



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 08:56 pm BST

PDB ID : 2BED
Title : Structure of FPT bound to inhibitor SCH207736
Authors : Strickland, C.
Deposited on : 2005-10-24
Resolution : 2.70 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

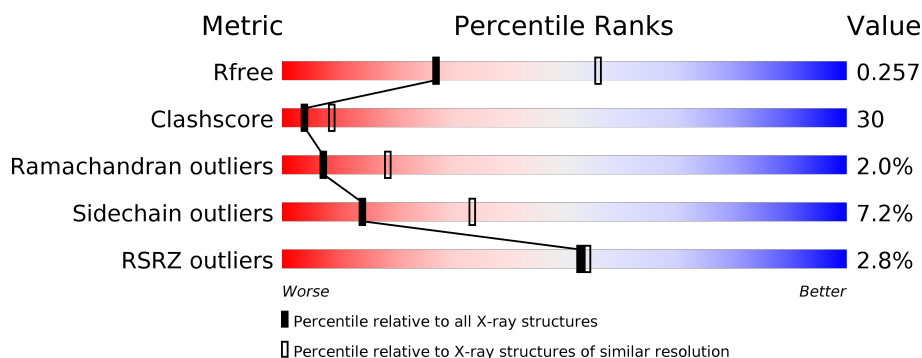
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 respectively. 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	313	<div> <div>3%</div> <div> <div></div> <div>51%</div> <div>45%</div> <div>• •</div> </div> </div>
2	B	401	<div> <div>3%</div> <div> <div></div> <div>53%</div> <div>41%</div> <div>6%</div> </div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6264 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/geranylgeranyltransferase type I alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2665	1697	466	497	5			

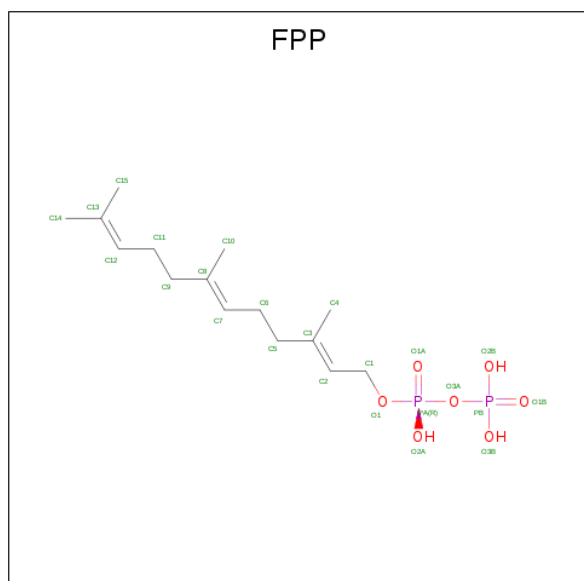
- Molecule 2 is a protein called Protein farnesyltransferase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	401	Total	C	N	O	S	0	0	0
			3154	2016	543	572	23			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

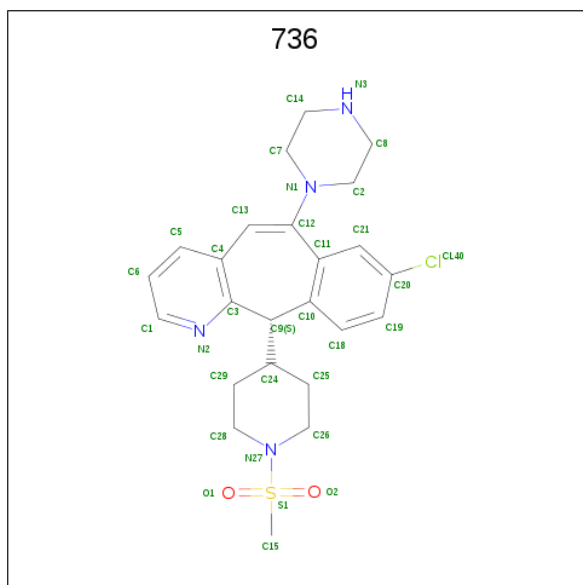
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is FARNESYL DIPHOSPHATE (three-letter code: FPP) (formula: C₁₅H₂₈O₇P₂).



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

- Molecule 5 is (11S)-8-CHLORO-11-[1-(METHYLSULFONYL)PIPERIDIN-4-YL]-6-PIPERAZIN-1-YL-11H-BENZO[5,6]CYCLOHEPTA[1,2-B]PYRIDINE (three-letter code: 736) (formula: C₂₄H₂₉ClN₄O₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	B	1	Total	C	Cl	N	O	S	0	0
			32	24	1	4	2	1		

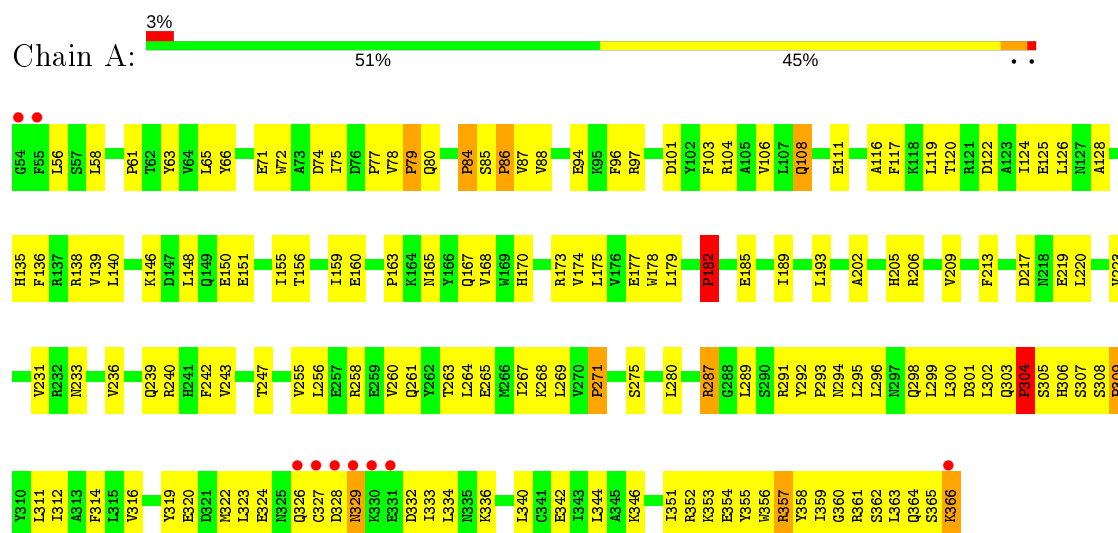
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	215	Total	O	0	0
			215	215		
6	B	173	Total	O	0	0
			173	173		

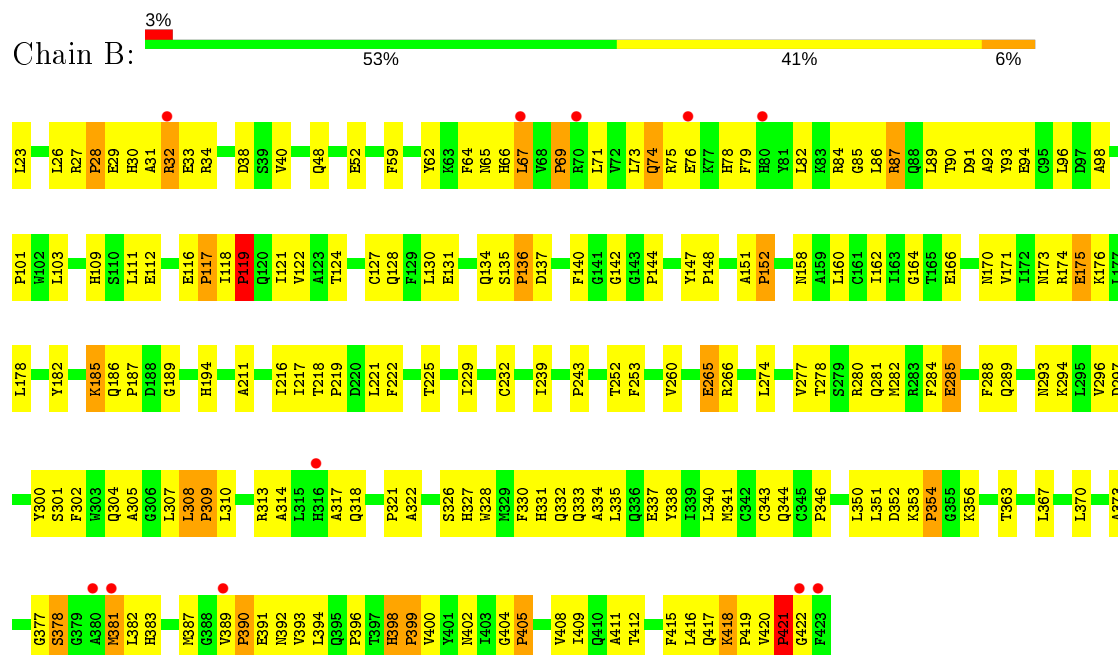
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Protein farnesyltransferase/geranylgeranyltransferase type I alpha subunit



- Molecule 2: Protein farnesyltransferase beta subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	174.02Å 174.02Å 69.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.70 25.23 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.2 (50.00-2.70) 98.4 (25.23-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.72Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.223 , 0.262 0.215 , 0.257	Depositor DCC
R_{free} test set	1662 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	55.8	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 73.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.033 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6264	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: ZN, FPP, 736

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	0/2730	0.77	1/3705 (0.0%)
2	B	0.38	0/3239	0.75	3/4397 (0.1%)
All	All	0.38	0/5969	0.76	4/8102 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	308	LEU	CA-CB-CG	-6.94	99.35	115.30
1	A	357	ARG	NE-CZ-NH2	-5.74	117.43	120.30
2	B	32	ARG	CG-CD-NE	5.59	123.54	111.80
2	B	32	ARG	N-CA-CB	-5.24	101.18	110.60

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	2665	0	2588	157	0
2	B	3154	0	3085	200	0
3	B	1	0	0	0	0
4	B	24	0	25	1	0
5	B	32	0	29	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	215	0	0	5	0
6	B	173	0	0	7	0
All	All	6264	0	5727	353	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 30.

All (353) 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)
2:B:282:MET:HG3	2:B:296:VAL:HG22	1.39	1.04
2:B:266:ARG:HH12	2:B:317:ALA:HB3	1.24	0.97
1:A:359:ILE:O	1:A:363:LEU:HD13	1.64	0.97
1:A:357:ARG:O	1:A:361:ARG:HD3	1.64	0.96
2:B:87:ARG:HG3	2:B:87:ARG:HH11	1.33	0.94
1:A:56:LEU:HD21	1:A:63:TYR:HA	1.51	0.93
2:B:151:ALA:HB3	2:B:152:PRO:HD3	1.53	0.90
2:B:285:GLU:HG2	2:B:296:VAL:HG21	1.54	0.90
1:A:342:GLU:OE1	1:A:346:LYS:HE3	1.71	0.90
2:B:66:HIS:HB3	2:B:67:LEU:HD13	1.54	0.89
2:B:189:GLY:O	2:B:225:THR:HG22	1.73	0.88
2:B:412:THR:O	2:B:416:LEU:HD13	1.76	0.86
1:A:294:ASN:O	1:A:298:GLN:HG2	1.76	0.85
2:B:124:THR:HG22	2:B:128:GLN:HE21	1.40	0.85
2:B:313:ARG:HH11	2:B:313:ARG:HG3	1.43	0.84
1:A:303:GLN:O	1:A:307:SER:HB2	1.77	0.82
2:B:331:HIS:CD2	2:B:334:ALA:H	1.98	0.80
2:B:389:VAL:HG22	2:B:391:GLU:OE1	1.82	0.80
2:B:67:LEU:HD13	2:B:67:LEU:H	1.43	0.80
1:A:80:GLN:HB2	1:A:104:ARG:NH2	1.97	0.79
2:B:266:ARG:NH1	2:B:317:ALA:HB3	1.97	0.79
2:B:131:GLU:HG2	2:B:171:VAL:HG13	1.65	0.78
2:B:308:LEU:HD12	2:B:330:PHE:CD2	2.18	0.78
1:A:351:ILE:HG23	1:A:352:ARG:HG2	1.65	0.77
1:A:316:VAL:O	1:A:320:GLU:HG3	1.83	0.77
1:A:106:VAL:HG13	1:A:111:GLU:HB3	1.67	0.77
1:A:108:GLN:HE21	1:A:108:GLN:CA	1.98	0.76
2:B:87:ARG:NH1	2:B:87:ARG:HG3	1.93	0.76
1:A:296:LEU:O	1:A:300:LEU:HD13	1.85	0.75
1:A:78:VAL:O	1:A:104:ARG:HD2	1.87	0.74
1:A:357:ARG:O	1:A:361:ARG:CD	2.36	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:301:SER:O	2:B:305:ALA:HB3	1.89	0.72
2:B:353:LYS:HB2	2:B:354:PRO:HD2	1.71	0.72
2:B:27:ARG:HG2	2:B:29:GLU:OE2	1.88	0.72
1:A:261:GLN:O	1:A:265:GLU:HG2	1.90	0.72
2:B:29:GLU:N	2:B:29:GLU:OE2	2.23	0.71
1:A:56:LEU:HD22	1:A:63:TYR:HD1	1.56	0.70
2:B:178:LEU:HD21	2:B:182:TYR:CE1	2.25	0.70
1:A:263:THR:O	1:A:267:ILE:HG13	1.91	0.70
1:A:80:GLN:HB2	1:A:104:ARG:CZ	2.23	0.69
1:A:365:SER:OG	1:A:366:LYS:NZ	2.22	0.69
2:B:185:LYS:HD3	2:B:186:GLN:N	2.08	0.69
2:B:33:GLU:O	2:B:284:PHE:HB2	1.93	0.68
1:A:304:PRO:HG2	1:A:305:SER:H	1.60	0.67
2:B:26:LEU:HD13	2:B:59:PHE:HB3	1.77	0.67
2:B:285:GLU:CG	2:B:296:VAL:HG21	2.24	0.66
2:B:381:MET:C	2:B:382:LEU:HD22	2.16	0.66
2:B:93:TYR:HD2	2:B:96:LEU:HD12	1.60	0.66
2:B:121:ILE:N	2:B:121:ILE:HD12	2.10	0.66
1:A:319:TYR:HD2	1:A:322:MET:HE2	1.62	0.65
1:A:289:LEU:HD12	1:A:289:LEU:H	1.60	0.65
1:A:56:LEU:HD22	1:A:63:TYR:CD1	2.33	0.64
2:B:253:PHE:HA	2:B:307:LEU:HD21	1.81	0.63
2:B:331:HIS:HD2	2:B:334:ALA:H	1.46	0.63
1:A:108:GLN:HE21	1:A:108:GLN:HA	1.63	0.63
1:A:96:PHE:HA	1:A:126:LEU:HD13	1.78	0.63
2:B:91:ASP:O	2:B:94:GLU:HG3	1.98	0.63
1:A:302:LEU:HD22	1:A:306:HIS:HD2	1.63	0.62
1:A:94:GLU:HG3	1:A:97:ARG:NH1	2.14	0.62
2:B:377:GLY:O	2:B:378:SER:HB2	1.98	0.62
2:B:266:ARG:HH11	2:B:318:GLN:HG3	1.64	0.62
2:B:78:HIS:O	2:B:82:LEU:HG	1.99	0.62
2:B:420:VAL:O	2:B:422:GLY:N	2.33	0.62
2:B:67:LEU:HD13	2:B:67:LEU:N	2.14	0.62
2:B:64:PHE:O	2:B:66:HIS:N	2.34	0.61
1:A:289:LEU:HD12	1:A:289:LEU:N	2.13	0.61
2:B:124:THR:CG2	2:B:128:GLN:HE21	2.13	0.61
2:B:158:ASN:O	2:B:162:ILE:HG13	2.00	0.61
1:A:302:LEU:HD22	1:A:306:HIS:CD2	2.35	0.61
2:B:31:ALA:O	2:B:34:ARG:HB2	2.00	0.61
2:B:412:THR:O	2:B:416:LEU:CD1	2.48	0.61
1:A:359:ILE:O	1:A:363:LEU:CD1	2.45	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:ARG:HD2	1:A:174:VAL:HG11	1.83	0.60
2:B:266:ARG:N	2:B:266:ARG:HD2	2.17	0.60
2:B:93:TYR:CD2	2:B:96:LEU:HD12	2.36	0.60
2:B:328:TRP:H	2:B:332:GLN:NE2	2.00	0.60
2:B:23:LEU:HD11	2:B:71:LEU:HD11	1.83	0.60
1:A:223:VAL:HG11	1:A:240:ARG:HB2	1.82	0.59
1:A:260:VAL:O	1:A:264:LEU:HG	2.03	0.58
2:B:34:ARG:HH11	2:B:34:ARG:HG2	1.68	0.58
2:B:23:LEU:CB	2:B:26:LEU:HD23	2.34	0.58
2:B:282:MET:HG3	2:B:296:VAL:CG2	2.25	0.58
1:A:255:VAL:HG13	1:A:258:ARG:NH2	2.18	0.58
2:B:173:ASN:ND2	2:B:176:LYS:HE2	2.19	0.58
1:A:327:CYS:SG	1:A:333:ILE:HD12	2.43	0.58
2:B:89:LEU:HD11	2:B:103:LEU:HD13	1.85	0.58
2:B:76:GLU:HG2	6:B:2108:HOH:O	2.04	0.58
2:B:178:LEU:C	2:B:178:LEU:HD23	2.24	0.58
2:B:302:PHE:CZ	2:B:402:ASN:HB2	2.39	0.58
2:B:178:LEU:HD13	2:B:421:PRO:HB2	1.85	0.58
2:B:350:LEU:HD22	2:B:350:LEU:N	2.19	0.57
1:A:264:LEU:O	1:A:268:LYS:HG3	2.03	0.57
4:B:1001:FPP:H102	5:B:2001:736:H71	1.84	0.57
2:B:331:HIS:NE2	2:B:333:GLN:HB3	2.19	0.57
2:B:337:GLU:HB3	2:B:341:MET:HE2	1.86	0.57
1:A:106:VAL:HG11	1:A:116:ALA:CB	2.35	0.57
2:B:266:ARG:HH12	2:B:317:ALA:CB	2.08	0.57
2:B:418:LYS:HD2	2:B:418:LYS:N	2.19	0.57
2:B:34:ARG:HE	2:B:52:GLU:HB3	1.69	0.57
2:B:29:GLU:O	2:B:32:ARG:HB2	2.04	0.57
2:B:337:GLU:HB3	2:B:341:MET:CE	2.35	0.57
2:B:174:ARG:HD2	2:B:415:PHE:HD2	1.69	0.57
1:A:312:ILE:CG2	1:A:344:LEU:HD21	2.35	0.57
1:A:56:LEU:CD2	1:A:63:TYR:HA	2.32	0.56
2:B:313:ARG:CG	2:B:313:ARG:HH11	2.16	0.56
2:B:66:HIS:HB3	2:B:67:LEU:CD1	2.32	0.56
1:A:104:ARG:O	1:A:108:GLN:HB2	2.06	0.56
1:A:108:GLN:NE2	1:A:108:GLN:CA	2.68	0.56
1:A:108:GLN:NE2	1:A:108:GLN:HA	2.20	0.56
2:B:127:CYS:O	2:B:131:GLU:HG3	2.06	0.56
1:A:58:LEU:HD21	1:A:122:ASP:HB3	1.86	0.56
1:A:156:THR:O	1:A:160:GLU:HG2	2.06	0.56
2:B:79:PHE:HD1	2:B:116:GLU:HG3	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ILE:HG21	1:A:206:ARG:HB2	1.88	0.55
2:B:338:TYR:CE2	2:B:343:CYS:SG	2.99	0.55
2:B:118:ILE:N	2:B:118:ILE:HD12	2.21	0.55
2:B:297:ASP:HB3	2:B:300:TYR:CD1	2.42	0.55
2:B:75:ARG:NH2	2:B:393:VAL:O	2.35	0.55
2:B:335:LEU:HD23	2:B:373:ALA:HB2	1.88	0.55
2:B:398:HIS:ND1	2:B:399:PRO:HD2	2.22	0.55
2:B:175:GLU:H	2:B:175:GLU:CD	2.10	0.54
2:B:151:ALA:HB3	2:B:152:PRO:CD	2.34	0.54
2:B:281:GLN:HB2	2:B:288:PHE:CE2	2.42	0.54
2:B:331:HIS:CD2	2:B:333:GLN:HB3	2.43	0.54
2:B:296:VAL:HG23	6:B:2002:HOH:O	2.07	0.54
1:A:106:VAL:HG11	1:A:116:ALA:HB1	1.89	0.54
1:A:148:LEU:HB2	1:A:179:LEU:HD21	1.88	0.54
2:B:308:LEU:HB2	2:B:309:PRO:HD3	1.90	0.54
2:B:398:HIS:CD2	2:B:408:VAL:HG21	2.43	0.54
2:B:166:GLU:O	2:B:170:ASN:OD1	2.26	0.53
1:A:170:HIS:O	1:A:174:VAL:HG23	2.08	0.53
1:A:319:TYR:CE1	1:A:336:LYS:HG2	2.44	0.53
1:A:289:LEU:CD1	1:A:289:LEU:H	2.20	0.53
2:B:173:ASN:HD21	2:B:176:LYS:HE2	1.73	0.53
1:A:155:ILE:O	1:A:159:ILE:HG13	2.08	0.53
1:A:304:PRO:HG2	1:A:305:SER:N	2.23	0.53
2:B:174:ARG:HD2	2:B:415:PHE:CD2	2.44	0.53
2:B:265:GLU:CD	2:B:265:GLU:H	2.13	0.53
2:B:296:VAL:HG12	2:B:297:ASP:N	2.24	0.53
1:A:300:LEU:HD12	1:A:300:LEU:N	2.24	0.52
2:B:135:SER:OG	2:B:136:PRO:HD2	2.08	0.52
1:A:239:GLN:O	1:A:243:VAL:HG23	2.09	0.52
2:B:310:LEU:HD21	2:B:400:VAL:HG22	1.90	0.52
1:A:135:HIS:O	1:A:139:VAL:HG23	2.10	0.52
1:A:365:SER:O	1:A:366:LYS:HE3	2.08	0.52
1:A:120:THR:O	1:A:124:ILE:HG13	2.08	0.52
2:B:280:ARG:NE	2:B:280:ARG:HA	2.25	0.52
1:A:354:GLU:OE2	2:B:327:HIS:HB2	2.10	0.52
1:A:329:ASN:HB3	1:A:332:ASP:HB3	1.91	0.52
1:A:94:GLU:HG3	1:A:97:ARG:HH12	1.75	0.52
2:B:82:LEU:HB2	2:B:111:LEU:HD21	1.92	0.52
2:B:134:GLN:HB2	2:B:140:PHE:CE2	2.44	0.51
1:A:151:GLU:HG3	1:A:175:LEU:HD11	1.92	0.51
1:A:362:SER:O	1:A:366:LYS:HD2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:89:LEU:CD1	2:B:103:LEU:HD13	2.40	0.51
1:A:65:LEU:HD11	6:A:581:HOH:O	2.11	0.51
1:A:322:MET:HB3	1:A:327:CYS:SG	2.51	0.51
1:A:84:PRO:C	1:A:86:PRO:HD3	2.31	0.50
1:A:165:ASN:HB3	1:A:168:VAL:HG22	1.94	0.50
1:A:319:TYR:O	1:A:323:LEU:HD13	2.11	0.50
1:A:117:PHE:CE2	1:A:146:LYS:HE2	2.47	0.50
1:A:287:ARG:HG2	1:A:291:ARG:HD2	1.93	0.50
1:A:364:GLN:NE2	1:A:364:GLN:HA	2.26	0.50
2:B:333:GLN:HG3	2:B:387:MET:HE2	1.93	0.50
2:B:98:ALA:O	2:B:142:GLY:HA3	2.12	0.50
1:A:182:PRO:HG3	1:A:213:PHE:CD2	2.47	0.50
1:A:220:LEU:HB2	1:A:243:VAL:HG11	1.94	0.50
2:B:265:GLU:HG2	2:B:314:ALA:HB2	1.94	0.50
2:B:326:SER:HB2	2:B:383:HIS:ND1	2.27	0.50
2:B:260:VAL:HA	2:B:265:GLU:OE2	2.11	0.49
1:A:296:LEU:O	1:A:300:LEU:CD1	2.57	0.49
1:A:357:ARG:O	1:A:361:ARG:CG	2.60	0.49
2:B:124:THR:HG22	2:B:128:GLN:NE2	2.19	0.49
2:B:27:ARG:O	2:B:30:HIS:HB3	2.13	0.49
1:A:312:ILE:HG21	1:A:344:LEU:HD21	1.94	0.49
1:A:319:TYR:HE1	1:A:336:LYS:HG2	1.76	0.49
2:B:119:PRO:HB2	2:B:122:VAL:HG12	1.94	0.49
1:A:243:VAL:O	1:A:247:THR:HG23	2.13	0.49
1:A:56:LEU:CD2	1:A:63:TYR:HD1	2.25	0.49
1:A:351:ILE:O	2:B:331:HIS:HB2	2.13	0.49
2:B:74:GLN:HA	2:B:74:GLN:NE2	2.28	0.49
2:B:62:TYR:C	2:B:64:PHE:H	2.16	0.49
2:B:71:LEU:N	2:B:71:LEU:HD12	2.27	0.49
2:B:239:ILE:HB	2:B:252:THR:HA	1.94	0.49
2:B:194:HIS:HB2	6:B:2021:HOH:O	2.12	0.49
2:B:121:ILE:N	2:B:121:ILE:CD1	2.76	0.48
2:B:73:LEU:HD12	2:B:344:GLN:OE1	2.13	0.48
1:A:301:ASP:O	1:A:304:PRO:HD2	2.14	0.48
1:A:135:HIS:CD2	2:B:147:TYR:HE1	2.32	0.48
2:B:178:LEU:HD11	2:B:218:THR:HG23	1.96	0.48
2:B:69:PRO:HG2	6:B:2074:HOH:O	2.13	0.48
1:A:319:TYR:HD2	1:A:322:MET:CE	2.25	0.48
1:A:328:ASP:O	1:A:329:ASN:HB2	2.14	0.48
1:A:361:ARG:NH1	2:B:322:ALA:O	2.47	0.48
2:B:23:LEU:HB3	2:B:26:LEU:HD23	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:GLY:O	1:A:364:GLN:HB2	2.13	0.48
2:B:117:PRO:O	2:B:119:PRO:HD3	2.14	0.48
2:B:27:ARG:HG2	2:B:29:GLU:CD	2.33	0.48
2:B:64:PHE:O	2:B:66:HIS:ND1	2.42	0.48
1:A:193:LEU:HD21	1:A:202:ALA:HB3	1.95	0.47
2:B:308:LEU:CD1	2:B:330:PHE:HD2	2.28	0.47
1:A:265:GLU:O	1:A:269:LEU:HD23	2.14	0.47
1:A:74:ASP:OD1	1:A:75:ILE:HG13	2.13	0.47
1:A:287:ARG:HD3	1:A:292:TYR:OH	2.15	0.47
1:A:239:GLN:HA	1:A:239:GLN:OE1	2.14	0.47
1:A:275:SER:OG	2:B:293:ASN:ND2	2.48	0.47
2:B:308:LEU:CD1	2:B:330:PHE:CD2	2.93	0.47
1:A:103:PHE:HE1	1:A:120:THR:HG22	1.80	0.47
1:A:302:LEU:CD2	1:A:306:HIS:HD2	2.27	0.47
1:A:362:SER:C	1:A:364:GLN:H	2.17	0.47
2:B:134:GLN:HE22	2:B:173:ASN:H	1.62	0.47
1:A:185:GLU:O	1:A:189:ILE:HG13	2.15	0.47
1:A:233:ASN:CG	1:A:236:VAL:HG23	2.35	0.47
2:B:404:GLY:O	2:B:408:VAL:HG23	2.15	0.47
2:B:304:GLN:HA	2:B:304:GLN:OE1	2.15	0.46
2:B:308:LEU:HD23	2:B:308:LEU:HA	1.53	0.46
2:B:277:VAL:O	2:B:280:ARG:HB2	2.15	0.46
2:B:90:THR:C	2:B:92:ALA:H	2.18	0.46
2:B:75:ARG:NE	2:B:391:GLU:O	2.46	0.46
1:A:223:VAL:HG13	1:A:236:VAL:CG1	2.46	0.46
1:A:66:TYR:CE2	1:A:119:LEU:HD13	2.50	0.46
1:A:336:LYS:O	1:A:340:LEU:HG	2.16	0.46
1:A:344:LEU:HD13	1:A:356:TRP:CE2	2.51	0.46
1:A:167:GLN:H	1:A:167:GLN:CD	2.19	0.46
2:B:211:ALA:HA	2:B:216:ILE:HG12	1.98	0.46
2:B:66:HIS:CB	2:B:67:LEU:HD13	2.37	0.46
1:A:319:TYR:HA	1:A:322:MET:HE2	1.98	0.46
2:B:353:LYS:HB2	2:B:354:PRO:CD	2.45	0.46
2:B:281:GLN:HB2	2:B:288:PHE:CZ	2.52	0.45
2:B:310:LEU:CD2	2:B:400:VAL:HG22	2.46	0.45
1:A:231:VAL:HG22	1:A:231:VAL:O	2.15	0.45
1:A:328:ASP:O	1:A:329:ASN:CB	2.63	0.45
2:B:119:PRO:O	2:B:122:VAL:HG12	2.15	0.45
2:B:26:LEU:N	2:B:26:LEU:HD22	2.31	0.45
2:B:87:ARG:CG	2:B:87:ARG:NH1	2.69	0.45
1:A:303:GLN:HA	1:A:303:GLN:OE1	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LEU:H	1:A:300:LEU:CD1	2.30	0.45
1:A:173:ARG:O	1:A:177:GLU:HG3	2.16	0.45
2:B:370:LEU:HD23	2:B:394:LEU:HD11	1.98	0.45
2:B:96:LEU:HD13	2:B:103:LEU:HD21	1.99	0.45
2:B:130:LEU:HD12	2:B:160:LEU:HD21	1.99	0.45
2:B:266:ARG:HD2	2:B:266:ARG:H	1.81	0.45
2:B:313:ARG:CG	2:B:313:ARG:NH1	2.79	0.45
2:B:84:ARG:O	2:B:85:GLY:C	2.55	0.45
2:B:75:ARG:NH2	2:B:391:GLU:O	2.49	0.44
1:A:268:LYS:O	1:A:271:PRO:HD3	2.17	0.44
1:A:87:VAL:O	1:A:88:VAL:C	2.56	0.44
2:B:178:LEU:O	2:B:178:LEU:HD23	2.17	0.44
2:B:23:LEU:CD1	2:B:71:LEU:HD11	2.45	0.44
2:B:389:VAL:CG2	2:B:391:GLU:OE1	2.60	0.44
2:B:71:LEU:HB3	2:B:340:LEU:HD13	1.98	0.44
1:A:287:ARG:HE	1:A:291:ARG:CD	2.30	0.44
1:A:320:GLU:O	1:A:324:GLU:HG3	2.17	0.44
1:A:336:LYS:HB2	1:A:336:LYS:HE3	1.70	0.44
2:B:266:ARG:HE	2:B:318:GLN:NE2	2.16	0.44
1:A:346:LYS:HA	6:A:496:HOH:O	2.17	0.44
2:B:151:ALA:CB	2:B:152:PRO:HD3	2.37	0.44
2:B:84:ARG:O	2:B:87:ARG:N	2.43	0.44
2:B:389:VAL:CG2	2:B:390:PRO:HD2	2.48	0.44
1:A:56:LEU:HD21	1:A:63:TYR:CA	2.33	0.44
2:B:232:CYS:CB	2:B:239:ILE:HG23	2.47	0.44
1:A:342:GLU:HA	1:A:342:GLU:OE2	2.18	0.44
2:B:232:CYS:HB2	2:B:239:ILE:HG23	2.00	0.44
2:B:381:MET:O	2:B:382:LEU:HD22	2.17	0.44
1:A:85:SER:N	1:A:86:PRO:HD3	2.33	0.43
2:B:23:LEU:HB2	2:B:26:LEU:HD23	1.99	0.43
2:B:34:ARG:HG2	2:B:34:ARG:NH1	2.32	0.43
2:B:73:LEU:O	2:B:75:ARG:N	2.42	0.43
1:A:319:TYR:HA	1:A:322:MET:CE	2.48	0.43
1:A:219:GLU:O	1:A:223:VAL:HG23	2.18	0.43
1:A:296:LEU:CD1	1:A:300:LEU:HD11	2.47	0.43
1:A:300:LEU:HD12	1:A:300:LEU:H	1.83	0.43
2:B:412:THR:C	2:B:416:LEU:HD13	2.39	0.43
1:A:223:VAL:HG13	1:A:236:VAL:HG12	2.00	0.43
2:B:405:PRO:O	2:B:409:ILE:HG13	2.17	0.43
2:B:412:THR:HG21	6:B:2173:HOH:O	2.18	0.43
1:A:219:GLU:HA	1:A:219:GLU:OE1	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:350:LEU:HD23	2:B:363:THR:HG23	2.01	0.43
1:A:205:HIS:O	1:A:209:VAL:HG23	2.18	0.43
1:A:66:TYR:CZ	1:A:119:LEU:HD13	2.53	0.43
1:A:136:PHE:CE2	1:A:140:LEU:HD11	2.54	0.43
2:B:284:PHE:HB3	2:B:285:GLU:OE2	2.18	0.43
2:B:29:GLU:HA	2:B:32:ARG:HD3	2.00	0.43
2:B:389:VAL:HG13	2:B:392:ASN:ND2	2.34	0.43
1:A:71:GLU:HG2	1:A:72:TRP:CD1	2.54	0.42
2:B:351:LEU:HD22	2:B:356:LYS:O	2.20	0.42
1:A:289:LEU:N	1:A:289:LEU:CD1	2.80	0.42
1:A:353:LYS:O	1:A:357:ARG:HB2	2.20	0.42
2:B:82:LEU:CB	2:B:111:LEU:HD21	2.48	0.42
2:B:178:LEU:HB2	2:B:216:ILE:HD12	2.01	0.42
2:B:327:HIS:HB3	2:B:332:GLN:NE2	2.34	0.42
2:B:327:HIS:HB3	2:B:332:GLN:HE21	1.84	0.42
1:A:287:ARG:HE	1:A:291:ARG:HH21	1.66	0.42
1:A:295:LEU:O	1:A:296:LEU:C	2.58	0.42
1:A:299:LEU:HB3	1:A:311:LEU:HD11	2.01	0.42
1:A:319:TYR:O	1:A:323:LEU:CD1	2.67	0.42
2:B:274:LEU:O	2:B:278:THR:HG23	2.19	0.42
2:B:411:ALA:HB2	6:B:2105:HOH:O	2.18	0.42
2:B:86:LEU:CD2	2:B:122:VAL:HG21	2.49	0.42
1:A:300:LEU:HA	1:A:303:GLN:HG2	2.02	0.42
2:B:351:LEU:HD13	2:B:353:LYS:O	2.20	0.42
1:A:355:TYR:O	1:A:358:TYR:HB3	2.19	0.42
1:A:322:MET:HB3	1:A:327:CYS:HB3	2.01	0.42
2:B:48:GLN:NE2	2:B:48:GLN:HA	2.34	0.42
1:A:256:LEU:O	1:A:260:VAL:HG23	2.20	0.42
2:B:90:THR:C	2:B:92:ALA:N	2.73	0.42
1:A:319:TYR:CD2	1:A:322:MET:CE	3.03	0.41
1:A:56:LEU:C	1:A:56:LEU:HD23	2.40	0.41
2:B:135:SER:C	2:B:137:ASP:H	2.23	0.41
2:B:350:LEU:N	2:B:350:LEU:CD2	2.83	0.41
1:A:289:LEU:HD23	1:A:314:PHE:CE2	2.55	0.41
1:A:323:LEU:HD21	1:A:334:LEU:HD13	2.00	0.41
1:A:96:PHE:CD2	1:A:96:PHE:C	2.93	0.41
1:A:213:PHE:CD1	1:A:213:PHE:N	2.88	0.41
2:B:117:PRO:C	2:B:118:ILE:HD12	2.41	0.41
2:B:216:ILE:HA	2:B:420:VAL:HG13	2.02	0.41
2:B:109:HIS:O	2:B:112:GLU:HB3	2.21	0.41
2:B:67:LEU:H	2:B:67:LEU:CD1	2.25	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:285:GLU:HG2	2:B:296:VAL:CG2	2.38	0.41
2:B:38:ASP:O	2:B:40:VAL:HG23	2.20	0.41
2:B:221:LEU:HD23	2:B:222:PHE:CE1	2.56	0.41
1:A:79:PRO:HA	1:A:101:ASP:OD1	2.20	0.41
2:B:74:GLN:HA	2:B:74:GLN:HE21	1.85	0.41
1:A:263:THR:HG21	1:A:280:LEU:HB2	2.02	0.41
1:A:294:ASN:C	1:A:298:GLN:HG2	2.38	0.41
1:A:366:LYS:HD2	1:A:366:LYS:N	2.35	0.41
2:B:294:LYS:HD2	2:B:294:LYS:HA	1.79	0.41
2:B:73:LEU:H	2:B:392:ASN:ND2	2.18	0.41
2:B:398:HIS:HD2	2:B:408:VAL:HG21	1.82	0.41
1:A:138:ARG:HH11	1:A:138:ARG:HD3	1.70	0.41
1:A:138:ARG:HG3	1:A:178:TRP:CH2	2.55	0.41
2:B:266:ARG:CD	2:B:266:ARG:H	2.35	0.41
1:A:125:GLU:HB2	6:A:404:HOH:O	2.20	0.40
1:A:242:PHE:HA	6:A:372:HOH:O	2.20	0.40
1:A:300:LEU:N	1:A:300:LEU:CD1	2.84	0.40
2:B:377:GLY:O	2:B:378:SER:CB	2.64	0.40
1:A:287:ARG:NH1	6:A:478:HOH:O	2.54	0.40
1:A:333:ILE:O	1:A:336:LYS:HB3	2.21	0.40
2:B:86:LEU:HD21	2:B:122:VAL:HG21	2.03	0.40
2:B:225:THR:O	2:B:229:ILE:HG13	2.21	0.40
2:B:266:ARG:CD	2:B:266:ARG:N	2.82	0.40
2:B:280:ARG:O	2:B:289:GLN:HG2	2.21	0.40
2:B:417:GLN:HB2	2:B:418:LYS:HD2	2.03	0.40
2:B:67:LEU:N	2:B:67:LEU:CD1	2.82	0.40
1:A:326:GLN:N	1:A:326:GLN:CD	2.74	0.40
1:A:308:SER:HB2	1:A:309:PRO:HD2	2.04	0.40
2:B:217:ILE:HB	6:B:2087:HOH:O	2.22	0.40
2:B:351:LEU:HD12	2:B:351:LEU:C	2.42	0.40
1:A:362:SER:C	1:A:364:GLN:N	2.75	0.40
1:A:80:GLN:HB2	1:A:104:ARG:HH21	1.81	0.40
2:B:29:GLU:HB3	2:B:32:ARG:NH1	2.37	0.40
2:B:327:HIS:HA	2:B:332:GLN:HE22	1.85	0.40
2:B:363:THR:O	2:B:367:LEU:HG	2.22	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	311/313 (99%)	273 (88%)	32 (10%)	6 (2%)	8	20
2	B	399/401 (100%)	359 (90%)	32 (8%)	8 (2%)	7	19
All	All	710/714 (99%)	632 (89%)	64 (9%)	14 (2%)	7	19

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	217	ASP
1	A	287	ARG
1	A	329	ASN
2	B	65	ASN
2	B	378	SER
2	B	421	PRO
2	B	74	GLN
2	B	352	ASP
1	A	128	ALA
1	A	304	PRO
2	B	119	PRO
1	A	182	PRO
2	B	28	PRO
2	B	164	GLY

5.3.2 Protein sidechains [i](#)

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	291/291 (100%)	277 (95%)	14 (5%)	25	53
2	B	338/338 (100%)	307 (91%)	31 (9%)	9	21
All	All	629/629 (100%)	584 (93%)	45 (7%)	14	34

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	PRO
1	A	77	PRO
1	A	79	PRO
1	A	84	PRO
1	A	86	PRO
1	A	108	GLN
1	A	150	GLU
1	A	163	PRO
1	A	182	PRO
1	A	271	PRO
1	A	293	PRO
1	A	304	PRO
1	A	309	PRO
1	A	366	LYS
2	B	28	PRO
2	B	67	LEU
2	B	69	PRO
2	B	87	ARG
2	B	101	PRO
2	B	117	PRO
2	B	119	PRO
2	B	136	PRO
2	B	144	PRO
2	B	148	PRO
2	B	152	PRO
2	B	175	GLU
2	B	185	LYS
2	B	187	PRO
2	B	219	PRO
2	B	243	PRO
2	B	265	GLU
2	B	285	GLU
2	B	309	PRO
2	B	321	PRO
2	B	346	PRO

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Mol	Chain	Res	Type
2	B	354	PRO
2	B	381	MET
2	B	390	PRO
2	B	396	PRO
2	B	398	HIS
2	B	399	PRO
2	B	405	PRO
2	B	418	LYS
2	B	419	PRO
2	B	421	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	GLN
1	A	135	HIS
1	A	149	GLN
1	A	246	ASN
1	A	285	GLN
1	A	294	ASN
1	A	306	HIS
1	A	325	ASN
1	A	364	GLN
2	B	48	GLN
2	B	74	GLN
2	B	88	GLN
2	B	128	GLN
2	B	134	GLN
2	B	146	GLN
2	B	275	GLN
2	B	293	ASN
2	B	318	GLN
2	B	327	HIS
2	B	331	HIS
2	B	332	GLN
2	B	333	GLN
2	B	392	ASN
2	B	410	GLN
2	B	417	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 3 ligands modelled in this entry, 1 is 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
5	736	B	2001	-	36,36,36	2.22	11 (30%)	40,53,53	1.36	6 (15%)
4	FPP	B	1001	-	21,23,23	0.67	1 (4%)	27,31,31	1.10	3 (11%)

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
5	736	B	2001	-	-	2/10/48/48	0/5/5/5
4	FPP	B	1001	-	-	9/25/25/25	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	2001	736	S1-N27	8.30	1.73	1.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	2001	736	C11-C10	4.31	1.45	1.40
5	B	2001	736	C11-C12	4.25	1.51	1.47
5	B	2001	736	C12-N1	3.43	1.48	1.37
5	B	2001	736	C21-C20	2.68	1.42	1.38
5	B	2001	736	C18-C10	2.65	1.43	1.39
5	B	2001	736	C19-C20	2.55	1.42	1.38
5	B	2001	736	C4-C13	2.47	1.50	1.46
5	B	2001	736	C13-C12	2.46	1.38	1.35
5	B	2001	736	C21-C11	2.25	1.43	1.39
4	B	1001	FPP	O1-C1	2.23	1.47	1.43
5	B	2001	736	C19-C18	2.19	1.42	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	2001	736	C26-N27-S1	3.11	120.13	116.30
5	B	2001	736	C28-N27-S1	3.06	120.07	116.30
5	B	2001	736	O2-S1-N27	-3.05	104.47	107.03
5	B	2001	736	O1-S1-N27	-2.92	104.58	107.03
5	B	2001	736	C1-N2-C3	2.71	122.50	116.78
4	B	1001	FPP	C6-C7-C8	-2.34	122.03	127.66
4	B	1001	FPP	C15-C13-C14	2.23	119.53	114.60
4	B	1001	FPP	C10-C8-C9	2.21	118.99	115.27
5	B	2001	736	C6-C1-N2	-2.00	120.16	123.43

There are no chirality outliers.

All (11) torsion outliers are listed below:

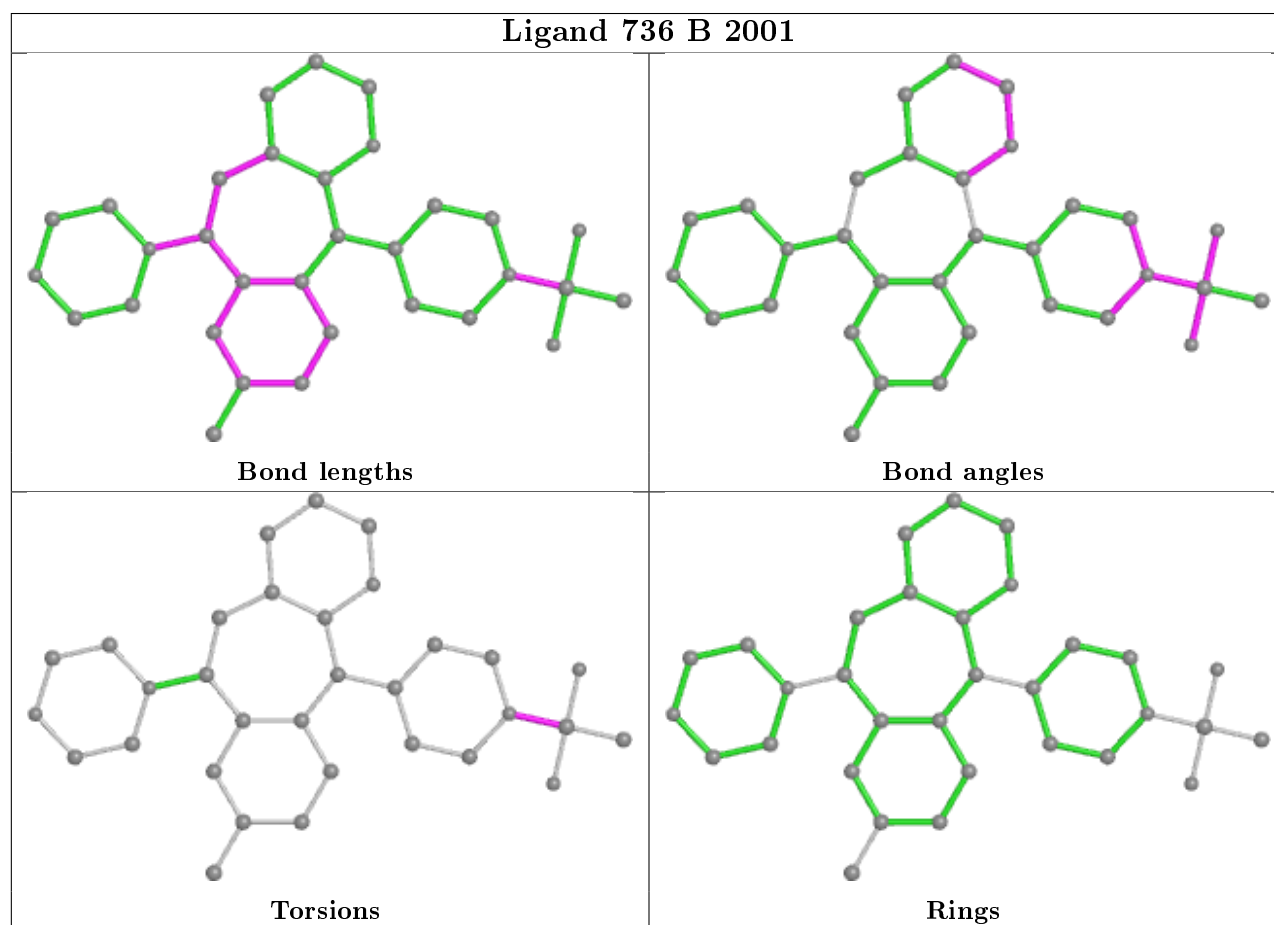
Mol	Chain	Res	Type	Atoms
5	B	2001	736	C28-N27-S1-C15
5	B	2001	736	C28-N27-S1-O2
4	B	1001	FPP	C1-C2-C3-C4
4	B	1001	FPP	C1-C2-C3-C5
4	B	1001	FPP	C6-C7-C8-C10
4	B	1001	FPP	C6-C7-C8-C9
4	B	1001	FPP	PB-O3A-PA-O2A
4	B	1001	FPP	C4-C3-C5-C6
4	B	1001	FPP	C2-C3-C5-C6
4	B	1001	FPP	PB-O3A-PA-O1A
4	B	1001	FPP	C1-O1-PA-O1A

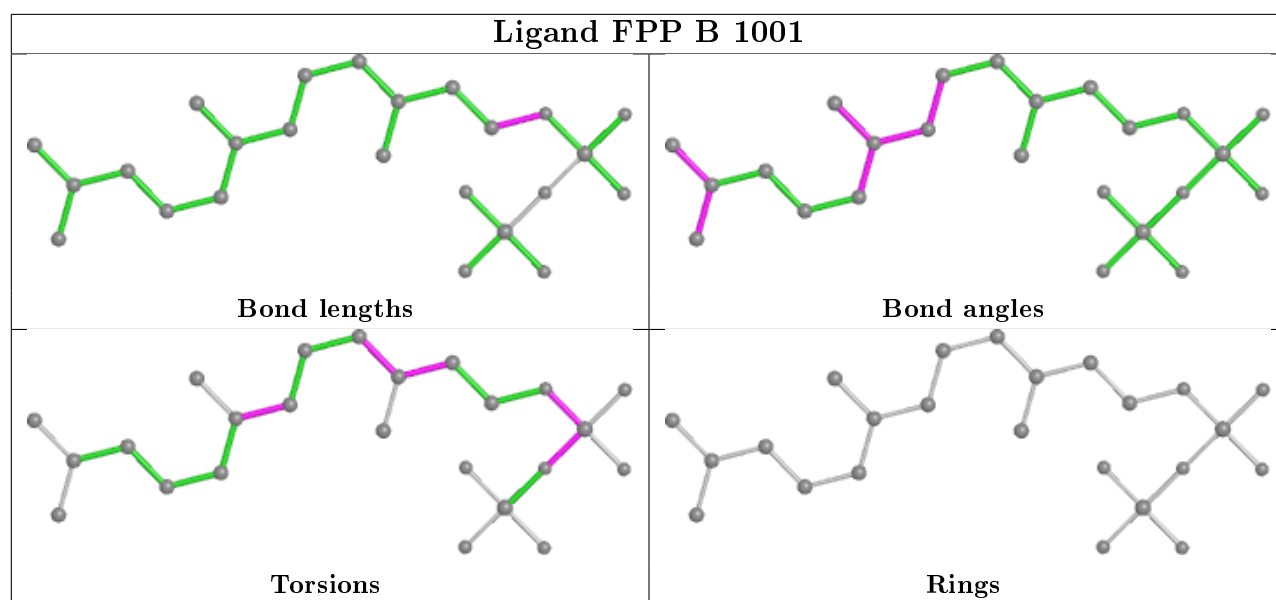
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	2001	736	1	0
4	B	1001	FPP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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

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	313/313 (100%)	-0.20	9 (2%)	51	52	26, 53, 94, 111	0
2	B	401/401 (100%)	-0.02	11 (2%)	54	55	24, 51, 80, 110	0
All	All	714/714 (100%)	-0.10	20 (2%)	53	54	24, 52, 85, 111	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	423	PHE	8.4
1	A	54	GLY	5.6
1	A	326	GLN	4.1
2	B	381	MET	3.7
1	A	330	LYS	3.3
2	B	422	GLY	3.3
2	B	389	VAL	3.2
1	A	55	PHE	3.2
1	A	327	CYS	2.9
2	B	380	ALA	2.8
1	A	331	GLU	2.7
2	B	70	ARG	2.7
1	A	328	ASP	2.5
1	A	366	LYS	2.4
2	B	32	ARG	2.4
1	A	329	ASN	2.2
2	B	67	LEU	2.2
2	B	316	HIS	2.1
2	B	76	GLU	2.1
2	B	80	HIS	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](#)

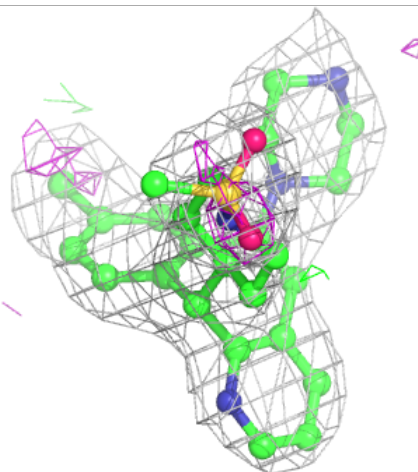
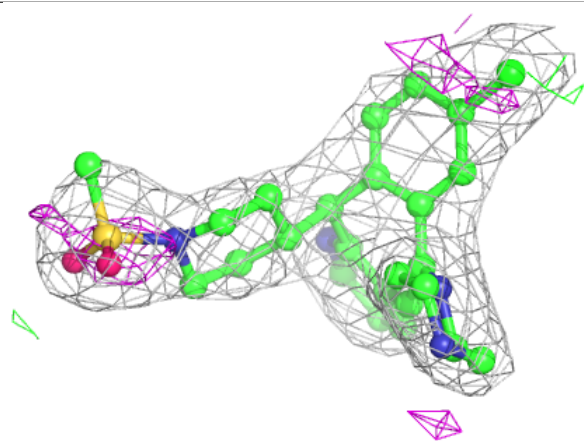
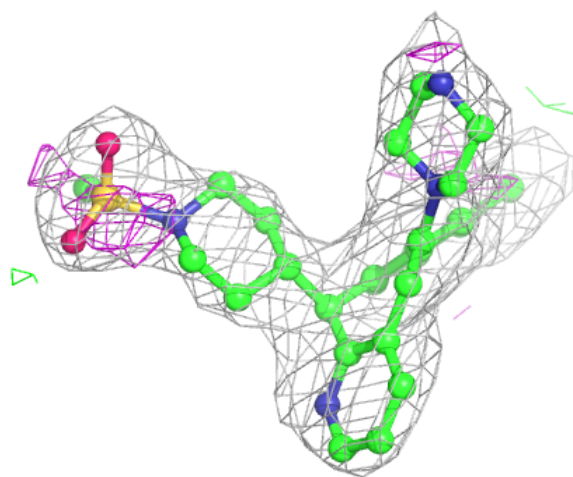
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

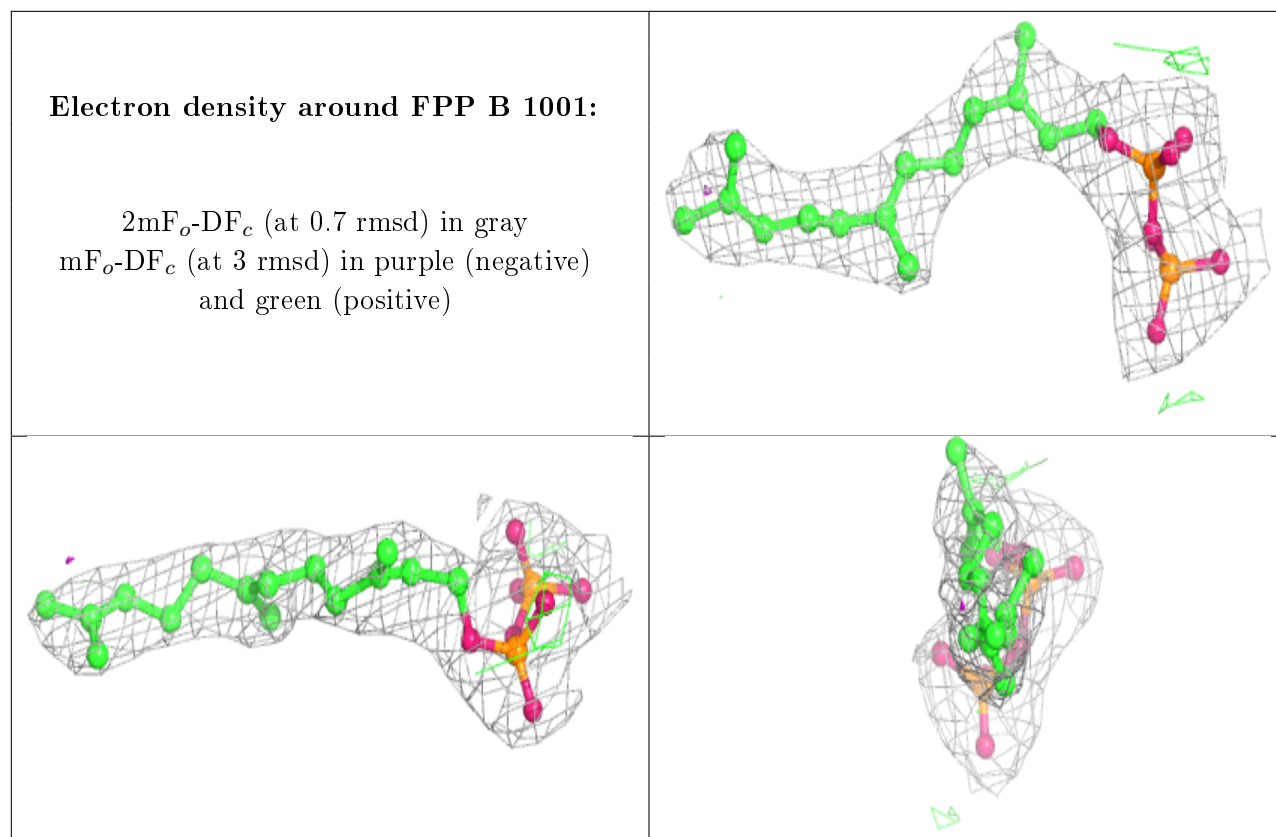
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	736	B	2001	32/32	0.95	0.18	39,52,58,59	0
3	ZN	B	1	1/1	0.98	0.06	66,66,66,66	0
4	FPP	B	1001	24/24	0.98	0.11	15,36,41,43	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 736 B 2001:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.