



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:14 PM GMT

PDB ID : 1JCQ
Title : CRYSTAL STRUCTURE OF HUMAN PROTEIN FARNESYLTRANSFERASE COMPLEXED WITH FARNESYL DIPHOSPHATE AND THE PEPTIDOMIMETIC INHIBITOR L-739,750
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Deposited on : 2001-06-11
Resolution : 2.30 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

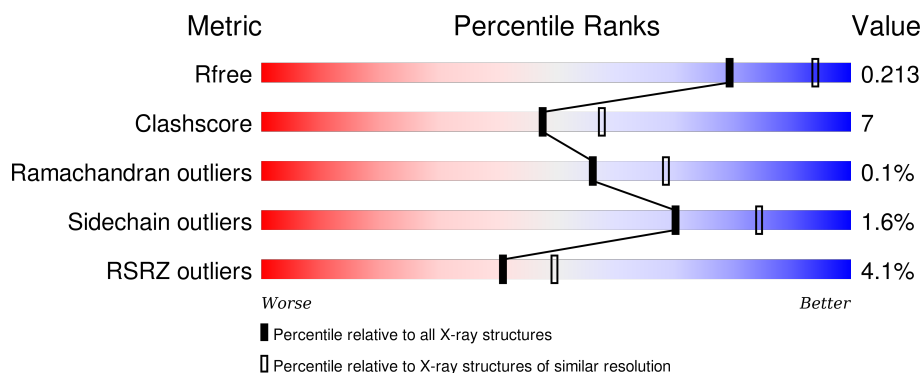
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	382	
2	B	437	

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
7	ACY	A	3002	-	-	X	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6318 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 PROTEIN FARNESYLTRANSFERASE, ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2670	1704	465	496	5			

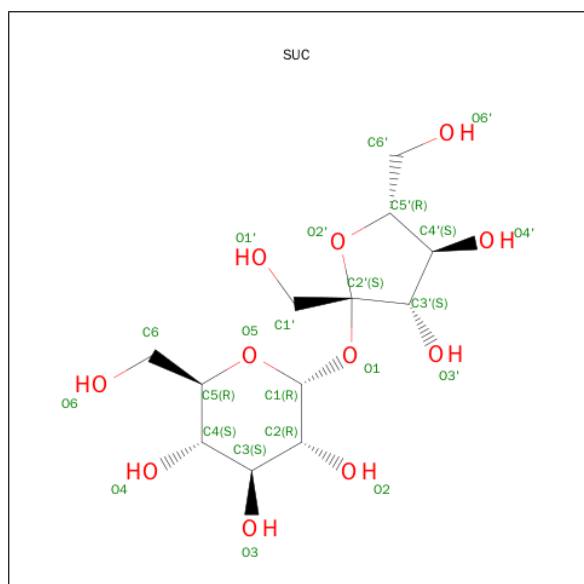
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	380	GLU	-	INSERTION	UNP P49354
A	381	GLU	-	INSERTION	UNP P49354
A	382	PHE	-	INSERTION	UNP P49354

- Molecule 2 is a protein called PROTEIN FARNESYLTRANSFERASE, BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	410	Total	C	N	O	S	0	0	0
			3228	2065	552	589	22			

- Molecule 3 is SUGAR (SUCROSE) (three-letter code: SUC) (formula: C₁₂H₂₂O₁₁).

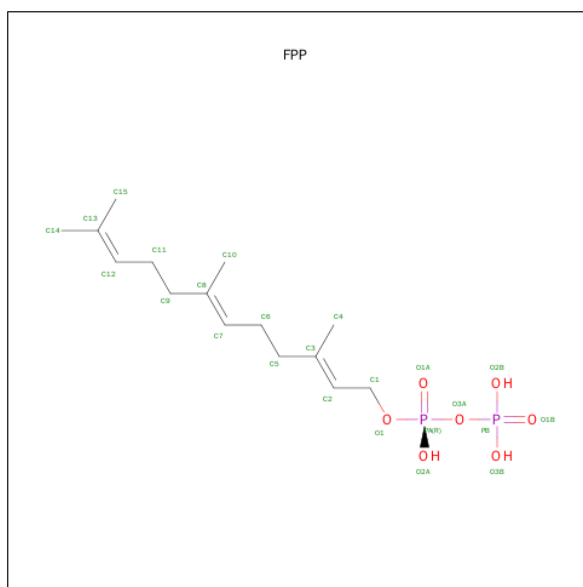


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			23	12	11		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

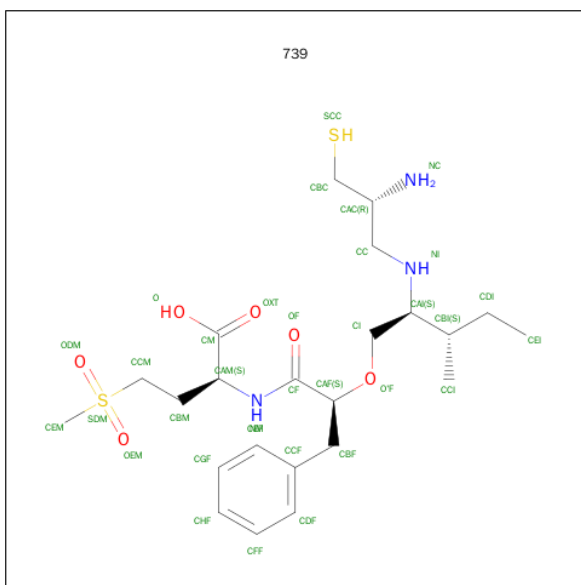
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		

- Molecule 5 is FARNESYL DIPHOSPHATE (three-letter code: FPP) (formula: C₁₅H₂₈O₇P₂).



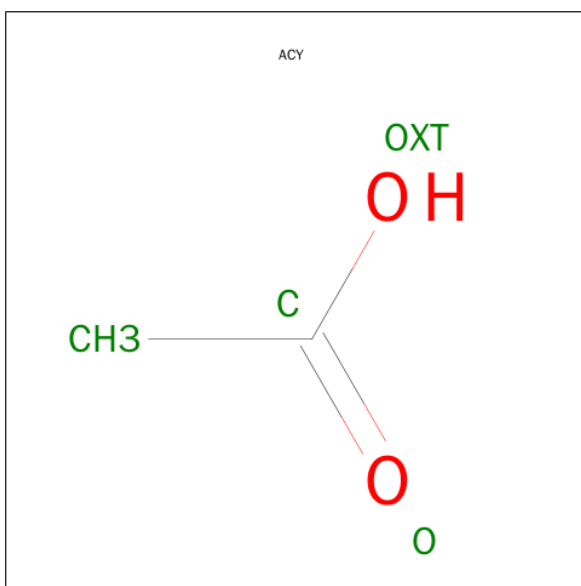
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	O	P	0	0
			24	15	7	2		

- Molecule 6 is 2(S)-{2(S)-[2(R)-AMINO-3-MERCAPTO]PROPYLAMINO-3(S)-METHYL}PENTYLOXY-3-PHENYLPROPYNYLMETHIONINE SULFONE (three-letter code: 739) (formula: C₂₃H₃₉N₃O₆S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	S	0	0
			34	23	3	6	2		

- Molecule 7 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	156	Total 156	O 156	0	0
8	B	178	Total 178	O 178	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	178.48 Å 178.48 Å 64.84 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.30 31.11 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.0 (50.00-2.30) 95.8 (31.11-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 2.29 Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.179 , 0.209 0.183 , 0.213	Depositor DCC
R_{free} test set	2527 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 56.7	EDS
Estimated twinning fraction	0.026 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 50719 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6318	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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: 739, ZN, SUC, FPP, ACY

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	0.43	0/2737	0.59	0/3717
2	B	0.46	0/3317	0.66	1/4508 (0.0%)
All	All	0.44	0/6054	0.63	1/8225 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	308	LEU	CA-CB-CG	-6.26	100.91	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2670	0	2589	41	0
2	B	3228	0	3147	48	0
3	B	23	0	22	0	0
4	B	1	0	0	0	0
5	B	24	0	25	1	0
6	B	34	0	37	2	0
7	A	4	0	3	6	0
8	A	156	0	0	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	B	178	0	0	5	0
All	All	6318	0	5823	87	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 7.

All (87) 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:A:303:GLN:O	1:A:307:SER:HB2	1.81	0.80
2:B:131:GLU:HG3	8:B:1471:HOH:O	1.82	0.80
1:A:81:ASN:HB3	8:A:1361:HOH:O	1.81	0.78
8:A:1213:HOH:O	2:B:275:GLN:HG2	1.84	0.76
2:B:152:PRO:HD3	6:B:3012:739:OEM	1.86	0.75
2:B:134:GLN:HE22	2:B:173:ASN:H	1.34	0.72
2:B:417:GLN:HG3	8:B:1223:HOH:O	1.89	0.71
1:A:300:LEU:O	1:A:303:GLN:HB2	1.91	0.71
1:A:170:HIS:HD2	7:A:3002:ACY:O	1.74	0.69
2:B:280:ARG:HE	2:B:289:GLN:HE21	1.41	0.68
2:B:74:GLN:H	2:B:344:GLN:HE22	1.45	0.63
2:B:135:SER:HB3	2:B:148:PRO:HG3	1.80	0.63
1:A:239:GLN:HA	1:A:239:GLN:HE21	1.64	0.62
1:A:346:LYS:HB2	1:A:346:LYS:NZ	2.14	0.62
2:B:135:SER:CB	2:B:148:PRO:HG3	2.30	0.62
1:A:189:ILE:HD11	1:A:205:HIS:HD2	1.64	0.61
2:B:301:SER:O	2:B:305:ALA:HB3	2.02	0.60
2:B:78:HIS:HD1	2:B:349:GLY:H	1.51	0.59
2:B:135:SER:OG	2:B:137:GLU:HG2	2.03	0.58
2:B:280:ARG:HE	2:B:289:GLN:NE2	2.00	0.58
1:A:170:HIS:HA	7:A:3002:ACY:CH3	2.34	0.57
2:B:253:PHE:HA	2:B:307:LEU:HD21	1.86	0.56
1:A:207:GLN:HE21	1:A:239:GLN:NE2	2.04	0.55
2:B:124:THR:HG22	2:B:167:GLU:OE1	2.06	0.55
2:B:312:HIS:ND1	8:B:1669:HOH:O	2.34	0.55
1:A:148:LEU:HB2	1:A:179:LEU:HD21	1.89	0.54
1:A:304:PRO:HG2	1:A:305:SER:H	1.73	0.54
2:B:239:ILE:HB	2:B:252:THR:HA	1.90	0.54
1:A:271:PRO:HG3	1:A:306:HIS:HD2	1.73	0.53
2:B:377:GLY:HA2	2:B:381:MET:O	2.09	0.53
2:B:77:LYS:HD3	2:B:346:PRO:O	2.09	0.53
2:B:86:LEU:HB2	2:B:107:ILE:HG21	1.89	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:ASN:O	1:A:298:GLN:HG3	2.09	0.52
1:A:170:HIS:CD2	7:A:3002:ACY:O	2.59	0.52
2:B:325:MET:HG2	8:B:1495:HOH:O	2.11	0.51
2:B:338:TYR:CE2	2:B:343:CYS:SG	3.04	0.50
1:A:318:ILE:O	1:A:322:MET:HG3	2.11	0.50
1:A:57:SER:HB3	1:A:60:SER:OG	2.11	0.50
1:A:329:ASN:HB3	1:A:332:ASP:HB3	1.93	0.50
1:A:346:LYS:HZ2	1:A:346:LYS:HB2	1.75	0.50
2:B:78:HIS:HD1	2:B:349:GLY:N	2.09	0.50
1:A:170:HIS:HA	7:A:3002:ACY:H3	1.93	0.50
2:B:93:TYR:CE2	2:B:96:LEU:HD12	2.48	0.49
2:B:148:PRO:HB2	2:B:180:TYR:CZ	2.47	0.49
2:B:93:TYR:CD2	2:B:96:LEU:HD12	2.48	0.49
2:B:306:GLY:O	2:B:309:PRO:HD2	2.12	0.48
1:A:239:GLN:CA	1:A:239:GLN:HE21	2.27	0.48
1:A:138:ARG:HD2	1:A:174:VAL:HG11	1.94	0.48
2:B:308:LEU:HD13	2:B:330:PHE:HB3	1.95	0.47
1:A:239:GLN:NE2	1:A:239:GLN:HA	2.29	0.47
2:B:423:PHE:CE2	2:B:424:GLU:HG2	2.49	0.47
1:A:91:ILE:O	2:B:91:ASP:HB3	2.14	0.46
2:B:175:GLU:HA	2:B:421:PRO:HB3	1.98	0.46
2:B:332:GLN:O	2:B:336:GLN:HG3	2.16	0.46
1:A:170:HIS:HA	7:A:3002:ACY:H2	1.98	0.46
2:B:350:LEU:HB2	2:B:363:THR:HA	1.98	0.45
2:B:148:PRO:HB2	2:B:180:TYR:CE1	2.52	0.45
1:A:221:GLN:HG3	8:A:1180:HOH:O	2.16	0.45
1:A:170:HIS:HE1	2:B:196:GLY:O	1.99	0.45
1:A:86:PRO:HG2	1:A:89:GLN:NE2	2.32	0.44
1:A:170:HIS:CA	7:A:3002:ACY:H2	2.48	0.44
2:B:238:GLY:HA3	2:B:247:ALA:HB1	1.99	0.44
1:A:286:ASP:HB2	8:A:1601:HOH:O	2.17	0.44
2:B:134:GLN:HB2	2:B:140:PHE:CE2	2.53	0.44
1:A:273:ASN:HB3	8:A:1205:HOH:O	2.17	0.43
2:B:370:LEU:HD23	2:B:394:LEU:HD11	2.00	0.43
1:A:227:LEU:HG	1:A:236:VAL:HG11	2.00	0.43
2:B:99:SER:O	2:B:102:TRP:HB2	2.19	0.43
2:B:202:ARG:HD2	5:B:3011:FPP:H142	2.01	0.43
1:A:170:HIS:CE1	8:A:1345:HOH:O	2.72	0.42
2:B:345:CYS:HB3	2:B:348:GLY:O	2.20	0.42
1:A:227:LEU:HG	1:A:236:VAL:CG1	2.49	0.42
1:A:96:PHE:HA	1:A:126:LEU:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:THR:HG23	1:A:188:PHE:CD1	2.55	0.42
1:A:171:HIS:O	1:A:175:LEU:HG	2.21	0.41
2:B:192:LEU:HD23	2:B:199:VAL:CG2	2.50	0.41
2:B:211:ALA:HA	2:B:216:ILE:HG12	2.01	0.41
2:B:260:VAL:HG22	2:B:265:GLU:HG2	2.02	0.41
1:A:327:CYS:O	1:A:330:LYS:HG3	2.21	0.41
1:A:225:GLN:HG2	8:A:1286:HOH:O	2.21	0.41
2:B:271:LYS:HD3	8:B:1574:HOH:O	2.20	0.40
2:B:140:PHE:HB2	2:B:153:THR:HA	2.03	0.40
1:A:271:PRO:HD2	2:B:38:ASP:CG	2.42	0.40
1:A:263:THR:HG21	1:A:280:LEU:HB2	2.02	0.40
1:A:112:ARG:O	1:A:144:LEU:HD21	2.22	0.40
6:B:3012:739:HDI2	6:B:3012:739:O'F	2.21	0.40
2:B:378:SER:O	2:B:381:MET:HB2	2.21	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	311/382 (81%)	294 (94%)	17 (6%)	0	100	100
2	B	408/437 (93%)	395 (97%)	12 (3%)	1 (0%)	52	64
All	All	719/819 (88%)	689 (96%)	29 (4%)	1 (0%)	56	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	74	GLN

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	292/344 (85%)	288 (99%)	4 (1%)	74	86
2	B	346/370 (94%)	340 (98%)	6 (2%)	68	83
All	All	638/714 (89%)	628 (98%)	10 (2%)	70	84

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	ARG
1	A	239	GLN
1	A	303	GLN
1	A	348	LYS
2	B	87	ARG
2	B	90	THR
2	B	99	SER
2	B	223	GLU
2	B	271	LYS
2	B	351	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	A	170	HIS
1	A	218	ASN
1	A	221	GLN
1	A	239	GLN
1	A	246	ASN
1	A	261	GLN
1	A	303	GLN
1	A	306	HIS
2	B	30	HIS
2	B	56	GLN
2	B	134	GLN

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Mol	Chain	Res	Type
2	B	289	GLN
2	B	318	GLN
2	B	344	GLN
2	B	375	HIS
2	B	410	GLN

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 5 ligands modelled in this entry, 1 is 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$
7	ACY	A	3002	-	1,3,3	2.26	1 (100%)	0,3,3	0.00	-
3	SUC	B	3010	-	24,24,24	1.29	1 (4%)	36,36,36	0.95	2 (5%)
5	FPP	B	3011	-	21,23,23	0.62	0	27,31,31	0.80	0
6	739	B	3012	4	31,34,34	2.33	6 (19%)	32,45,45	1.94	7 (21%)

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
7	ACY	A	3002	-	-	0/0/0/0	0/0/0/0
3	SUC	B	3010	-	-	0/12/51/51	0/2/2/2
5	FPP	B	3011	-	-	0/25/25/25	0/0/0/0
6	739	B	3012	4	-	0/34/40/40	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	3012	739	CCM-SDM	-6.10	1.70	1.78
6	B	3012	739	CC-NI	-2.42	1.40	1.47
7	A	3002	ACY	CH3-C	2.26	1.52	1.48
6	B	3012	739	O'F-CI	2.88	1.48	1.43
6	B	3012	739	CBI-CAI	3.50	1.59	1.53
6	B	3012	739	ODM-SDM	3.62	1.54	1.44
3	B	3010	SUC	C3-C2	4.58	1.64	1.52
6	B	3012	739	OEM-SDM	8.53	1.66	1.44

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	3012	739	ODM-SDM-CEM	-4.98	103.42	108.92
6	B	3012	739	CBI-CAI-NI	-3.74	105.67	111.15
6	B	3012	739	CAM-NM-CF	-2.32	119.81	123.43
6	B	3012	739	CI-O'F-CAF	-2.27	110.65	113.46
3	B	3010	SUC	C1-C2-C3	-2.26	105.52	109.97
3	B	3010	SUC	O3-C3-C2	-2.10	105.60	110.34
6	B	3012	739	ODM-SDM-CCM	2.46	109.86	108.28
6	B	3012	739	OEM-SDM-CEM	3.21	112.48	108.92
6	B	3012	739	OEM-SDM-CCM	5.97	112.10	108.28

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	3002	ACY	6	0
5	B	3011	FPP	1	0
6	B	3012	739	2	0

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	313/382 (81%)	-0.11	8 (2%) 59 68	26, 44, 65, 83	0
2	B	410/437 (93%)	0.11	22 (5%) 29 38	25, 37, 60, 86	0
All	All	723/819 (88%)	0.01	30 (4%) 41 50	25, 40, 62, 86	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	380	ALA	4.9
2	B	381	MET	4.0
2	B	120	GLN	4.0
1	A	55	PHE	4.0
2	B	15	SER	3.7
1	A	304	PRO	3.6
2	B	121	ILE	3.5
2	B	379	GLY	3.3
1	A	109	ARG	3.1
2	B	151	ALA	3.1
1	A	329	ASN	3.1
2	B	90	THR	3.0
1	A	91	ILE	2.8
2	B	424	GLU	2.5
2	B	152	PRO	2.4
2	B	346	PRO	2.4
2	B	154	TYR	2.4
2	B	155	ALA	2.4
1	A	301	ASP	2.3
2	B	277	VAL	2.3
2	B	91	ASP	2.3
1	A	84	PRO	2.2
2	B	118	ILE	2.2
2	B	32	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	378	SER	2.2
1	A	306	HIS	2.1
2	B	153	THR	2.1
2	B	101	PRO	2.1
2	B	382	LEU	2.0
2	B	105	TYR	2.0

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
7	ACY	A	3002	4/4	0.85	0.26	5.63	53,54,55,57	0
3	SUC	B	3010	23/23	0.89	0.17	1.46	49,54,57,59	0
6	739	B	3012	34/34	0.97	0.19	0.35	33,36,41,42	0
5	FPP	B	3011	24/24	0.98	0.18	0.01	22,31,33,34	0
4	ZN	B	1001	1/1	0.99	0.09	-	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.