



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:20 AM GMT

PDB ID : 2WPV
Title : CRYSTAL STRUCTURE OF S. CEREVISIAE GET4-GET5 COMPLEX
Authors : Chang, Y.-W.; Chuang, Y.-C.; Ho, Y.-C.; Cheng, M.-Y.; Sun, Y.-J.; Hsiao, C.-D.; Wang, C.
Deposited on : 2009-08-11
Resolution : 1.99 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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

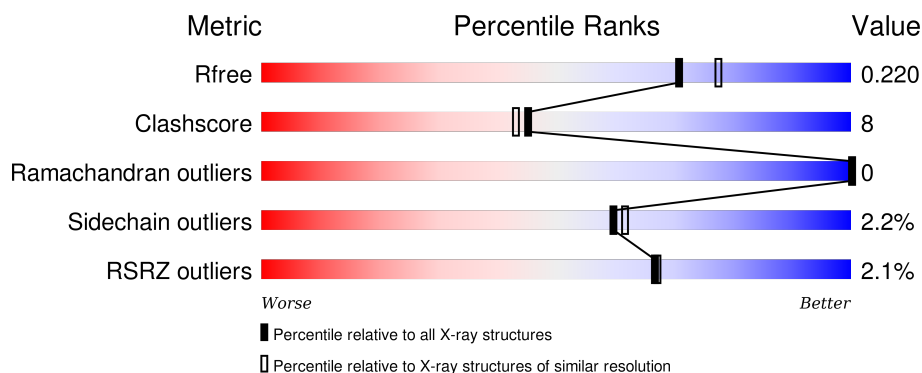
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	312	<div> <div>0%</div> <div> <div></div> <div>75%</div> <div>13%</div> <div>10%</div> </div> </div>
1	C	312	<div> <div>3%</div> <div> <div></div> <div>71%</div> <div>19%</div> <div>10%</div> </div> </div>
1	E	312	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>12%</div> <div>10%</div> </div> </div>
1	G	312	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>14%</div> <div>10%</div> </div> </div>
2	B	59	<div> <div>3%</div> <div> <div></div> <div>66%</div> <div>15%</div> <div>19%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	D	59	<div><div><div></div><div></div><div></div></div><div><div>3%</div><div>73%</div><div>7%</div><div>•</div><div>19%</div></div></div>
2	F	59	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>71%</div><div>8%</div><div></div><div>20%</div></div></div>
2	H	59	<div><div><div></div><div></div><div></div></div><div><div>3%</div><div>64%</div><div>17%</div><div></div><div>19%</div></div></div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 12001 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 UPF0363 PROTEIN YOR164C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2334	1521	369	439	5			
1	C	282	Total	C	N	O	S	0	0	0
			2318	1512	366	435	5			
1	E	280	Total	C	N	O	S	0	0	0
			2303	1502	360	436	5			
1	G	281	Total	C	N	O	S	0	0	0
			2329	1519	367	438	5			

- Molecule 2 is a protein called UBIQUITIN-LIKE PROTEIN MDY2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	48	Total	C	N	O	0	0	0
			374	249	58	67			
2	D	48	Total	C	N	O	0	0	0
			382	254	59	69			
2	F	47	Total	C	N	O	0	0	0
			373	249	57	67			
2	H	48	Total	C	N	O	0	0	0
			382	254	59	69			

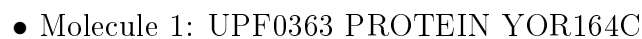
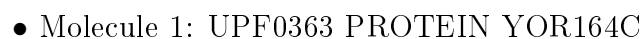
- Molecule 3 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

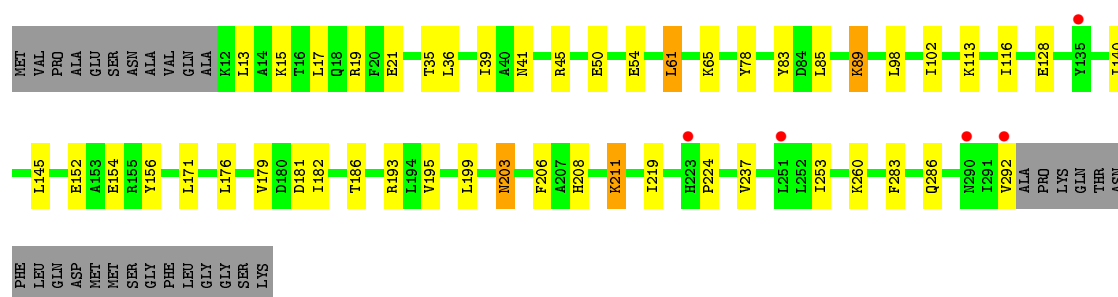
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Hg	0	0
			1	1		
3	A	1	Total	Hg	0	0
			1	1		
3	C	1	Total	Hg	0	0
			1	1		
3	E	1	Total	Hg	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	254	Total 254	O 254	0	0
4	B	44	Total 44	O 44	0	0
4	C	237	Total 237	O 237	0	0
4	D	55	Total 55	O 55	0	0
4	E	253	Total 253	O 253	0	0
4	F	56	Total 56	O 56	0	0
4	G	265	Total 265	O 265	0	0
4	H	38	Total 38	O 38	0	0

● Molecule 1: UPF0363 PROTEIN YOR164C





• Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



• Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



• Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



• Molecule 2: UBIQUITIN-LIKE PROTEIN MDY2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.27Å 118.77Å 168.38Å 90.00° 95.17° 90.00°	Depositor
Resolution (Å)	25.59 – 1.99 25.59 – 1.99	Depositor EDS
% Data completeness (in resolution range)	80.0 (25.59-1.99) 97.5 (25.59-1.99)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 1.99Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.175 , 0.206 0.190 , 0.220	Depositor DCC
R_{free} test set	5893 reflections (4.69%)	DCC
Wilson B-factor (Å ²)	18.4	Xtriage
Anisotropy	0.664	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 239485 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12001	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.35% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HG

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	0.34	0/2390	0.54	1/3226 (0.0%)
1	C	0.33	0/2375	0.49	0/3213
1	E	0.34	0/2360	0.49	0/3193
1	G	0.33	0/2386	0.50	0/3224
2	B	0.32	0/385	0.58	0/526
2	D	0.32	0/393	0.60	0/535
2	F	0.33	0/384	0.58	0/523
2	H	0.31	0/393	0.58	0/535
All	All	0.33	0/11066	0.52	1/14975 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	ASP	CB-CG-OD1	8.12	125.61	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2334	0	2275	36	0
1	C	2318	0	2231	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2303	0	2206	27	0
1	G	2329	0	2256	37	0
2	B	374	0	385	10	0
2	D	382	0	400	10	0
2	F	373	0	392	7	0
2	H	382	0	400	8	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
3	G	1	0	0	0	0
4	A	254	0	0	6	0
4	B	44	0	0	2	0
4	C	237	0	0	3	0
4	D	55	0	0	0	0
4	E	253	0	0	3	0
4	F	56	0	0	0	0
4	G	265	0	0	4	0
4	H	38	0	0	1	0
All	All	12001	0	10545	166	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 8.

All (166) 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:E:258:LYS:HD3	2:F:26:LYS:HD3	1.50	0.91
1:E:113:LYS:HE2	1:E:152:GLU:HG3	1.57	0.87
1:C:193:ARG:NH2	2:D:46:PRO:O	2.08	0.87
1:C:205:SER:O	1:C:209:GLU:HG2	1.80	0.81
1:C:113:LYS:HE2	1:C:152:GLU:HG2	1.62	0.80
2:D:24:GLU:CD	2:D:24:GLU:H	1.91	0.72
1:G:113:LYS:HE2	1:G:152:GLU:HG2	1.70	0.71
1:C:116:ILE:HD12	1:C:140:ILE:HD12	1.71	0.70
1:A:260:LYS:HG2	1:A:291:ILE:HD11	1.73	0.70
1:A:290:ASN:HB2	4:A:2253:HOH:O	1.96	0.66
1:G:253:ILE:N	1:G:253:ILE:HD12	2.12	0.64
2:H:53:LYS:O	2:H:54:GLN:HB2	1.97	0.64
1:E:175:TRP:O	1:E:178:GLN:HB2	1.98	0.63
2:H:11:GLU:O	2:H:15:LYS:HG2	1.97	0.63
1:A:175:TRP:CH2	2:B:48:LEU:HD12	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:15:LYS:O	1:C:19:ARG:HG2	2.00	0.62
1:A:177:CYS:HB3	4:A:2193:HOH:O	2.00	0.61
1:G:193:ARG:NH2	2:H:46:PRO:O	2.34	0.60
1:E:278:LYS:HD2	4:E:2243:HOH:O	2.01	0.59
1:G:41:ASN:O	1:G:45:ARG:HG3	2.01	0.59
4:A:2150:HOH:O	2:B:36:LYS:HD3	2.02	0.59
1:G:61:LEU:HD22	1:G:65:LYS:HE2	1.85	0.58
1:A:175:TRP:O	1:A:178:GLN:HB2	2.04	0.58
1:G:286:GLN:OE1	1:G:292:VAL:HA	2.03	0.58
1:C:291:ILE:HG22	1:C:292:VAL:N	2.18	0.58
2:F:35:LEU:O	2:F:35:LEU:HD23	2.04	0.57
1:G:35:THR:O	1:G:39:ILE:HG12	2.05	0.57
1:E:89:LYS:NZ	4:E:2077:HOH:O	2.37	0.56
1:A:283:PHE:CZ	2:B:42:GLY:HA2	2.41	0.56
1:E:283:PHE:CZ	2:F:42:GLY:HA2	2.41	0.56
1:E:171:LEU:HD23	1:E:171:LEU:C	2.27	0.56
1:G:182:ILE:HG23	1:G:182:ILE:O	2.06	0.55
1:C:102:ILE:HD11	1:C:119:MET:SD	2.45	0.55
1:G:128:GLU:HG3	2:H:36:LYS:HB2	1.87	0.55
1:E:144:LEU:HD13	1:E:152:GLU:HB3	1.88	0.55
1:E:50:GLU:HG2	1:E:51:HIS:N	2.21	0.55
1:C:165:MET:HE3	1:C:169:VAL:HG23	1.89	0.55
1:G:171:LEU:C	1:G:171:LEU:HD23	2.27	0.55
1:C:191:PHE:CE1	1:C:215:LEU:HD11	2.42	0.54
1:G:253:ILE:H	1:G:253:ILE:HD12	1.72	0.54
1:A:175:TRP:CZ3	2:B:48:LEU:HD12	2.43	0.54
1:C:171:LEU:C	1:C:171:LEU:HD23	2.28	0.54
1:A:283:PHE:O	1:A:286:GLN:HG2	2.07	0.54
1:A:171:LEU:C	1:A:171:LEU:HD23	2.29	0.54
1:C:113:LYS:CE	1:C:152:GLU:HG2	2.34	0.53
1:G:203:ASN:ND2	1:G:206:PHE:H	2.05	0.53
1:G:45:ARG:HD3	4:G:2045:HOH:O	2.09	0.53
1:A:61:LEU:HD12	1:A:65:LYS:HE2	1.90	0.53
2:F:11:GLU:OE2	2:F:15:LYS:HG3	2.10	0.52
1:A:23:LYS:HG2	1:E:182:ILE:HD11	1.92	0.51
1:G:283:PHE:CZ	2:H:42:GLY:HA2	2.45	0.51
1:C:253:ILE:HD11	4:C:2195:HOH:O	2.08	0.51
1:A:225:LYS:HA	1:A:225:LYS:HE3	1.91	0.51
1:G:203:ASN:C	1:G:203:ASN:HD22	2.13	0.51
1:C:260:LYS:HB3	1:C:260:LYS:NZ	2.26	0.51
1:C:291:ILE:HG22	1:C:292:VAL:H	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:129:TYR:CD1	2:F:35:LEU:HD22	2.46	0.51
1:A:190:PHE:CE1	2:B:48:LEU:HD13	2.46	0.50
1:A:234:TYR:CD1	1:A:253:ILE:HD12	2.46	0.50
1:G:195:VAL:O	1:G:199:LEU:HD13	2.11	0.50
1:E:47:LYS:NZ	1:E:47:LYS:HB2	2.26	0.50
1:A:179:VAL:O	1:A:180:ASP:C	2.50	0.50
1:A:35:THR:HG21	4:A:2015:HOH:O	2.11	0.50
1:G:208:HIS:O	1:G:211:LYS:HG3	2.12	0.50
1:C:283:PHE:CZ	2:D:42:GLY:HA2	2.46	0.50
1:G:140:ILE:HG21	1:G:156:TYR:CE2	2.47	0.50
1:C:179:VAL:HG12	1:C:180:ASP:N	2.27	0.50
1:A:87:GLU:OE2	1:C:278:LYS:HD2	2.11	0.50
1:G:89:LYS:HA	1:G:89:LYS:HE3	1.92	0.49
1:G:176:LEU:O	1:G:179:VAL:HG22	2.12	0.49
1:G:36:LEU:C	1:G:36:LEU:HD23	2.33	0.49
1:C:41:ASN:O	1:C:45:ARG:HG2	2.13	0.49
1:G:128:GLU:O	1:G:128:GLU:HG3	2.12	0.49
1:A:203:ASN:HD22	1:A:203:ASN:C	2.16	0.49
1:G:253:ILE:HD11	4:G:2230:HOH:O	2.12	0.49
1:E:113:LYS:CE	1:E:152:GLU:HG3	2.37	0.49
1:C:283:PHE:O	1:C:286:GLN:HG2	2.13	0.49
1:A:41:ASN:ND2	1:A:78:TYR:OH	2.46	0.49
1:C:116:ILE:HG23	1:C:140:ILE:HG23	1.96	0.48
1:G:15:LYS:O	1:G:19:ARG:HG3	2.13	0.48
1:A:84:ASP:CG	2:D:46:PRO:HB3	2.34	0.48
1:A:31:GLU:O	1:A:35:THR:HG22	2.13	0.48
1:C:190:PHE:CE1	2:D:48:LEU:HD23	2.49	0.48
1:C:175:TRP:CH2	2:D:48:LEU:HD22	2.49	0.48
1:C:46:SER:O	1:C:47:LYS:HB2	2.13	0.48
1:C:17:LEU:O	1:C:21:GLU:HG2	2.14	0.48
1:A:203:ASN:ND2	1:A:206:PHE:H	2.12	0.47
1:C:260:LYS:HG3	4:C:2211:HOH:O	2.14	0.47
1:C:239:PHE:CB	1:C:242:TYR:HB2	2.44	0.47
1:C:36:LEU:HD23	1:C:36:LEU:C	2.35	0.47
1:C:148:ASP:HB2	2:D:54:GLN:NE2	2.28	0.47
1:G:50:GLU:O	1:G:54:GLU:HG3	2.15	0.47
1:C:154:GLU:OE1	1:C:193:ARG:NH1	2.48	0.47
1:C:230:ASP:OD1	1:C:235:GLU:HB2	2.15	0.47
2:B:46:PRO:HB3	1:C:84:ASP:CG	2.35	0.47
1:C:90:VAL:HG21	1:C:127:SER:HB3	1.97	0.47
1:E:20:PHE:CZ	1:E:24:ILE:HD11	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LEU:CD1	1:A:65:LYS:HE2	2.45	0.47
2:B:23:THR:HG22	4:B:2006:HOH:O	2.15	0.47
1:G:253:ILE:H	1:G:253:ILE:CD1	2.28	0.46
1:G:253:ILE:N	1:G:253:ILE:CD1	2.78	0.46
1:A:90:VAL:HG21	1:A:123:SER:HA	1.97	0.46
1:A:46:SER:O	1:A:47:LYS:HB2	2.16	0.46
1:C:112:LEU:O	1:C:116:ILE:HG12	2.15	0.46
1:C:219:ILE:HD11	1:C:239:PHE:CE1	2.50	0.46
1:E:102:ILE:HD13	1:E:140:ILE:HG22	1.98	0.46
1:C:180:ASP:HA	1:G:19:ARG:HD3	1.97	0.45
4:B:2040:HOH:O	1:C:37:ARG:HD3	2.17	0.45
1:G:219:ILE:HA	1:G:224:PRO:HD2	1.99	0.45
1:C:260:LYS:HG2	1:C:289:PHE:HB3	1.99	0.45
1:C:165:MET:HE2	1:C:166:ILE:HA	1.99	0.45
1:C:260:LYS:HB3	1:C:260:LYS:HZ3	1.81	0.45
1:A:114:ASP:OD2	1:C:150:VAL:HG12	2.17	0.45
1:A:14:ALA:HB3	4:A:2001:HOH:O	2.16	0.45
1:E:46:SER:O	1:E:47:LYS:HB2	2.17	0.44
2:B:24:GLU:HA	2:B:25:PRO:HD3	1.89	0.44
1:C:179:VAL:HG12	1:C:181:ASP:H	1.82	0.44
2:F:11:GLU:O	2:F:15:LYS:HG2	2.18	0.44
1:E:154:GLU:HG3	1:E:168:TYR:CE1	2.53	0.44
2:H:15:LYS:NZ	2:H:15:LYS:HB3	2.33	0.43
1:C:260:LYS:HE2	1:C:291:ILE:HG12	2.00	0.43
1:C:175:TRP:CZ2	2:D:48:LEU:HD22	2.53	0.43
1:G:260:LYS:HG2	4:G:2238:HOH:O	2.18	0.43
1:E:23:LYS:HB3	1:E:32:ALA:HB2	2.01	0.43
1:G:17:LEU:O	1:G:21:GLU:HG2	2.19	0.43
1:A:263:PHE:HB2	1:A:289:PHE:CE2	2.53	0.43
1:G:98:LEU:O	1:G:102:ILE:HG13	2.19	0.43
1:E:205:SER:O	1:E:209:GLU:HG3	2.19	0.43
1:E:155:ARG:HA	1:E:158:MET:HE2	2.01	0.43
2:H:7:GLY:N	2:H:8:PRO:HD2	2.34	0.43
2:H:8:PRO:HG2	4:H:2002:HOH:O	2.19	0.42
1:E:239:PHE:CB	1:E:242:TYR:HB2	2.49	0.42
1:C:211:LYS:O	1:C:215:LEU:HD13	2.20	0.42
1:G:260:LYS:HB2	4:G:2239:HOH:O	2.19	0.42
1:E:17:LEU:O	1:E:21:GLU:HG2	2.19	0.42
1:A:175:TRP:CZ2	2:B:48:LEU:HD12	2.54	0.42
1:C:50:GLU:H	1:C:50:GLU:CD	2.22	0.42
1:G:154:GLU:OE1	1:G:193:ARG:NH1	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:242:TYR:HB3	1:C:245:LEU:HB2	2.02	0.42
1:G:181:ASP:HB3	1:G:186:THR:HG21	2.01	0.42
1:C:140:ILE:HG21	1:C:156:TYR:CE2	2.55	0.41
1:C:190:PHE:CE1	2:D:48:LEU:CD2	3.03	0.41
1:E:20:PHE:CE1	1:E:32:ALA:HB1	2.55	0.41
1:A:102:ILE:HD13	1:A:140:ILE:HG22	2.01	0.41
1:G:116:ILE:HG23	1:G:140:ILE:HG23	2.02	0.41
1:C:175:TRP:CZ3	2:D:48:LEU:HD22	2.55	0.41
1:C:208:HIS:NE2	1:C:256:GLN:NE2	2.69	0.41
1:A:61:LEU:O	1:A:61:LEU:HD13	2.21	0.41
1:G:219:ILE:HD11	1:G:237:VAL:HG11	2.03	0.41
1:C:165:MET:HE2	1:C:165:MET:O	2.20	0.41
1:C:215:LEU:O	1:C:219:ILE:HG12	2.20	0.41
1:E:89:LYS:HD3	4:E:2082:HOH:O	2.20	0.41
1:A:215:LEU:O	1:A:219:ILE:HG13	2.21	0.41
1:A:12:LYS:N	4:A:2001:HOH:O	2.53	0.41
1:A:242:TYR:HB3	1:A:245:LEU:HD22	2.02	0.41
1:C:209:GLU:HG2	1:C:209:GLU:H	1.70	0.41
1:C:286:GLN:HB2	4:C:2237:HOH:O	2.20	0.40
1:E:41:ASN:ND2	1:E:78:TYR:OH	2.53	0.40
1:C:197:ASN:HD22	1:C:197:ASN:HA	1.72	0.40
2:F:35:LEU:HD23	2:F:35:LEU:C	2.42	0.40
1:C:13:LEU:HD21	1:C:55:LEU:HB2	2.04	0.40
1:C:116:ILE:HG23	1:C:140:ILE:CG2	2.52	0.40
1:A:144:LEU:HD11	1:A:156:TYR:CE2	2.57	0.40
1:E:23:LYS:HD2	1:E:32:ALA:HA	2.04	0.40
1:A:130:LYS:HD3	2:B:39:THR:HA	2.04	0.40
1:E:36:LEU:HD23	1:E:36:LEU:C	2.42	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	279/312 (89%)	272 (98%)	7 (2%)	0	100	100
1	C	280/312 (90%)	272 (97%)	8 (3%)	0	100	100
1	E	278/312 (89%)	270 (97%)	8 (3%)	0	100	100
1	G	279/312 (89%)	272 (98%)	7 (2%)	0	100	100
2	B	46/59 (78%)	46 (100%)	0	0	100	100
2	D	46/59 (78%)	45 (98%)	1 (2%)	0	100	100
2	F	45/59 (76%)	45 (100%)	0	0	100	100
2	H	46/59 (78%)	45 (98%)	1 (2%)	0	100	100
All	All	1299/1484 (88%)	1267 (98%)	32 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	254/279 (91%)	249 (98%)	5 (2%)	63	65
1	C	248/279 (89%)	242 (98%)	6 (2%)	57	58
1	E	247/279 (88%)	244 (99%)	3 (1%)	78	81
1	G	252/279 (90%)	243 (96%)	9 (4%)	42	39
2	B	42/54 (78%)	41 (98%)	1 (2%)	57	58
2	D	44/54 (82%)	43 (98%)	1 (2%)	58	60
2	F	43/54 (80%)	43 (100%)	0	100	100
2	H	44/54 (82%)	43 (98%)	1 (2%)	58	60
All	All	1174/1332 (88%)	1148 (98%)	26 (2%)	60	62

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	LEU
1	A	83	TYR

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Mol	Chain	Res	Type
1	A	203	ASN
1	A	225	LYS
1	A	260	LYS
2	B	22	LEU
1	C	12	LYS
1	C	50	GLU
1	C	78	TYR
1	C	83	TYR
1	C	165	MET
1	C	245	LEU
2	D	24	GLU
1	E	54	GLU
1	E	83	TYR
1	E	260	LYS
1	G	13	LEU
1	G	61	LEU
1	G	78	TYR
1	G	83	TYR
1	G	85	LEU
1	G	89	LYS
1	G	145	LEU
1	G	203	ASN
1	G	211	LYS
2	H	41	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	HIS
1	A	41	ASN
1	A	51	HIS
1	A	58	GLN
1	A	203	ASN
1	A	268	ASN
1	A	275	GLN
1	C	22	ASN
1	C	34	GLN
1	C	197	ASN
1	C	256	GLN
1	C	275	GLN
1	C	286	GLN
2	D	54	GLN

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Mol	Chain	Res	Type
1	E	33	HIS
1	E	34	GLN
1	E	41	ASN
1	E	223	HIS
1	E	232	ASN
1	G	58	GLN
1	G	138	ASN
1	G	203	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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 ⓘ

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	281/312 (90%)	-0.02	3 (1%) 82 83	9, 20, 35, 48	0
1	C	282/312 (90%)	0.04	8 (2%) 56 57	11, 20, 37, 52	0
1	E	280/312 (89%)	0.00	5 (1%) 71 72	9, 18, 42, 52	0
1	G	281/312 (90%)	-0.00	5 (1%) 71 72	8, 19, 33, 46	0
2	B	48/59 (81%)	0.23	2 (4%) 40 41	15, 24, 41, 49	0
2	D	48/59 (81%)	0.19	2 (4%) 40 41	14, 21, 37, 44	0
2	F	47/59 (79%)	-0.08	1 (2%) 67 67	11, 19, 34, 44	0
2	H	48/59 (81%)	0.30	2 (4%) 40 41	13, 24, 38, 48	0
All	All	1315/1484 (88%)	0.03	28 (2%) 67 67	8, 20, 37, 52	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	292	VAL	4.3
1	C	292	VAL	4.1
1	A	292	VAL	4.0
1	C	135	TYR	4.0
1	E	292	VAL	3.3
2	H	54	GLN	3.2
1	C	290	ASN	3.1
1	C	11	ALA	3.0
1	G	251	LEU	2.9
1	E	22	ASN	2.6
2	H	24	GLU	2.6
1	E	24	ILE	2.6
2	B	8	PRO	2.5
2	D	54	GLN	2.4
1	C	19	ARG	2.4
1	G	290	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	135	TYR	2.4
2	D	22	LEU	2.3
1	C	286	GLN	2.3
2	B	7	GLY	2.2
1	G	223	HIS	2.2
1	A	223	HIS	2.2
2	F	7	GLY	2.2
1	A	286	GLN	2.2
1	G	135	TYR	2.2
1	C	45	ARG	2.1
1	C	28	ASP	2.1
1	E	28	ASP	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	HG	C	1293	1/1	0.96	0.07	-1.15	47,47,47,47	0
3	HG	G	1293	1/1	0.98	0.07	-1.36	51,51,51,51	0
3	HG	E	1293	1/1	0.97	0.06	-1.82	44,44,44,44	0
3	HG	A	1293	1/1	0.96	0.05	-2.44	53,53,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.