



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 03:41 PM GMT

PDB ID : 3RDH
Title : X-ray induced covalent inhibition of 14-3-3
Authors : Horton, J.R.; Upadhyay, A.K.; Fu, H.; Cheng, X.
Deposited on : 2011-04-01
Resolution : 2.39 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

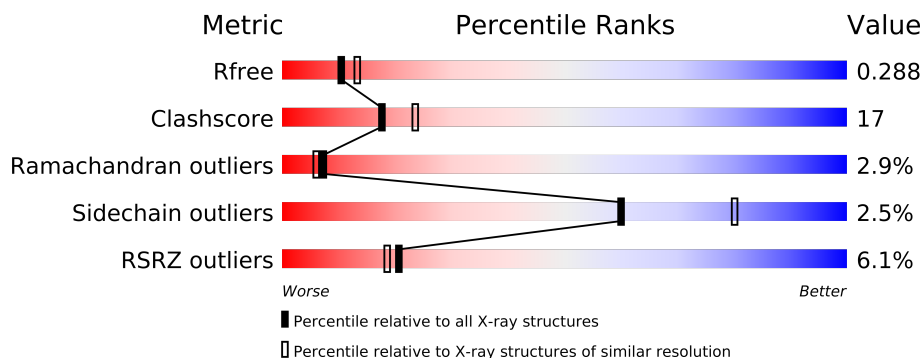
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.39 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	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (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. 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	248	
1	B	248	
1	C	248	
1	D	248	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	NI	A	247	-	X
3	NI	B	247	-	X
3	NI	C	247	-	X
3	NI	D	247	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7471 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.

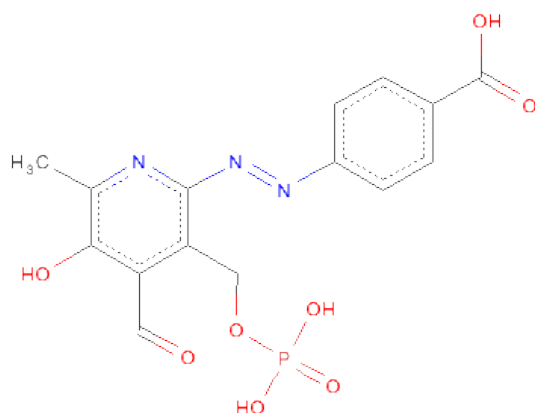
- Molecule 1 is a protein called 14-3-3 protein zeta/delta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	3	0
			1862	1163	316	373	10			
1	B	233	Total	C	N	O	S	0	1	0
			1815	1134	307	364	10			
1	C	233	Total	C	N	O	S	0	0	0
			1820	1136	309	365	10			
1	D	233	Total	C	N	O	S	0	0	0
			1816	1138	304	364	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P63104
A	-1	SER	-	EXPRESSION TAG	UNP P63104
A	0	HIS	-	EXPRESSION TAG	UNP P63104
B	-2	GLY	-	EXPRESSION TAG	UNP P63104
B	-1	SER	-	EXPRESSION TAG	UNP P63104
B	0	HIS	-	EXPRESSION TAG	UNP P63104
C	-2	GLY	-	EXPRESSION TAG	UNP P63104
C	-1	SER	-	EXPRESSION TAG	UNP P63104
C	0	HIS	-	EXPRESSION TAG	UNP P63104
D	-2	GLY	-	EXPRESSION TAG	UNP P63104
D	-1	SER	-	EXPRESSION TAG	UNP P63104
D	0	HIS	-	EXPRESSION TAG	UNP P63104

- Molecule 2 is 4-[(E)-{4-FORMYL-5-HYDROXY-6-METHYL-3-[(PHOSPHONOOXY)METHYL]PYRIDIN-2-YL}DIAZENYL]BENZOICACID (three-letter code: 3RD) (formula: C₁₅H₁₄N₃O₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			17	8	2	6	1		
2	B	1	Total	C	N	O	P	0	0
			17	8	2	6	1		
2	C	1	Total	C	N	O	P	0	0
			17	8	2	6	1		
2	D	1	Total	C	N	O	P	0	0
			17	8	2	6	1		

- Molecule 3 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ni	0	0
			1	1		
3	A	1	Total	Ni	0	0
			1	1		
3	D	1	Total	Ni	0	0
			1	1		
3	C	1	Total	Ni	0	0
			1	1		

- Molecule 4 is water.

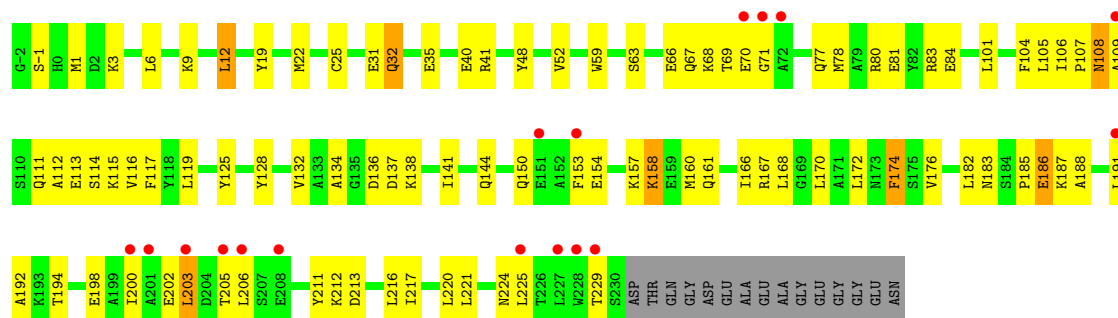
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	21	Total	O	0	0
			21	21		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	24	Total 24	O 24	0	0
4	C	21	Total 21	O 21	0	0
4	D	20	Total 20	O 20	0	0

- Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	94.75Å 94.75Å 237.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.77 – 2.39 33.77 – 2.38	Depositor EDS
% Data completeness (in resolution range)	93.8 (33.77-2.39) 93.6 (33.77-2.38)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.39Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.235 , 0.286 0.243 , 0.288	Depositor DCC
R_{free} test set	2230 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.451	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 33.5	EDS
Estimated twinning fraction	0.217 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	1 of 47686 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7471	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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: NI, 3RD

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	0/1888	0.56	0/2543
1	B	0.38	0/1841	0.55	0/2483
1	C	0.36	0/1845	0.54	0/2486
1	D	0.37	0/1842	0.55	0/2485
All	All	0.37	0/7416	0.55	0/9997

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1862	0	1811	75	0
1	B	1815	0	1738	59	0
1	C	1820	0	1763	57	0
1	D	1816	0	1752	75	0
2	A	17	0	9	0	0
2	B	17	0	9	0	0
2	C	17	0	9	0	0
2	D	17	0	9	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	21	0	0	1	0
4	B	24	0	0	0	0
4	C	21	0	0	1	0
4	D	20	0	0	0	0
All	All	7471	0	7100	251	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 17.

All (251) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:113:GLU:HA	1:D:160:MET:HE3	1.55	0.88
1:C:3:LYS:HA	1:C:6:LEU:HD12	1.63	0.80
1:A:200:ILE:HD13	1:A:203:LEU:HD21	1.62	0.80
1:D:108:ASN:HD22	1:D:108:ASN:N	1.81	0.78
1:B:5:GLU:HG3	1:C:78:MET:HE3	1.66	0.77
1:B:69:THR:O	1:B:76:GLN:HB2	1.87	0.75
1:C:196:PHE:O	1:C:200:ILE:HG13	1.87	0.74
1:D:168:LEU:HB3	1:D:217:ILE:HG21	1.69	0.74
1:C:196:PHE:CE1	1:C:200:ILE:HD11	2.24	0.72
1:A:49:LYS:HD3	1:A:49:LYS:C	2.11	0.71
1:A:226:THR:HA	1:A:229:THR:HG22	1.72	0.70
1:D:202:GLU:O	1:D:205:THR:HG22	1.92	0.70
1:A:127:ARG:HD2	4:A:262:HOH:O	1.91	0.70
1:B:51:VAL:HG12	1:B:93:ILE:HD13	1.73	0.69
1:A:83:ARG:O	1:A:87:GLU:HG3	1.93	0.69
1:A:146:GLN:OE1	1:A:182:LEU:HD11	1.94	0.68
1:C:58:SER:O	1:C:62:VAL:HG23	1.95	0.66
1:D:69:THR:O	1:D:71:GLY:N	2.29	0.65
1:C:138:LYS:HE3	1:C:142:VAL:HG21	1.77	0.65
1:A:212:LYS:O	1:A:216:LEU:HB2	1.96	0.65
1:D:67:GLN:C	1:D:69:THR:H	1.99	0.65
1:D:200:ILE:O	1:D:203:LEU:HB2	1.96	0.65
1:D:192:ALA:HB3	1:D:225:LEU:HD13	1.80	0.64
1:D:80:ARG:O	1:D:84:GLU:HG3	1.98	0.63
1:B:146:GLN:OE1	1:B:182:LEU:HD11	2.00	0.61
1:C:210:SER:HA	1:C:213:ASP:HB2	1.83	0.61
1:A:62:VAL:HG22	1:D:12:LEU:HD21	1.82	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:202:GLU:HB3	1:A:205:THR:CG2	2.32	0.60
1:A:172:LEU:O	1:A:176:VAL:HG23	2.03	0.59
1:C:197:ASP:HA	1:C:200:ILE:HD12	1.84	0.58
1:A:55:ARG:HB3	1:A:90:LEU:HG	1.86	0.58
1:C:37:SER:OG	1:C:40:GLU:HG3	2.04	0.58
1:D:67:GLN:O	1:D:69:THR:N	2.37	0.58
1:C:208:GLU:C	1:C:210:SER:H	2.06	0.57
1:D:216:LEU:O	1:D:220:LEU:HG	2.05	0.57
1:C:49:LYS:HE3	1:C:49:LYS:HA	1.87	0.57
1:A:106:ILE:HB	1:A:107:PRO:HD3	1.85	0.57
1:A:51:VAL:HG12	1:A:93:ILE:HD13	1.86	0.57
1:D:172:LEU:O	1:D:176:VAL:HG23	2.05	0.57
1:A:178:TYR:CD1	1:A:182:LEU:HD12	2.40	0.57
1:B:78:MET:HE1	1:C:5:GLU:HB3	1.87	0.57
1:A:150:GLN:O	1:A:154:GLU:HG2	2.05	0.57
1:D:66:GLU:CD	1:D:83:ARG:HE	2.08	0.57
1:B:221:LEU:O	1:B:221:LEU:HD23	2.05	0.57
1:C:101:LEU:HD23	1:C:105:LEU:HD12	1.87	0.57
1:B:220[A]:LEU:HA	1:B:223:ASP:OD2	2.06	0.56
1:D:112:ALA:O	1:D:116:VAL:HG23	2.04	0.56
1:D:108:ASN:ND2	1:D:108:ASN:N	2.54	0.55
1:D:153:PHE:HA	1:D:170:LEU:HD21	1.89	0.55
1:D:48:TYR:O	1:D:52:VAL:HG12	2.06	0.55
1:A:202:GLU:HB3	1:A:205:THR:HB	1.88	0.55
1:A:157:LYS:HD2	1:A:167:ARG:NH2	2.21	0.55
1:D:77:GLN:O	1:D:81:GLU:HG3	2.05	0.55
1:B:64:SER:O	1:B:65:ILE:C	2.45	0.55
1:C:174:PHE:O	1:C:177:PHE:HB3	2.08	0.54
1:B:38:ASN:OD1	1:B:39:GLU:N	2.34	0.54
1:C:1:MET:HE2	4:C:259:HOH:O	2.07	0.54
1:C:65:ILE:HG22	1:C:65:ILE:O	2.08	0.54
1:B:73:GLU:O	1:B:77:GLN:HB2	2.07	0.54
1:B:77:GLN:O	1:B:81:GLU:HG3	2.07	0.54
1:A:207:SER:O	1:A:211:TYR:N	2.41	0.53
1:B:106:ILE:HB	1:B:107:PRO:HD3	1.91	0.53
1:D:211:TYR:CE2	1:D:212:LYS:HG3	2.43	0.53
1:A:3:LYS:HA	1:A:6:LEU:HD12	1.90	0.53
1:B:162:PRO:HG3	1:B:167:ARG:NH2	2.22	0.53
1:D:41:ARG:NH1	1:D:114:SER:HB3	2.24	0.53
1:D:221:LEU:O	1:D:225:LEU:HB2	2.07	0.53
1:D:203:LEU:O	1:D:206:LEU:HG	2.09	0.53
1:C:196:PHE:CD1	1:C:200:ILE:HD11	2.44	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:150:GLN:O	1:D:154:GLU:HG2	2.09	0.53
1:A:2:ASP:OD1	1:A:5:GLU:HB2	2.09	0.53
1:D:67:GLN:C	1:D:69:THR:N	2.63	0.53
1:A:67:GLN:C	1:A:69:THR:H	2.13	0.52
1:D:35:GLU:OE2	1:D:109:ALA:HA	2.08	0.52
1:A:215:THR:HA	1:A:218:MET:HB2	1.90	0.52
1:D:9:LYS:HG3	1:D:25:CYS:SG	2.49	0.52
1:C:163:THR:CG2	1:C:202:GLU:HB2	2.40	0.52
1:A:202:GLU:O	1:A:205:THR:HG22	2.09	0.52
1:B:224:ASN:HB3	1:B:228:TRP:CZ2	2.45	0.52
1:A:208:GLU:HA	1:A:211:TYR:HD2	1.75	0.52
1:A:200:ILE:HA	1:A:203:LEU:HG	1.91	0.52
1:B:220[B]:LEU:HA	1:B:223:ASP:OD2	2.10	0.52
1:C:74:LYS:O	1:C:77:GLN:HB3	2.10	0.52
1:C:104:PHE:O	1:C:108:ASN:HB2	2.10	0.52
1:D:106:ILE:HB	1:D:107:PRO:HD3	1.92	0.52
1:B:36:LEU:HB3	1:B:40:GLU:HB2	1.92	0.51
1:D:185:PRO:HA	1:D:188:ALA:HB3	1.92	0.51
1:C:212:LYS:O	1:C:216:LEU:HD13	2.10	0.51
1:D:211:TYR:CZ	1:D:212:LYS:HG3	2.45	0.51
1:B:4:ASN:O	1:B:8:GLN:HG2	2.11	0.51
1:A:208:GLU:OE1	1:A:208:GLU:N	2.41	0.51
1:A:4:ASN:O	1:A:8[A]:GLN:HG3	2.10	0.51
1:D:59:TRP:CE2	1:D:132:VAL:HG12	2.46	0.51
1:B:78:MET:HE3	1:C:5:GLU:HG2	1.92	0.51
1:D:107:PRO:C	1:D:108:ASN:HD22	2.15	0.51
1:D:185:PRO:HG2	1:D:186:GLU:OE2	2.11	0.50
1:A:210:SER:HA	1:A:213:ASP:OD2	2.11	0.50
1:B:185:PRO:HA	1:B:188:ALA:HB3	1.92	0.50
1:B:41:ARG:HH11	1:B:114:SER:HB3	1.77	0.50
1:B:63:SER:O	1:B:67:GLN:HG2	2.12	0.50
1:D:157:LYS:O	1:D:157:LYS:HG2	2.12	0.50
1:D:158:LYS:HG2	1:D:158:LYS:O	2.11	0.50
1:A:80:ARG:HG3	1:A:83:ARG:HH21	1.76	0.50
1:D:160:MET:O	1:D:167:ARG:NH1	2.44	0.49
1:D:168:LEU:HB3	1:D:217:ILE:CG2	2.39	0.49
1:B:64:SER:O	1:B:67:GLN:N	2.45	0.49
1:A:113:GLU:OE1	1:A:164:HIS:HE1	1.94	0.49
1:A:200:ILE:O	1:A:203:LEU:HG	2.13	0.49
1:A:80:ARG:O	1:A:84:GLU:HG3	2.11	0.49
1:B:77:GLN:HE21	1:B:80:ARG:NH2	2.11	0.49
1:C:119:LEU:HB3	1:C:152:ALA:HB2	1.93	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:0:HIS:CD2	1:C:1:MET:HG3	2.48	0.49
1:D:3:LYS:HB2	1:D:32:GLN:HG2	1.95	0.49
1:A:14:GLU:HB2	1:A:22:MET:SD	2.53	0.49
1:A:151:GLU:O	1:A:155:ILE:HG13	2.13	0.49
1:B:5:GLU:HG3	1:C:78:MET:CE	2.41	0.48
1:B:57:SER:O	1:B:61:VAL:HG23	2.13	0.48
1:C:157:LYS:HG3	1:C:157:LYS:O	2.12	0.48
1:C:51:VAL:HG13	1:C:55:ARG:NH1	2.28	0.48
1:D:161:GLN:O	1:D:167:ARG:HD3	2.13	0.48
1:B:78:MET:CE	1:C:5:GLU:HB3	2.43	0.48
1:B:73:GLU:OE1	1:B:73:GLU:HA	2.13	0.48
1:B:202:GLU:C	1:B:204:ASP:H	2.16	0.48
1:B:202:GLU:C	1:B:204:ASP:N	2.67	0.48
1:A:164:HIS:CD2	1:A:165:PRO:HD2	2.48	0.48
1:A:215:THR:O	1:A:219:GLN:HG3	2.14	0.48
1:B:153:PHE:HA	1:B:170:LEU:HD21	1.96	0.48
1:D:3:LYS:NZ	1:D:40:GLU:OE2	2.39	0.47
1:A:42:ASN:O	1:A:46:VAL:HG23	2.14	0.47
1:D:80:ARG:HG3	1:D:83:ARG:NH2	2.28	0.47
1:A:181:ILE:C	1:A:183[A]:ASN:H	2.17	0.47
1:D:172:LEU:HD11	1:D:224:ASN:ND2	2.29	0.47
1:C:210:SER:O	1:C:211:TYR:C	2.53	0.47
1:B:200:ILE:C	1:B:202:GLU:H	2.18	0.47
1:B:3:LYS:HA	1:B:6:LEU:HD12	1.96	0.47
1:B:14:GLU:HB2	1:B:22:MET:SD	2.54	0.47
1:B:101:LEU:HD23	1:B:105:LEU:HD12	1.96	0.47
1:A:49:LYS:O	1:A:49:LYS:HD3	2.13	0.47
1:C:55:ARG:N	1:C:55:ARG:HD2	2.30	0.47
1:A:163:THR:HB	1:A:206:LEU:HD23	1.97	0.47
1:A:104:PHE:CZ	1:C:103:LYS:HG3	2.50	0.47
1:A:221:LEU:O	1:A:225:LEU:HB2	2.15	0.47
1:B:164:HIS:HD2	1:B:166:ILE:H	1.63	0.47
1:C:98:LEU:HA	1:C:101:LEU:HD12	1.95	0.46
1:B:80:ARG:HG2	1:B:84:GLU:OE2	2.15	0.46
1:B:83:ARG:NH1	1:B:87:GLU:OE2	2.47	0.46
1:B:36:LEU:HD22	1:B:40:GLU:HB3	1.96	0.46
1:A:178:TYR:HD1	1:A:182:LEU:HD12	1.80	0.46
1:C:210:SER:HB3	1:C:214:SER:OG	2.15	0.46
1:A:150:GLN:HE21	1:A:154:GLU:CD	2.19	0.46
1:C:227:LEU:C	1:C:229:THR:H	2.18	0.46
1:D:200:ILE:HG23	1:D:203:LEU:HD12	1.97	0.46
1:A:93:ILE:O	1:A:97:VAL:HG23	2.15	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:203:LEU:HD21	1:B:215:THR:CG2	2.46	0.46
1:D:125:TYR:HA	1:D:128:TYR:CD2	2.51	0.46
1:C:160:MET:HE3	1:C:164:HIS:CE1	2.50	0.46
1:C:48:TYR:O	1:C:52:VAL:HG12	2.15	0.46
1:A:182:LEU:O	1:A:183[B]:ASN:C	2.54	0.46
1:A:67:GLN:O	1:A:69:THR:N	2.49	0.46
1:B:66:GLU:OE2	1:B:80:ARG:HG3	2.16	0.45
1:A:3:LYS:NZ	1:A:40:GLU:OE2	2.47	0.45
1:D:170:LEU:O	1:D:174:PHE:HB2	2.16	0.45
1:A:227:LEU:O	1:A:227:LEU:HD12	2.16	0.45
1:C:208:GLU:C	1:C:210:SER:N	2.69	0.45
1:D:115:LYS:HB2	1:D:115:LYS:HE3	1.77	0.45
1:C:3:LYS:O	1:C:6:LEU:HB2	2.17	0.45
1:A:62:VAL:HG22	1:D:12:LEU:CD2	2.46	0.45
1:B:174:PHE:CD1	1:B:174:PHE:O	2.70	0.45
1:C:3:LYS:NZ	1:C:40:GLU:OE2	2.49	0.45
1:D:41:ARG:HH11	1:D:114:SER:HB3	1.82	0.45
1:C:67:GLN:C	1:C:69:THR:H	2.20	0.45
1:A:78:MET:SD	1:D:12:LEU:HD12	2.57	0.45
1:A:202:GLU:HB3	1:A:205:THR:CB	2.47	0.45
1:D:1:MET:HE3	1:D:6:LEU:HD11	1.99	0.45
1:A:102:GLU:OE1	1:A:122:LYS:HE3	2.17	0.45
1:C:71:GLY:O	1:C:73:GLU:N	2.50	0.44
1:B:103:LYS:HE3	1:D:31:GLU:O	2.17	0.44
1:A:177:PHE:O	1:A:181:ILE:HB	2.17	0.44
1:B:153:PHE:CA	1:B:170:LEU:HD21	2.47	0.44
1:A:119:LEU:CB	1:A:152:ALA:HB2	2.47	0.44
1:A:119:LEU:HB3	1:A:152:ALA:HB2	1.98	0.44
1:B:27:LYS:HG3	1:B:100:LEU:HD11	1.98	0.44
1:A:87:GLU:O	1:A:91:ARG:HG3	2.17	0.44
1:A:157:LYS:HG3	1:A:157:LYS:O	2.18	0.44
1:A:8[A]:GLN:HB2	1:D:78:MET:HE2	2.00	0.44
1:D:101:LEU:HD23	1:D:105:LEU:HD12	2.00	0.44
1:A:181:ILE:C	1:A:183[B]:ASN:H	2.19	0.44
1:B:164:HIS:CD2	1:B:165:PRO:HD2	2.53	0.44
1:C:106:ILE:HB	1:C:107:PRO:HD3	2.00	0.43
1:D:66:GLU:OE2	1:D:80:ARG:HA	2.18	0.43
1:B:66:GLU:CD	1:B:83:ARG:HE	2.20	0.43
1:A:196:PHE:O	1:A:200:ILE:HG12	2.18	0.43
1:A:182:LEU:O	1:A:183[A]:ASN:C	2.55	0.43
1:D:117:PHE:HB2	1:D:166:ILE:HD13	1.99	0.43
1:C:208:GLU:HA	1:C:211:TYR:CB	2.48	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:194:THR:HG22	1:D:198:GLU:OE2	2.18	0.43
1:D:113:GLU:HG3	1:D:160:MET:CE	2.48	0.43
1:A:202:GLU:O	1:A:205:THR:N	2.43	0.43
1:D:182:LEU:HA	1:D:182:LEU:HD23	1.87	0.43
1:B:119:LEU:HD12	1:B:155:ILE:HD12	2.01	0.43
1:B:7:VAL:HG12	1:B:11:LYS:HE3	2.01	0.43
1:C:150:GLN:O	1:C:153:PHE:HB3	2.19	0.43
1:A:163:THR:HG21	1:A:205:THR:HG22	2.00	0.43
1:B:150:GLN:O	1:B:154:GLU:HG2	2.18	0.43
1:C:178:TYR:CD1	1:C:182:LEU:HD12	2.54	0.42
1:C:125:TYR:HA	1:C:128:TYR:CD2	2.54	0.42
1:A:200:ILE:HA	1:A:203:LEU:CD2	2.49	0.42
1:C:119:LEU:CB	1:C:152:ALA:HB2	2.49	0.42
1:B:211:TYR:CG	1:B:212:LYS:N	2.86	0.42
1:C:91:ARG:HA	1:C:129:LEU:HD21	2.01	0.42
1:D:101:LEU:HA	1:D:105:LEU:HB2	2.01	0.42
1:B:226:THR:HA	1:B:229:THR:OG1	2.20	0.42
1:D:229:THR:O	1:D:229:THR:HG22	2.20	0.42
1:A:207:SER:C	1:A:209:GLU:N	2.73	0.42
1:D:134:ALA:O	1:D:138:LYS:HB2	2.20	0.42
1:D:141:ILE:O	1:D:144:GLN:HB2	2.19	0.42
1:A:0:HIS:HD2	1:C:95:ASN:OD1	2.03	0.42
1:C:45:SER:O	1:C:49:LYS:HB2	2.20	0.42
1:C:52:VAL:HG13	1:C:53:GLY:N	2.35	0.42
1:C:27:LYS:O	1:C:31:GLU:HG3	2.19	0.42
1:D:203:LEU:O	1:D:203:LEU:HD23	2.19	0.42
1:D:41:ARG:HH11	1:D:114:SER:CB	2.33	0.42
1:A:67:GLN:C	1:A:69:THR:N	2.73	0.41
1:B:8:GLN:HB2	1:C:78:MET:HE1	2.02	0.41
1:D:106:ILE:HD13	1:D:119:LEU:HD23	2.02	0.41
1:B:8:GLN:HB2	1:C:78:MET:CE	2.51	0.41
1:D:168:LEU:CB	1:D:217:ILE:HG21	2.44	0.41
1:D:153:PHE:CA	1:D:170:LEU:HD21	2.51	0.41
1:B:151:GLU:O	1:B:155:ILE:HG13	2.19	0.41
1:D:19:TYR:HA	1:D:22:MET:HB3	2.02	0.41
1:A:164:HIS:HD2	1:A:166:ILE:H	1.69	0.41
1:C:124:ASP:O	1:C:127:ARG:HB3	2.20	0.41
1:D:1:MET:HG3	1:D:32:GLN:HE22	1.86	0.41
1:D:32:GLN:OE1	1:D:32:GLN:HA	2.20	0.41
1:A:22:MET:HG2	1:A:47:ALA:HB2	2.02	0.41
1:B:124:ASP:O	1:B:127:ARG:HB3	2.21	0.41
1:A:163:THR:HB	1:A:206:LEU:CD2	2.51	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:224:ASN:HB3	1:B:228:TRP:CE2	2.56	0.41
1:C:83:ARG:HG2	1:C:83:ARG:HH11	1.87	0.40
1:B:161:GLN:NE2	1:B:163:THR:OG1	2.54	0.40
1:A:75:LYS:O	1:A:78:MET:HB3	2.21	0.40
1:D:211:TYR:OH	1:D:212:LYS:HE2	2.21	0.40
1:D:31:GLU:HA	1:D:104:PHE:CE2	2.56	0.40
1:B:58:SER:HB3	1:B:86:ILE:HD13	2.03	0.40
1:C:157:LYS:CG	1:C:157:LYS:O	2.69	0.40
1:A:70:GLU:O	1:A:71:GLY:C	2.60	0.40
1:A:226:THR:CA	1:A:229:THR:HG22	2.45	0.40
1:A:12:LEU:HD12	1:D:78:MET:SD	2.62	0.40
1:B:203:LEU:HD12	1:B:203:LEU:HA	1.93	0.40
1:D:63:SER:O	1:D:66:GLU:HB3	2.22	0.40
1:D:186:GLU:HB2	1:D:187:LYS:H	1.79	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	234/248 (94%)	203 (87%)	24 (10%)	7 (3%)	7	5
1	B	232/248 (94%)	207 (89%)	18 (8%)	7 (3%)	7	5
1	C	231/248 (93%)	209 (90%)	17 (7%)	5 (2%)	10	11
1	D	231/248 (93%)	208 (90%)	15 (6%)	8 (4%)	6	4
All	All	928/992 (94%)	827 (89%)	74 (8%)	27 (3%)	7	6

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	-1	SER
1	C	69	THR
1	C	72	ALA
1	A	-1	SER

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Mol	Chain	Res	Type
1	A	68	LYS
1	A	135	GLY
1	A	188	ALA
1	B	-1	SER
1	B	183	ASN
1	B	201	ALA
1	D	-1	SER
1	D	68	LYS
1	D	70	GLU
1	D	158	LYS
1	D	186	GLU
1	A	186	GLU
1	A	210	SER
1	B	208	GLU
1	C	73	GLU
1	C	183	ASN
1	B	65	ILE
1	D	136	ASP
1	A	182	LEU
1	B	192	ALA
1	D	203	LEU
1	D	183	ASN
1	B	200	ILE

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	198/211 (94%)	195 (98%)	3 (2%)	76	90
1	B	188/211 (89%)	187 (100%)	1 (0%)	94	98
1	C	191/211 (90%)	184 (96%)	7 (4%)	45	66
1	D	190/211 (90%)	182 (96%)	8 (4%)	40	60
All	All	767/844 (91%)	748 (98%)	19 (2%)	60	80

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	174	PHE
1	A	205	THR
1	B	12	LEU
1	C	12	LEU
1	C	49	LYS
1	C	90	LEU
1	C	121	MET
1	C	174	PHE
1	C	183	ASN
1	C	222	ARG
1	D	12	LEU
1	D	32	GLN
1	D	108	ASN
1	D	111	GLN
1	D	137	ASP
1	D	174	PHE
1	D	191	LEU
1	D	213	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	108	ASN
1	A	150	GLN
1	A	164	HIS
1	B	32	GLN
1	B	77	GLN
1	B	108	ASN
1	B	161	GLN
1	B	164	HIS
1	B	224	ASN
1	C	32	GLN
1	C	108	ASN
1	C	111	GLN
1	C	146	GLN
1	C	150	GLN
1	C	164	HIS
1	D	77	GLN
1	D	108	ASN
1	D	224	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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$
2	3RD	A	246	1	16,17,28	2.28	6 (37%)	23,25,40	2.30	5 (21%)
2	3RD	B	246	1	16,17,28	2.22	6 (37%)	23,25,40	2.20	5 (21%)
2	3RD	C	246	1	16,17,28	2.12	6 (37%)	23,25,40	2.11	6 (26%)
2	3RD	D	246	1	16,17,28	2.21	6 (37%)	23,25,40	2.17	6 (26%)

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	3RD	A	246	1	-	0/8/8/17	0/1/1/2
2	3RD	B	246	1	-	0/8/8/17	0/1/1/2
2	3RD	C	246	1	-	0/8/8/17	0/1/1/2
2	3RD	D	246	1	-	0/8/8/17	0/1/1/2

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	246	3RD	C5-C6	4.41	1.57	1.46
2	C	246	3RD	C5-C6	4.36	1.56	1.46
2	B	246	3RD	C5-C6	4.20	1.56	1.46
2	D	246	3RD	C5-C6	4.17	1.56	1.46
2	B	246	3RD	C5-C1	3.70	1.45	1.40
2	B	246	3RD	C1-C2	3.69	1.43	1.40
2	A	246	3RD	C5-C1	3.60	1.45	1.40
2	D	246	3RD	C1-C2	3.57	1.43	1.40
2	D	246	3RD	C5-C1	3.30	1.44	1.40
2	A	246	3RD	C1-C2	3.27	1.43	1.40
2	C	246	3RD	C1-C2	3.24	1.43	1.40
2	A	246	3RD	C5-C4	3.22	1.46	1.41
2	D	246	3RD	C8-C4	3.12	1.56	1.51
2	D	246	3RD	C5-C4	3.05	1.46	1.41
2	C	246	3RD	C5-C1	2.98	1.44	1.40
2	C	246	3RD	C5-C4	2.92	1.45	1.41
2	A	246	3RD	C8-C4	2.84	1.56	1.51
2	B	246	3RD	C8-C4	2.76	1.55	1.51
2	B	246	3RD	C5-C4	2.70	1.45	1.41
2	C	246	3RD	C8-C4	2.67	1.55	1.51
2	A	246	3RD	P1-O5	2.07	1.62	1.54
2	B	246	3RD	P1-O5	2.03	1.62	1.54
2	C	246	3RD	P1-O5	2.03	1.62	1.54
2	D	246	3RD	C7-C2	2.02	1.54	1.50

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	246	3RD	O3-C8-C4	-7.01	95.76	109.22
2	B	246	3RD	O3-C8-C4	-6.37	96.99	109.22
2	D	246	3RD	O3-C8-C4	-5.97	97.75	109.22
2	C	246	3RD	O3-C8-C4	-5.79	98.10	109.22
2	D	246	3RD	C3-N2-C2	5.27	123.60	116.90
2	C	246	3RD	C3-N2-C2	5.21	123.52	116.90
2	A	246	3RD	C3-N2-C2	5.11	123.40	116.90
2	B	246	3RD	C3-N2-C2	5.05	123.32	116.90
2	A	246	3RD	P1-O3-C8	3.79	134.90	121.22
2	B	246	3RD	P1-O3-C8	3.33	133.22	121.22
2	D	246	3RD	P1-O3-C8	3.19	132.72	121.22
2	C	246	3RD	P1-O3-C8	2.96	131.90	121.22
2	B	246	3RD	O1-C6-C5	-2.94	118.96	125.32
2	D	246	3RD	O1-C6-C5	-2.89	119.09	125.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	246	3RD	C5-C1-C2	-2.88	118.24	120.18
2	A	246	3RD	C5-C1-C2	-2.85	118.27	120.18
2	B	246	3RD	C5-C1-C2	-2.77	118.32	120.18
2	D	246	3RD	C5-C1-C2	-2.68	118.38	120.18
2	C	246	3RD	O1-C6-C5	-2.64	119.63	125.32
2	A	246	3RD	O1-C6-C5	-2.50	119.91	125.32
2	C	246	3RD	C4-C3-N2	-2.24	120.83	123.13
2	D	246	3RD	C4-C3-N2	-2.02	121.05	123.13

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, 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	233/248 (93%)	0.60	8 (3%) 43 41	26, 50, 85, 93	0
1	B	233/248 (93%)	0.61	20 (8%) 11 9	27, 56, 100, 115	0
1	C	233/248 (93%)	0.60	12 (5%) 26 24	34, 54, 94, 113	0
1	D	233/248 (93%)	0.67	17 (7%) 15 13	32, 58, 103, 112	0
All	All	932/992 (93%)	0.62	57 (6%) 21 18	26, 55, 96, 115	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	206	LEU	5.8
1	A	71	GLY	5.1
1	A	209	GLU	4.9
1	B	210	SER	4.7
1	C	207	SER	4.5
1	D	205	THR	4.2
1	D	225	LEU	3.9
1	D	228	TRP	3.8
1	B	209	GLU	3.8
1	C	70	GLU	3.6
1	D	200	ILE	3.5
1	D	72	ALA	3.4
1	D	203	LEU	3.4
1	A	211	TYR	3.4
1	C	211	TYR	3.3
1	B	191	LEU	3.1
1	B	203	LEU	3.1
1	A	203	LEU	3.1
1	C	195	ALA	3.0
1	B	229	THR	2.9
1	A	196	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	73	GLU	2.9
1	D	153	PHE	2.9
1	D	71	GLY	2.9
1	C	199	ALA	2.9
1	B	225	LEU	2.8
1	B	74	LYS	2.7
1	C	227	LEU	2.7
1	B	134	ALA	2.7
1	C	34	ALA	2.7
1	B	220[A]	LEU	2.6
1	D	227	LEU	2.6
1	B	160	MET	2.6
1	D	229	THR	2.5
1	C	216	LEU	2.5
1	D	70	GLU	2.5
1	D	201	ALA	2.4
1	B	228	TRP	2.4
1	A	192	ALA	2.3
1	B	207	SER	2.3
1	B	69	THR	2.3
1	D	109	ALA	2.3
1	B	153	PHE	2.3
1	B	181	ILE	2.2
1	B	73	GLU	2.2
1	C	202	GLU	2.2
1	B	211	TYR	2.2
1	D	208	GLU	2.2
1	B	217	ILE	2.1
1	A	226	THR	2.1
1	D	191	LEU	2.1
1	C	191	LEU	2.1
1	B	178	TYR	2.1
1	C	161	GLN	2.1
1	D	151	GLU	2.1
1	A	200	ILE	2.0
1	B	196	PHE	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	NI	A	247	1/1	0.20	4.00	45,45,45,45	0
3	NI	B	247	1/1	0.20	2.99	50,50,50,50	0
3	NI	D	247	1/1	0.19	2.14	48,48,48,48	0
3	NI	C	247	1/1	0.18	2.13	42,42,42,42	0
2	3RD	C	246	17/27	0.15	-0.02	64,70,74,75	0
2	3RD	D	246	17/27	0.16	-0.19	65,71,78,78	0
2	3RD	B	246	17/27	0.14	-0.86	62,66,74,74	0
2	3RD	A	246	17/27	0.12	-1.27	58,62,69,70	0

6.5 Other polymers ⓘ

There are no such residues in this entry.