170:LEU:HD11	1.78	0.66
1:A:235:LEU:O	1:A:239:MET:HG2	1.96	0.65
1:B:1112:MET:SD	1:B:1168:LYS:HD2	2.38	0.64
1:A:277:LEU:HD12	1:A:285:VAL:HG22	1.78	0.64
1:A:291:MET:O	1:A:299:ARG:HD2	1.97	0.63
1:B:1130:ILE:O	1:B:1134:VAL:HG23	1.98	0.63
1:B:1109:MET:HE1	4:B:2002:HOH:O	1.98	0.63
1:A:185:PTR:P	1:A:192:ARG:HH22	2.21	0.63
1:B:1139:LYS:HD2	4:B:2143:HOH:O	1.98	0.62
1:B:1013:ARG:HG2	1:B:1022:GLU:HG3	1.81	0.62
1:A:290:LYS:HE3	1:B:1312:GLU:OE2	2.00	0.62
1:A:282:PRO:HB2	1:B:1282:PRO:HB3	1.80	0.62
1:A:130:ILE:O	1:A:134:VAL:HG23	2.00	0.62
1:A:92:VAL:HG12	1:A:351:PHE:CD2	2.35	0.62
1:B:1142:ARG:HH21	1:B:1306:LEU:HD13	1.65	0.61
1:B:1092:VAL:HG12	1:B:1351:PHE:CD2	2.35	0.61
1:B:1185:PTR:P	1:B:1192:ARG:HH22	2.22	0.61
1:A:26:VAL:O	1:A:28:ARG:HG3	2.01	0.60
1:A:142:ARG:HH21	1:A:306:LEU:HD13	1.65	0.60
1:B:1160:ALA:HB2	1:B:1170:LEU:HD11	1.82	0.60
1:B:1204:THR:O	1:B:1207:VAL:HG13	2.01	0.59
1:B:1026:VAL:O	1:B:1028:ARG:HG3	2.03	0.58
1:A:308:HIS:HD1	1:A:310:TYR:H	1.50	0.58
1:A:195:GLU:CD	1:A:299:ARG:HH22	2.06	0.58
1:B:1308:HIS:HD1	1:B:1310:TYR:H	1.51	0.57
1:A:102:PHE:O	1:A:341:LYS:HE3	2.05	0.57
1:B:1195:GLU:CD	1:B:1299:ARG:HH22	2.07	0.57
1:B:1102:PHE:O	1:B:1341:LYS:HE3	2.04	0.57
1:A:195:GLU:OE2	1:A:299:ARG:NH2	2.38	0.56
1:A:226:PHE:HE1	1:A:241:VAL:HG21	1.71	0.56
1:B:1026:VAL:CG2	1:B:1048:ARG:HG3	2.37	0.55
1:B:1226:PHE:HE1	1:B:1241:VAL:HG21	1.72	0.54
1:B:1279:ASN:ND2	1:B:1279:ASN:H	2.02	0.54
1:B:1195:GLU:OE2	1:B:1299:ARG:NH2	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:PHE:O	1:A:136:GLN:HG3	2.08	0.54
1:B:1236:LYS:HD3	1:B:1265:LEU:HD11	1.89	0.54
1:A:300:VAL:HG22	1:A:304:GLU:HB2	1.89	0.54
1:B:1152:ARG:O	1:B:1174:LEU:HD23	2.08	0.54
1:A:236:LYS:HD3	1:A:265:LEU:HD11	1.90	0.54
1:A:150:ILE:HG23	1:A:152:ARG:HD2	1.90	0.53
1:A:152:ARG:O	1:A:174:LEU:HD23	2.08	0.53
1:A:211:SER:O	1:A:215:ILE:HG13	2.07	0.53
1:B:1211:SER:O	1:B:1215:ILE:HG13	2.08	0.53
1:B:1203:TYR:HA	4:B:2102:HOH:O	2.09	0.53
1:B:1194:PRO:HG3	1:B:1294:LEU:HD12	1.91	0.53
1:A:279:ASN:HD22	1:A:279:ASN:N	2.00	0.53
1:B:1300:VAL:HG22	1:B:1304:GLU:HB2	1.90	0.53
1:A:204:THR:O	1:A:207:VAL:HG13	2.07	0.53
1:B:1192:ARG:HD2	1:B:1196:VAL:HG11	1.91	0.53
1:A:26:VAL:CG2	1:A:48:ARG:HG3	2.39	0.52
1:B:1132:PHE:O	1:B:1136:GLN:HG3	2.10	0.52
1:A:279:ASN:ND2	1:A:279:ASN:H	2.04	0.51
1:A:194:PRO:HG3	1:A:294:LEU:HD12	1.93	0.51
1:B:1197:ILE:HD11	1:B:1231:HIS:HB2	1.93	0.50
1:B:1150:ILE:HG23	1:B:1152:ARG:HD2	1.92	0.50
1:A:250:VAL:O	1:A:259:LYS:HG3	2.13	0.49
1:B:1189:ARG:HG2	1:B:1234:GLN:OE1	2.13	0.49
1:A:176:ARG:NH1	1:A:177:GLN:O	2.45	0.48
1:B:1176:ARG:NH1	1:B:1177:GLN:O	2.47	0.48
1:B:1142:ARG:HE	1:B:1306:LEU:HD12	1.79	0.48
1:A:189:ARG:HG2	1:A:234:GLN:OE1	2.14	0.48
1:B:1242:THR:O	1:B:1270:LYS:HG2	2.14	0.48
1:A:242:THR:O	1:A:270:LYS:HG2	2.13	0.48
1:A:197:ILE:HD11	1:A:231:HIS:HB2	1.96	0.48
1:A:142:ARG:HE	1:A:306:LEU:HD12	1.80	0.47
1:A:278:THR:HA	4:A:2057:HOH:O	2.14	0.47
1:A:68:ALA:HB1	1:A:340:TRP:HB3	1.97	0.47
1:A:328:ASP:CG	1:A:329:SER:H	2.17	0.47
1:B:1068:ALA:HB1	1:B:1340:TRP:HB3	1.97	0.47
1:B:1250:VAL:O	1:B:1259:LYS:HG3	2.14	0.47
1:A:192:ARG:HD2	1:A:196:VAL:HG11	1.96	0.47
1:B:1328:ASP:CG	1:B:1329:SER:H	2.18	0.46
1:A:185:PTR:O3P	1:A:192:ARG:NH2	2.48	0.46
1:A:286:ASN:ND2	1:A:308:HIS:NE2	2.57	0.46
1:B:1015:GLU:HG3	1:B:1015:GLU:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:GLU:O	1:A:15:GLU:HG3	2.16	0.46
1:A:220:ILE:HG21	1:A:280:ALA:HB2	1.98	0.46
1:B:1220:ILE:HG21	1:B:1280:ALA:HB2	1.97	0.45
1:B:1150:ILE:HG12	1:B:1152:ARG:HG3	1.98	0.45
1:A:281:SER:O	1:A:285:VAL:HG23	2.16	0.45
1:A:142:ARG:HE	1:A:306:LEU:CD1	2.29	0.45
1:B:1064:SER:HB3	4:B:2033:HOH:O	2.15	0.45
1:A:309:PRO:HD3	1:B:1309:PRO:HG2	1.98	0.45
1:B:1286:ASN:ND2	1:B:1308:HIS:NE2	2.57	0.44
1:A:12:TYR:HD1	1:A:25:ALA:HA	1.81	0.44
1:B:1078:LEU:HB3	1:B:1089:LEU:CD2	2.45	0.44
1:A:99:LEU:HD12	1:A:99:LEU:O	2.17	0.44
1:B:1142:ARG:HE	1:B:1306:LEU:CD1	2.30	0.44
1:A:80:HIS:HB2	4:A:2058:HOH:O	2.17	0.44
1:A:161:VAL:HA	1:A:166:GLU:O	2.18	0.44
1:B:1185:PTR:O3P	1:B:1192:ARG:NH2	2.51	0.44
1:A:241:VAL:HG13	1:A:276:ILE:HD11	2.00	0.44
1:B:1236:LYS:HE3	4:B:2156:HOH:O	2.18	0.43
1:B:1016:VAL:CG2	1:B:1021:TRP:CD1	3.02	0.43
1:A:225:LEU:N	4:A:2172:HOH:O	2.51	0.43
1:A:160:ALA:O	1:A:167:LEU:HA	2.17	0.43
1:B:1241:VAL:HG13	1:B:1276:ILE:HD11	1.99	0.43
1:B:1250:VAL:HA	1:B:1253:LEU:HG	2.00	0.43
1:A:87:ILE:HG12	1:A:109:MET:HE3	2.00	0.43
1:B:1022:GLU:HB3	1:B:1095:PRO:HG3	2.01	0.43
1:B:1099:LEU:HD12	1:B:1099:LEU:O	2.18	0.43
1:A:250:VAL:HA	1:A:253:LEU:HG	2.01	0.43
1:A:152:ARG:CZ	1:A:176:ARG:HD2	2.48	0.42
1:B:1087:ILE:HG12	1:B:1109:MET:HE3	2.01	0.42
1:A:22:GLU:HB3	1:A:95:PRO:HG3	2.01	0.42
1:B:1281:SER:O	1:B:1285:VAL:HG23	2.20	0.42
1:B:1092:VAL:HG12	1:B:1351:PHE:CG	2.54	0.42
1:B:1160:ALA:O	1:B:1167:LEU:HA	2.20	0.42
1:A:218:GLU:HB2	4:A:2172:HOH:O	2.19	0.42
1:B:1244:THR:HG23	1:B:1268:LEU:HB2	2.02	0.42
1:B:1203:TYR:HB2	1:B:1207:VAL:CG1	2.50	0.42
1:A:150:ILE:HG12	1:A:152:ARG:HG3	2.02	0.42
1:B:1155:LYS:HB3	4:B:2048:HOH:O	2.19	0.42
1:B:1161:VAL:HA	1:B:1166:GLU:O	2.19	0.42
1:B:1193:ALA:HB2	1:B:1210:TRP:CB	2.50	0.42
1:A:300:VAL:HG22	1:A:304:GLU:CB	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1328:ASP:CG	1:B:1329:SER:N	2.73	0.41
1:B:1141:LEU:HA	1:B:1141:LEU:HD23	1.74	0.41
1:A:328:ASP:CG	1:A:329:SER:N	2.73	0.41
1:A:325:LYS:HB3	1:A:325:LYS:HE2	1.80	0.41
1:B:1218:GLU:HB2	4:B:2116:HOH:O	2.21	0.41
1:B:1152:ARG:CZ	1:B:1176:ARG:HD2	2.51	0.41
1:A:16:VAL:CG2	1:A:21:TRP:CD1	3.03	0.41
1:A:92:VAL:HG12	1:A:351:PHE:CG	2.54	0.41
1:A:242:THR:C	1:A:270:LYS:HG2	2.40	0.41
1:B:1293:VAL:HG12	1:B:1295:ASP:H	1.86	0.41
1:A:253:LEU:HA	1:A:253:LEU:HD23	1.85	0.41
1:A:308:HIS:HA	1:A:309:PRO:HD3	1.97	0.41
1:B:1012:TYR:HD1	1:B:1025:ALA:HA	1.85	0.41
1:B:1325:LYS:HB3	1:B:1325:LYS:HE2	1.79	0.41
1:A:193:ALA:HB2	1:A:210:TRP:CB	2.50	0.40
1:A:204:THR:HG22	4:A:2089:HOH:O	2.21	0.40
1:B:1242:THR:C	1:B:1270:LYS:HG2	2.41	0.40
1:B:1230:ASP:H	1:B:1233:ASP:HB3	1.86	0.40
1:A:58:LEU:O	1:A:61:PRO:HD3	2.21	0.40
1:A:279:ASN:N	1:A:279:ASN:ND2	2.68	0.40
1:A:203:TYR:HB2	1:A:207:VAL:CG1	2.51	0.40
1:A:259:LYS:HE2	1:A:259:LYS:HB3	1.89	0.40
1:B:1279:ASN:HD22	1:B:1279:ASN:N	1.99	0.40
1:B:1328:ASP:HB3	4:B:2105:HOH:O	2.21	0.40
1:B:1058:LEU:O	1:B:1061:PRO:HD3	2.21	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	319/367 (87%)	295 (92%)	21 (7%)	3 (1%)	21	30

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	319/367 (87%)	295 (92%)	22 (7%)	2 (1%)	30	43
All	All	638/734 (87%)	590 (92%)	43 (7%)	5 (1%)	24	35

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	187	VAL
1	B	1187	VAL
1	A	171	ASP
1	A	16	VAL
1	B	1016	VAL

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	265/316 (84%)	251 (95%)	14 (5%)	28	44
1	B	265/316 (84%)	251 (95%)	14 (5%)	28	44
All	All	530/632 (84%)	502 (95%)	28 (5%)	28	44

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	THR
1	A	19	THR
1	A	26	VAL
1	A	53	VAL
1	A	89	LEU
1	A	109	MET
1	A	129	ARG
1	A	141	LEU
1	A	250	VAL
1	A	254	GLN
1	A	270	LYS

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Mol	Chain	Res	Type
1	A	279	ASN
1	A	289	GLU
1	A	300	VAL
1	B	1017	THR
1	B	1019	THR
1	B	1026	VAL
1	B	1033	VAL
1	B	1053	VAL
1	B	1089	LEU
1	B	1109	MET
1	B	1141	LEU
1	B	1250	VAL
1	B	1254	GLN
1	B	1270	LYS
1	B	1279	ASN
1	B	1289	GLU
1	B	1300	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	GLN
1	A	260	ASN
1	A	279	ASN
1	A	286	ASN
1	B	1131	GLN
1	B	1260	ASN
1	B	1279	ASN
1	B	1286	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	183	1	8,10,11	1.29	0	7,14,16	0.98	0
1	PTR	A	185	1	14,16,17	1.55	3 (21%)	18,22,24	1.20	2 (11%)
1	TPO	B	1183	1	8,10,11	1.53	3 (37%)	7,14,16	1.07	1 (14%)
1	PTR	B	1185	1	14,16,17	1.40	2 (14%)	18,22,24	1.12	1 (5%)

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	183	1	1/1/3/4	0/8/11/13	0/0/0/0
1	PTR	A	185	1	-	0/9/11/13	0/1/1/1
1	TPO	B	1183	1	1/1/3/4	0/8/11/13	0/0/0/0
1	PTR	B	1185	1	-	0/9/11/13	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1185	PTR	P-O3P	-2.82	1.44	1.54
1	A	185	PTR	P-O3P	-2.69	1.45	1.54
1	B	1183	TPO	P-OG1	-2.34	1.52	1.60
1	B	1185	PTR	P-O2P	-2.31	1.46	1.54
1	B	1183	TPO	P-O3P	-2.17	1.46	1.54
1	A	185	PTR	CD1-CG	2.01	1.43	1.38
1	A	185	PTR	CE2-CZ	2.13	1.42	1.38
1	B	1183	TPO	CB-CA	2.15	1.57	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1185	PTR	CE2-CD2-CG	-2.23	117.98	121.04
1	A	185	PTR	CE2-CD2-CG	-2.11	118.15	121.04
1	A	185	PTR	CG-CB-CA	-2.00	109.69	114.21
1	B	1183	TPO	O3P-P-O2P	2.13	115.51	107.38

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	183	TPO	CB
1	B	1183	TPO	CB

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	183	TPO	1	0
1	A	185	PTR	2	0
1	B	1183	TPO	1	0
1	B	1185	PTR	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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
3	ANP	A	400	2	27,33,33	2.88	5 (18%)	30,52,52	2.45	12 (40%)
3	ANP	B	1400	2	27,33,33	2.70	8 (29%)	30,52,52	2.56	10 (33%)

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	ANP	A	400	2	-	2/12/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ANP	B	1400	2	-	2/12/38/38	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1400	ANP	PG-N3B	-3.93	1.52	1.63
3	A	400	ANP	PG-N3B	-3.72	1.53	1.63
3	A	400	ANP	PB-N3B	-3.52	1.54	1.63
3	B	1400	ANP	PB-N3B	-2.56	1.56	1.63
3	B	1400	ANP	PB-O3A	2.31	1.62	1.59
3	B	1400	ANP	C6-N6	2.48	1.42	1.34
3	B	1400	ANP	O4'-C1'	2.70	1.44	1.41
3	B	1400	ANP	C5-C4	2.89	1.47	1.40
3	A	400	ANP	C5-C4	3.59	1.48	1.40
3	B	1400	ANP	PB-O1B	6.98	1.54	1.46
3	A	400	ANP	PG-O1G	8.90	1.56	1.46
3	B	1400	ANP	PG-O1G	9.16	1.56	1.46
3	A	400	ANP	PB-O1B	9.33	1.56	1.46

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1400	ANP	N3-C2-N1	-7.72	122.98	128.89
3	A	400	ANP	N3-C2-N1	-6.80	123.68	128.89
3	B	1400	ANP	O1G-PG-N3B	-4.20	105.46	111.90
3	A	400	ANP	O1G-PG-N3B	-3.30	106.83	111.90
3	B	1400	ANP	O3G-PG-O1G	-2.93	105.71	113.49
3	A	400	ANP	O3G-PG-O1G	-2.87	105.87	113.49
3	A	400	ANP	O1B-PB-N3B	-2.41	108.20	111.90
3	A	400	ANP	PA-O3A-PB	-2.03	125.85	132.67
3	A	400	ANP	C2'-C3'-C4'	2.07	106.86	102.61
3	B	1400	ANP	O3A-PB-N3B	2.08	112.17	106.44
3	B	1400	ANP	C2'-C3'-C4'	2.09	106.90	102.61
3	A	400	ANP	N6-C6-N1	2.27	124.08	119.20
3	A	400	ANP	O3A-PA-O5'	2.42	109.36	102.94
3	A	400	ANP	O3A-PB-N3B	2.51	113.34	106.44
3	B	1400	ANP	O3A-PA-O5'	2.53	109.64	102.94
3	B	1400	ANP	N6-C6-N1	2.56	124.69	119.20
3	A	400	ANP	O2B-PB-O1B	3.69	117.70	110.00
3	B	1400	ANP	C2-N1-C6	3.89	125.72	118.77
3	A	400	ANP	C2-N1-C6	3.98	125.89	118.77
3	B	1400	ANP	O2B-PB-O1B	4.18	118.72	110.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1400	ANP	O4'-C1'-N9	5.26	119.12	108.10
3	A	400	ANP	O4'-C1'-N9	5.54	119.69	108.10

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1400	ANP	O1B-PB-N3B-PG
3	A	400	ANP	O1B-PB-N3B-PG
3	B	1400	ANP	O1G-PG-N3B-PB
3	A	400	ANP	O1G-PG-N3B-PB

There are no ring outliers.

No monomer is involved in short contacts.

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.