



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 12:07 PM BST

PDB ID : 1LBZ
Title : Crystal Structure of a complex (P32 crystal form) of dual activity FB-Pase/IMPase (AF2372) from Archaeoglobus fulgidus with 3 Calcium ions and Fructose-1,6 bisphosphate
Authors : Stieglitz, K.A.; Johnson, K.A.; Yang, H.; Roberts, M.F.; Seaton, B.A.; Head, J.F.; Stec, B.
Deposited on : 2002-04-04
Resolution : 2.20 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 : 2.13.1
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) : 2.13.1

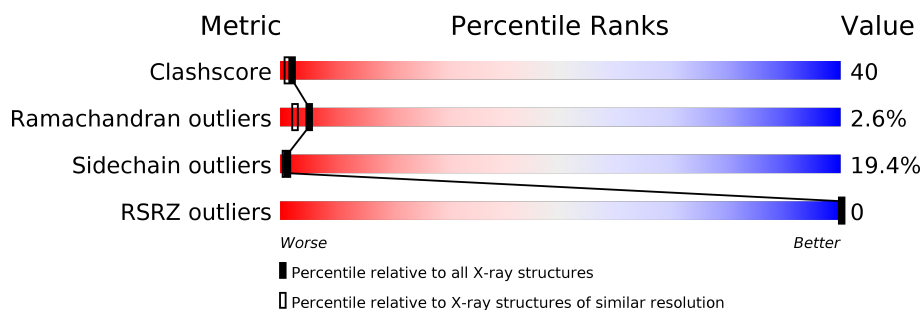
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	252	
1	B	252	

2 Entry composition [i](#)

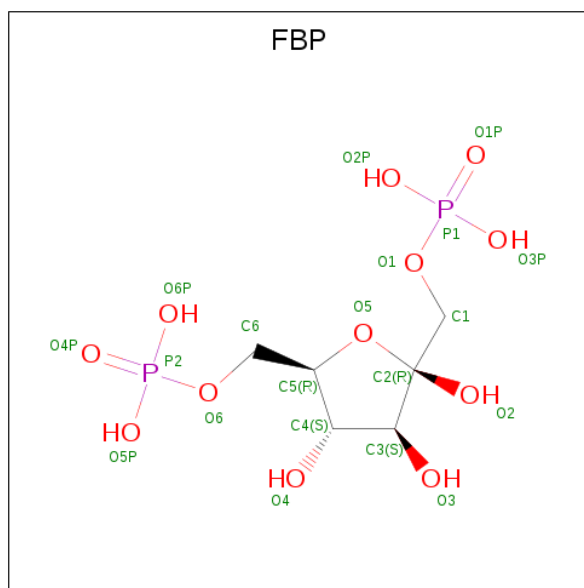
There are 4 unique types of molecules in this entry. The entry contains 4120 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 fructose 1,6-bisphosphatase/inositol monophosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	0	0
			1966	1243	333	380	10			
1	B	252	Total	C	N	O	S	0	0	0
			1966	1243	333	380	10			

- Molecule 2 is 1,6-di-O-phosphono-beta-D-fructofuranose (three-letter code: FBP) (formula: $C_6H_{14}O_{12}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			20	6	12	2		
2	B	1	Total	C	O	P	0	0
			20	6	12	2		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total 3	Ca 3	0	0
3	A	3	Total 3	Ca 3	0	0

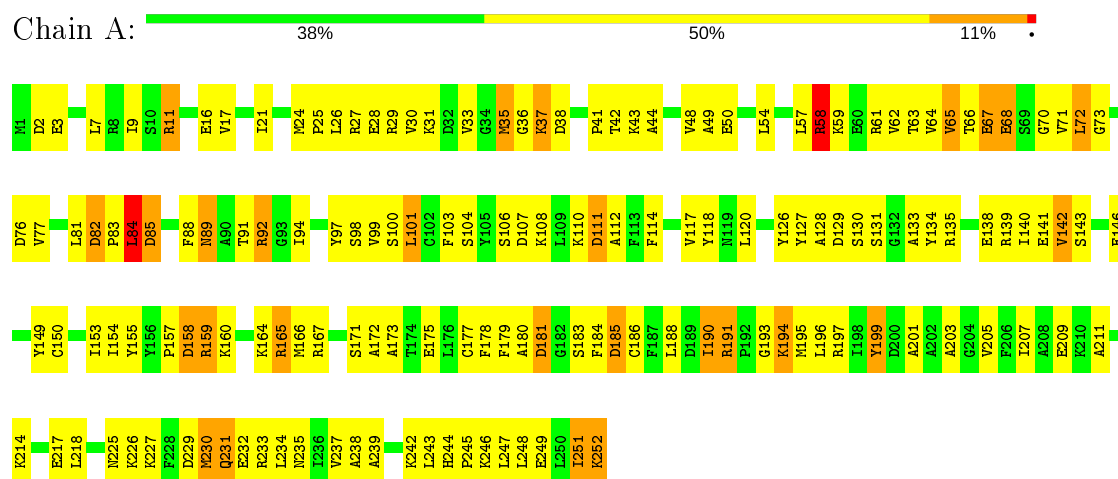
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	70	Total 70	O 70	0	0
4	B	72	Total 72	O 72	0	0

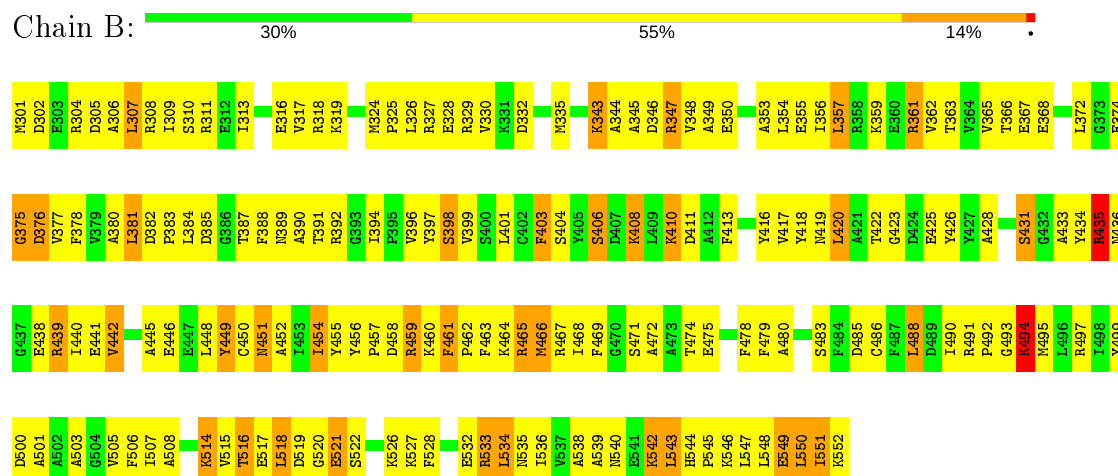
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: fructose 1,6-bisphosphatase/inositol monophosphatase



- Molecule 1: fructose 1,6-bisphosphatase/inositol monophosphatase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	89.61Å 89.61Å 102.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.20 44.80 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.20) 89.3 (44.80-2.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.20Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.213 , 0.250 0.216 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.232	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 57.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l 0.449 for h,-h-k,-l 0.000 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4120	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.82% 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: CA, FBP

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	0.31	0/1999	0.94	4/2688 (0.1%)
1	B	0.32	0/1999	0.99	4/2688 (0.1%)
All	All	0.32	0/3998	0.96	8/5376 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	435	ARG	NE-CZ-NH1	7.92	124.26	120.30
1	B	435	ARG	CD-NE-CZ	7.36	133.91	123.60
1	B	533	ARG	NE-CZ-NH1	6.78	123.69	120.30
1	A	85	ASP	CB-CG-OD2	6.63	124.27	118.30
1	B	533	ARG	CD-NE-CZ	6.39	132.55	123.60
1	A	165	ARG	NE-CZ-NH1	-5.68	117.46	120.30
1	A	185	ASP	CB-CG-OD1	5.07	122.86	118.30
1	A	82	ASP	CB-CG-OD2	5.01	122.81	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	1966	0	1949	142	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1966	0	1946	180	0
2	A	20	0	10	4	0
2	B	20	0	10	2	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
4	A	70	0	0	2	0
4	B	72	0	0	2	0
All	All	4120	0	3915	315	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 40.

All (315) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:448:LEU:HD13	1:B:550:LEU:HD22	1.47	0.94
1:A:70:GLY:HA2	1:A:230:MET:HE2	1.52	0.92
1:B:526:LYS:HE3	1:B:534:LEU:HB3	1.53	0.89
1:A:54:LEU:HB3	1:A:58:ARG:HE	1.37	0.89
1:A:62:VAL:HG12	1:A:77:VAL:HB	1.56	0.87
1:B:313:ILE:HG23	1:B:353:ALA:HA	1.61	0.83
1:A:89:ASN:HB3	1:A:94:ILE:HB	1.61	0.82
1:B:515:VAL:HG13	1:B:538:ALA:HB2	1.63	0.80
1:B:381:LEU:HB2	1:B:401:LEU:HD23	1.65	0.79
1:A:155:TYR:HB3	1:A:191:ARG:HD3	1.65	0.78
1:A:141:GLU:HG3	1:A:211:ALA:O	1.87	0.75
1:B:325:PRO:HG2	1:B:328:GLU:HG2	1.68	0.74
1:B:547:LEU:O	1:B:551:ILE:HD13	1.86	0.74
1:A:159:ARG:HA	1:B:458:ASP:O	1.87	0.73
1:B:435:ARG:HG2	1:B:436:ASN:HD22	1.52	0.73
1:B:540:ASN:OD1	1:B:542:LYS:HG3	1.88	0.73
1:A:157:PRO:HG2	1:B:457:PRO:HB2	1.72	0.72
1:A:28:GLU:HA	1:A:31:LYS:HE3	1.72	0.72
1:A:82:ASP:OD2	1:A:85:ASP:HB2	1.90	0.72
1:A:54:LEU:HB3	1:A:58:ARG:NE	2.05	0.71
1:A:190:ILE:HD12	1:A:251:ILE:HG21	1.72	0.70
1:A:85:ASP:HA	2:A:295:FBP:O1P	1.91	0.70
1:B:399:VAL:HG23	1:B:420:LEU:HD21	1.72	0.70
1:B:460:LYS:HE2	1:B:462:PRO:HD3	1.73	0.70
1:B:387:THR:O	1:B:391:THR:HG23	1.91	0.69
1:B:302:ASP:O	1:B:305:ASP:HB2	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:466:MET:O	1:B:467:ARG:HD3	1.92	0.69
1:B:330:VAL:O	1:B:343:LYS:HE2	1.92	0.69
1:A:33:VAL:HG22	1:A:42:THR:O	1.94	0.68
1:B:305:ASP:O	1:B:309:ILE:HD12	1.92	0.68
1:B:435:ARG:HG2	1:B:436:ASN:ND2	2.07	0.68
1:A:24:MET:SD	1:A:48:VAL:HG21	2.34	0.67
1:A:44:ALA:O	1:A:48:VAL:HG23	1.95	0.67
1:B:450:CYS:HB2	1:B:485:ASP:OD2	1.95	0.66
1:B:497:ARG:HD3	1:B:533:ARG:HG2	1.79	0.65
1:A:37:LYS:O	1:A:37:LYS:HD3	1.95	0.65
1:B:361:ARG:HH22	1:B:374:GLU:HB3	1.61	0.65
1:B:475:GLU:O	1:B:478:PHE:HB2	1.97	0.65
1:A:50:GLU:HG3	1:A:83:PRO:HG2	1.78	0.64
1:A:101:LEU:O	1:A:203:ALA:HB2	1.97	0.64
1:A:243:LEU:HD11	1:A:247:LEU:HD21	1.80	0.64
1:B:365:VAL:HG21	1:B:528:PHE:O	1.98	0.63
1:B:366:THR:HG22	1:B:381:LEU:HD23	1.80	0.63
1:B:516:THR:OG1	1:B:520:GLY:HA2	1.99	0.62
1:A:21:ILE:HD11	1:A:49:ALA:HB2	1.81	0.62
1:A:36:GLY:HA3	1:A:68:GLU:OE2	2.00	0.62
1:B:472:ALA:HB2	2:B:595:FBP:O4	2.00	0.61
1:A:164:LYS:O	1:A:164:LYS:HG2	1.99	0.61
1:B:514:LYS:HE3	1:B:516:THR:HG22	1.81	0.61
1:A:146:GLU:OE2	1:A:242:LYS:HE2	2.00	0.61
1:A:58:ARG:NH1	1:A:72:LEU:HD12	2.16	0.61
1:B:546:LYS:O	1:B:549:GLU:HB2	2.01	0.61
1:B:514:LYS:O	1:B:538:ALA:HA	2.01	0.61
1:A:83:PRO:HA	1:A:99:VAL:HG23	1.82	0.61
1:B:449:TYR:CZ	1:B:464:LYS:HG3	2.36	0.61
1:A:126:TYR:CE2	1:A:135:ARG:HD2	2.36	0.60
1:B:445:ALA:O	1:B:543:LEU:HD22	2.01	0.60
1:B:478:PHE:O	1:B:483:SER:HB2	2.01	0.60
1:A:108:LYS:HG2	1:A:227:LYS:HG2	1.82	0.60
1:A:85:ASP:OD2	2:A:295:FBP:O1	2.19	0.60
1:B:418:TYR:CZ	1:B:423:GLY:HA2	2.37	0.59
1:B:456:TYR:CE1	1:B:491:ARG:HD3	2.37	0.59
1:B:388:PHE:O	1:B:392:ARG:HG2	2.02	0.59
1:B:471:SER:O	1:B:475:GLU:HG3	2.03	0.59
1:A:126:TYR:CD2	1:A:177:CYS:HB3	2.37	0.59
1:A:25:PRO:HD2	1:A:28:GLU:HB3	1.84	0.59
1:B:426:TYR:OH	1:B:435:ARG:HD2	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:ALA:O	1:A:139:ARG:HA	2.02	0.58
1:B:376:ASP:HB2	4:B:1014:HOH:O	2.02	0.58
1:A:158:ASP:O	1:B:459:ARG:HA	2.04	0.58
1:B:419:ASN:HB3	1:B:422:THR:OG1	2.03	0.58
1:B:381:LEU:HB2	1:B:401:LEU:CD2	2.34	0.58
1:B:514:LYS:NZ	1:B:515:VAL:H	2.01	0.58
1:A:154:ILE:HD11	1:A:190:ILE:HD11	1.86	0.58
1:B:343:LYS:HB2	1:B:346:ASP:OD2	2.04	0.58
1:A:108:LYS:HD3	1:A:225:ASN:HA	1.85	0.57
1:A:81:LEU:HG	1:A:83:PRO:HD3	1.86	0.57
1:B:503:ALA:O	1:B:507:ILE:HD12	2.04	0.57
1:A:107:ASP:O	1:A:227:LYS:HA	2.05	0.57
1:A:28:GLU:HA	1:A:31:LYS:CE	2.33	0.57
1:A:84:LEU:HD12	1:A:120:LEU:CD1	2.34	0.57
1:A:214:LYS:HG2	1:A:244:HIS:CD2	2.40	0.57
1:B:349:ALA:HB1	1:B:383:PRO:HB2	1.87	0.57
1:A:191:ARG:HH12	2:A:295:FBP:H62	1.70	0.57
1:B:455:TYR:CZ	1:B:468:ILE:HG23	2.40	0.57
1:B:357:LEU:HD13	1:B:362:VAL:HG21	1.86	0.57
1:B:361:ARG:HH12	1:B:374:GLU:HB2	1.70	0.56
1:B:452:ALA:O	1:B:466:MET:HA	2.06	0.56
1:A:104:SER:HA	1:A:112:ALA:HA	1.87	0.56
1:A:159:ARG:HH22	1:A:251:ILE:HA	1.70	0.56
1:B:548:LEU:O	1:B:552:LYS:HB3	2.06	0.56
1:B:313:ILE:HD12	1:B:401:LEU:HD11	1.88	0.55
1:B:309:ILE:HG21	1:B:357:LEU:HD23	1.87	0.55
1:B:514:LYS:HZ2	1:B:514:LYS:HA	1.71	0.55
1:A:27:ARG:HG3	1:A:27:ARG:HH11	1.71	0.55
1:A:126:TYR:CE2	1:A:177:CYS:HB3	2.43	0.54
1:A:226:LYS:CE	1:A:234:LEU:HD22	2.38	0.54
1:A:217:GLU:OE2	1:A:235:ASN:HB2	2.08	0.54
1:A:64:VAL:O	1:A:71:VAL:HA	2.08	0.54
1:A:193:GLY:O	1:A:195:MET:HG3	2.08	0.54
1:A:196:LEU:O	1:A:233:ARG:HA	2.09	0.53
1:B:399:VAL:O	1:B:417:VAL:HA	2.09	0.53
1:A:194:LYS:O	1:A:233:ARG:HB3	2.08	0.53
1:B:328:GLU:O	1:B:344:ALA:HB3	2.08	0.53
1:B:440:ILE:HD12	1:B:480:ALA:O	2.09	0.52
1:A:7:LEU:HD11	1:A:11:ARG:CZ	2.40	0.52
1:A:183:SER:O	1:B:392:ARG:HD2	2.09	0.52
1:A:149:TYR:CZ	1:A:164:LYS:HG3	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:327:ARG:HG3	1:B:328:GLU:N	2.24	0.52
1:A:190:ILE:HD12	1:A:251:ILE:CG2	2.37	0.52
1:B:324:MET:CE	1:B:329:ARG:HA	2.38	0.52
1:B:547:LEU:HA	1:B:550:LEU:HB3	1.92	0.52
1:B:307:LEU:HD12	1:B:311:ARG:HG3	1.91	0.52
1:B:345:ALA:O	1:B:348:VAL:HG22	2.10	0.52
1:B:377:VAL:HG13	1:B:404:SER:O	2.09	0.52
1:B:494:LYS:NZ	1:B:535:ASN:HB2	2.25	0.51
1:B:517:GLU:OE2	1:B:518:LEU:HD12	2.10	0.51
1:A:244:HIS:HB3	1:A:245:PRO:HD3	1.92	0.51
1:B:324:MET:HE2	1:B:329:ARG:HA	1.91	0.51
1:A:194:LYS:HB3	1:A:233:ARG:HB3	1.92	0.51
1:B:497:ARG:HD3	1:B:533:ARG:HE	1.75	0.51
1:B:361:ARG:HH12	1:B:374:GLU:CB	2.23	0.51
1:B:514:LYS:HE3	1:B:516:THR:CG2	2.40	0.51
1:A:84:LEU:HD12	1:A:120:LEU:HD12	1.93	0.50
1:A:88:PHE:O	1:A:92:ARG:HG3	2.12	0.50
1:B:313:ILE:O	1:B:317:VAL:HG23	2.11	0.50
1:B:385:ASP:O	1:B:397:TYR:HA	2.12	0.50
1:A:166:MET:O	1:B:468:ILE:HG21	2.11	0.50
1:B:361:ARG:NH1	1:B:375:GLY:N	2.59	0.50
1:B:448:LEU:HD21	1:B:546:LYS:CB	2.41	0.50
1:A:179:PHE:CE1	1:A:238:ALA:HB1	2.46	0.50
1:B:350:GLU:OE1	1:B:368:GLU:HB2	2.11	0.50
1:B:490:ILE:O	1:B:492:PRO:HD3	2.12	0.50
1:A:38:ASP:OD1	1:A:67:GLU:O	2.30	0.50
1:B:367:GLU:HG3	1:B:499:TYR:CE1	2.46	0.50
1:B:518:LEU:HD11	1:B:535:ASN:HB2	1.93	0.50
1:A:58:ARG:HD2	1:A:72:LEU:HD12	1.94	0.50
1:A:177:CYS:HA	1:A:180:ALA:HB3	1.94	0.50
1:A:239:ALA:HB3	1:A:244:HIS:HA	1.93	0.50
1:A:157:PRO:HB3	1:A:166:MET:SD	2.52	0.49
1:B:362:VAL:CA	1:B:375:GLY:HA3	2.42	0.49
1:B:449:TYR:O	1:B:464:LYS:HE2	2.12	0.49
1:A:128:ALA:HB2	1:A:207:ILE:HA	1.95	0.49
1:A:3:GLU:HG2	1:A:129:ASP:HB2	1.93	0.49
1:A:153:ILE:HD13	1:A:175:GLU:HB3	1.93	0.49
1:A:26:LEU:O	1:A:29:ARG:HB2	2.13	0.49
1:B:325:PRO:HG2	1:B:328:GLU:CG	2.40	0.49
1:B:428:ALA:O	1:B:506:PHE:HE2	1.94	0.49
1:B:389:ASN:HB3	1:B:394:ILE:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:398:SER:HB3	1:B:418:TYR:O	2.11	0.49
1:B:469:PHE:HD1	1:B:475:GLU:OE1	1.95	0.49
1:A:100:SER:HA	1:A:117:VAL:HG22	1.95	0.49
1:B:493:GLY:HA3	1:B:495:MET:CE	2.43	0.49
1:A:229:ASP:OD1	1:A:231:GLN:HG2	2.13	0.48
1:A:88:PHE:CE2	1:A:92:ARG:HD2	2.48	0.48
1:B:330:VAL:O	1:B:330:VAL:HG22	2.13	0.48
1:B:439:ARG:HG3	1:B:440:ILE:N	2.27	0.48
1:B:501:ALA:O	1:B:505:VAL:HG23	2.13	0.48
1:A:146:GLU:HG2	1:A:242:LYS:HB2	1.95	0.48
1:B:507:ILE:HD12	1:B:507:ILE:H	1.78	0.48
1:B:514:LYS:HZ1	1:B:515:VAL:H	1.61	0.48
1:B:361:ARG:NH1	1:B:374:GLU:HB2	2.27	0.48
1:B:380:ALA:HB1	1:B:499:TYR:HA	1.95	0.48
1:B:471:SER:C	1:B:475:GLU:HG3	2.34	0.48
1:B:545:PRO:O	1:B:549:GLU:HG3	2.13	0.48
1:A:191:ARG:NH1	2:A:295:FBP:H62	2.28	0.48
1:B:399:VAL:CG2	1:B:420:LEU:HD21	2.40	0.48
1:A:28:GLU:O	1:A:31:LYS:HD2	2.14	0.47
1:A:205:VAL:O	1:A:209:GLU:HB2	2.13	0.47
1:A:217:GLU:OE2	1:A:235:ASN:O	2.32	0.47
1:B:327:ARG:HG3	1:B:328:GLU:OE2	2.14	0.47
1:A:65:VAL:HB	1:A:71:VAL:HG22	1.96	0.47
1:B:519:ASP:OD1	1:B:521:GLU:HB2	2.15	0.47
1:B:363:THR:HG22	1:B:363:THR:O	2.14	0.47
1:B:378:PHE:HB3	1:B:528:PHE:CD1	2.50	0.47
1:A:218:LEU:HB2	1:A:235:ASN:CB	2.45	0.47
1:A:146:GLU:HG2	1:A:242:LYS:CB	2.44	0.47
1:B:355:GLU:O	1:B:359:LYS:HG3	2.14	0.47
1:B:361:ARG:HH12	1:B:375:GLY:N	2.13	0.47
1:A:29:ARG:HB3	1:A:91:THR:HG22	1.96	0.47
1:B:488:LEU:HD11	1:B:547:LEU:HD22	1.97	0.46
1:A:181:ASP:OD1	1:A:183:SER:OG	2.31	0.46
1:A:3:GLU:CD	1:A:129:ASP:HB2	2.35	0.46
1:B:549:GLU:HG2	1:B:552:LYS:NZ	2.30	0.46
1:B:301:MET:HE2	1:B:413:PHE:HZ	1.81	0.46
1:B:439:ARG:HE	1:B:439:ARG:HB2	1.68	0.46
1:A:233:ARG:O	1:A:234:LEU:HD23	2.16	0.46
1:A:98:SER:HB3	1:A:118:TYR:O	2.15	0.46
1:B:361:ARG:HH22	1:B:374:GLU:CB	2.27	0.45
1:B:426:TYR:CZ	1:B:435:ARG:HD2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:456:TYR:HE1	1:B:491:ARG:HD3	1.82	0.45
1:A:67:GLU:OE2	1:A:197:ARG:NH1	2.50	0.45
1:B:357:LEU:CD1	1:B:362:VAL:HG21	2.46	0.45
1:A:242:LYS:HB3	1:A:242:LYS:HE3	1.57	0.45
1:A:3:GLU:HG2	4:A:1061:HOH:O	2.16	0.45
1:B:448:LEU:O	1:B:450:CYS:N	2.50	0.45
1:B:433:ALA:HB1	1:B:507:ILE:HG23	1.97	0.45
1:B:428:ALA:HB2	1:B:433:ALA:HA	1.97	0.45
1:B:451:ASN:OD1	1:B:465:ARG:NH1	2.50	0.45
1:A:126:TYR:CZ	1:A:135:ARG:HD2	2.52	0.45
1:B:306:ALA:HB1	1:B:403:PHE:CE1	2.52	0.45
1:A:111:ASP:OD2	1:A:111:ASP:N	2.50	0.45
1:A:166:MET:O	1:A:167:ARG:NH1	2.50	0.45
1:A:249:GLU:OE2	1:A:252:LYS:HE3	2.17	0.45
1:B:471:SER:O	1:B:474:THR:N	2.50	0.45
1:B:448:LEU:HD21	1:B:546:LYS:HB3	1.99	0.45
1:A:131:SER:CB	1:A:134:TYR:HH	2.30	0.45
1:A:157:PRO:HG2	1:B:457:PRO:CB	2.43	0.45
1:A:38:ASP:OD2	1:A:197:ARG:NH1	2.50	0.45
1:A:172:ALA:O	1:A:175:GLU:N	2.50	0.44
1:A:226:LYS:HE3	1:A:234:LEU:HD22	1.99	0.44
1:B:324:MET:HE1	1:B:344:ALA:CB	2.47	0.44
1:B:318:ARG:HA	1:B:420:LEU:HD22	1.99	0.44
1:A:197:ARG:NE	1:A:199:TYR:OH	2.50	0.44
1:A:59:LYS:NZ	4:A:1097:HOH:O	2.50	0.44
1:B:326:LEU:HD23	1:B:326:LEU:N	2.32	0.44
1:A:82:ASP:CG	1:A:85:ASP:HB2	2.37	0.44
1:B:391:THR:OG1	1:B:392:ARG:NH1	2.50	0.44
1:B:425:GLU:H	1:B:436:ASN:CG	2.21	0.44
1:B:383:PRO:O	1:B:399:VAL:HG22	2.18	0.44
1:A:243:LEU:O	1:A:246:LYS:HB2	2.16	0.44
1:A:3:GLU:OE2	1:A:130:SER:N	2.50	0.44
1:B:307:LEU:O	1:B:311:ARG:HG3	2.17	0.44
1:A:142:VAL:HG12	1:A:179:PHE:O	2.18	0.44
1:B:384:LEU:HG	1:B:384:LEU:O	2.17	0.44
1:B:527:LYS:H	1:B:532:GLU:HG2	1.83	0.44
1:A:3:GLU:CG	1:A:129:ASP:HB2	2.48	0.44
1:B:448:LEU:HD21	1:B:546:LYS:HB2	2.00	0.44
1:B:546:LYS:O	1:B:550:LEU:N	2.50	0.44
1:A:193:GLY:O	1:A:195:MET:N	2.51	0.43
1:A:126:TYR:O	1:A:127:TYR:HB3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:LEU:HD23	1:A:248:LEU:HA	1.87	0.43
1:B:356:ILE:O	1:B:359:LYS:N	2.50	0.43
1:B:381:LEU:HD13	1:B:401:LEU:HD21	2.01	0.43
1:B:347:ARG:HD3	1:B:347:ARG:HA	1.28	0.43
1:B:361:ARG:HH11	1:B:361:ARG:HG3	1.83	0.43
1:B:479:PHE:CE2	1:B:508:ALA:HB1	2.53	0.43
1:A:110:LYS:O	1:A:110:LYS:HG3	2.18	0.43
1:A:194:LYS:HB3	1:A:233:ARG:CB	2.48	0.43
1:A:27:ARG:HG3	1:A:27:ARG:NH1	2.32	0.43
1:A:72:LEU:N	1:A:72:LEU:HD23	2.33	0.43
1:B:431:SER:OG	1:B:431:SER:O	2.35	0.43
1:B:534:LEU:N	1:B:534:LEU:HD23	2.33	0.43
1:B:494:LYS:HZ2	1:B:535:ASN:H	1.67	0.43
1:A:164:LYS:O	1:A:165:ARG:HB3	2.19	0.43
1:A:184:PHE:CZ	1:B:388:PHE:HE2	2.37	0.43
1:B:350:GLU:O	1:B:354:LEU:N	2.50	0.43
1:B:406:SER:HB3	1:B:411:ASP:OD2	2.19	0.43
1:A:171:SER:O	1:A:175:GLU:HG3	2.19	0.43
1:B:324:MET:HA	1:B:325:PRO:HD2	1.71	0.43
1:B:361:ARG:HH12	1:B:374:GLU:CA	2.32	0.43
1:B:361:ARG:HH12	1:B:374:GLU:C	2.22	0.43
1:B:486:CYS:HA	1:B:538:ALA:O	2.18	0.43
1:B:544:HIS:N	1:B:545:PRO:HD2	2.34	0.43
1:A:35:MET:HE2	1:A:41:PRO:HA	2.00	0.43
1:B:408:LYS:HB2	1:B:410:LYS:HG2	2.00	0.43
1:A:17:VAL:O	1:A:21:ILE:HG12	2.19	0.42
1:B:425:GLU:N	1:B:436:ASN:OD1	2.50	0.42
2:B:595:FBP:O2	2:B:595:FBP:H61	2.19	0.42
1:B:461:PHE:HB3	1:B:463:PHE:CD1	2.54	0.42
1:B:494:LYS:HG3	1:B:535:ASN:OD1	2.19	0.42
1:A:7:LEU:O	1:A:11:ARG:HD3	2.19	0.42
1:A:31:LYS:O	1:A:43:LYS:HA	2.18	0.42
1:B:514:LYS:HD2	1:B:514:LYS:HA	1.71	0.42
1:A:63:THR:HA	1:A:73:GLY:HA3	2.01	0.42
1:B:310:SER:HB3	1:B:416:TYR:CD2	2.54	0.42
1:A:194:LYS:HG3	1:A:235:ASN:ND2	2.34	0.42
1:A:26:LEU:HD23	1:A:26:LEU:N	2.34	0.42
1:B:366:THR:CG2	1:B:381:LEU:HD23	2.45	0.42
1:B:455:TYR:CE2	1:B:468:ILE:HG23	2.55	0.42
1:A:183:SER:HB3	1:B:392:ARG:HD2	2.01	0.42
1:A:201:ALA:CB	1:A:234:LEU:HD12	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ARG:NH2	1:A:251:ILE:HA	2.34	0.42
1:A:57:LEU:O	1:A:59:LYS:N	2.52	0.42
1:B:410:LYS:HE2	1:B:410:LYS:HB2	1.50	0.42
1:B:494:LYS:NZ	1:B:518:LEU:HD11	2.35	0.42
1:A:26:LEU:HA	1:A:29:ARG:HD2	2.01	0.41
1:A:57:LEU:HA	1:A:57:LEU:HD23	1.82	0.41
1:B:396:VAL:HG12	1:B:396:VAL:O	2.20	0.41
1:A:178:PHE:O	1:A:183:SER:N	2.53	0.41
1:A:218:LEU:HB2	1:A:235:ASN:HB2	2.02	0.41
1:B:382:ASP:OD1	1:B:383:PRO:HD2	2.20	0.41
1:B:441:GLU:O	1:B:442:VAL:O	2.37	0.41
1:A:226:LYS:CD	1:A:234:LEU:HD22	2.50	0.41
1:A:217:GLU:OE2	1:A:217:GLU:HA	2.20	0.41
1:A:62:VAL:HG12	1:A:77:VAL:CB	2.37	0.41
1:B:389:ASN:O	1:B:394:ILE:N	2.49	0.41
1:B:418:TYR:CE1	1:B:423:GLY:HA2	2.56	0.41
1:B:539:ALA:HB1	1:B:543:LEU:HB3	2.02	0.41
1:A:89:ASN:N	1:A:89:ASN:OD1	2.52	0.41
1:A:150:CYS:HB2	1:A:185:ASP:OD2	2.19	0.41
1:A:188:LEU:HD23	1:A:237:VAL:HG13	2.02	0.41
1:A:30:VAL:HG22	1:A:30:VAL:O	2.21	0.41
1:B:301:MET:N	4:B:1031:HOH:O	2.53	0.41
1:B:497:ARG:HD3	1:B:533:ARG:NE	2.35	0.41
1:A:7:LEU:HA	1:A:114:PHE:CZ	2.56	0.41
1:A:84:LEU:HD11	1:A:97:TYR:CD1	2.56	0.41
1:B:350:GLU:HG2	1:B:354:LEU:HD12	2.02	0.41
1:B:497:ARG:HE	1:B:533:ARG:NH2	2.19	0.41
1:B:550:LEU:HG	1:B:551:ILE:N	2.36	0.41
1:A:167:ARG:HG3	1:A:184:PHE:CE2	2.56	0.41
1:B:390:ALA:HA	1:B:397:TYR:CE2	2.55	0.41
1:A:188:LEU:CD2	1:A:237:VAL:HG13	2.51	0.41
1:A:244:HIS:HE1	1:A:248:LEU:HD11	1.86	0.41
1:B:548:LEU:HD23	1:B:548:LEU:HA	1.90	0.41
1:B:546:LYS:HA	1:B:549:GLU:HB2	2.03	0.40
1:B:327:ARG:HD2	1:B:328:GLU:OE2	2.22	0.40
1:B:362:VAL:HA	1:B:375:GLY:HA3	2.02	0.40
1:B:454:ILE:HD13	1:B:454:ILE:HG21	1.94	0.40
1:B:526:LYS:HD2	1:B:532:GLU:CD	2.41	0.40
1:B:301:MET:HE2	1:B:413:PHE:CZ	2.56	0.40
1:B:451:ASN:HB3	1:B:465:ARG:HG3	2.03	0.40
1:B:469:PHE:HB2	1:B:475:GLU:OE1	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:426:TYR:HA	1:B:434:TYR:O	2.21	0.40
1:B:517:GLU:HB3	1:B:518:LEU:H	1.73	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	250/252 (99%)	212 (85%)	32 (13%)	6 (2%)	6	3
1	B	250/252 (99%)	217 (87%)	26 (10%)	7 (3%)	5	2
All	All	500/504 (99%)	429 (86%)	58 (12%)	13 (3%)	5	3

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	143	SER
1	B	442	VAL
1	B	549	GLU
1	B	550	LEU
1	A	58	ARG
1	A	194	LYS
1	B	522	SER
1	A	61	ARG
1	B	449	TYR
1	A	84	LEU
1	A	173	ALA
1	B	494	LYS
1	B	375	GLY

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	204/204 (100%)	168 (82%)	36 (18%)	2	1
1	B	204/204 (100%)	161 (79%)	43 (21%)	1	1
All	All	408/408 (100%)	329 (81%)	79 (19%)	1	1

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	9	ILE
1	A	11	ARG
1	A	16	GLU
1	A	35	MET
1	A	37	LYS
1	A	58	ARG
1	A	65	VAL
1	A	66	THR
1	A	67	GLU
1	A	68	GLU
1	A	72	LEU
1	A	76	ASP
1	A	84	LEU
1	A	89	ASN
1	A	92	ARG
1	A	101	LEU
1	A	103	PHE
1	A	106	SER
1	A	111	ASP
1	A	138	GLU
1	A	140	ILE
1	A	142	VAL
1	A	158	ASP
1	A	159	ARG
1	A	160	LYS
1	A	181	ASP

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Mol	Chain	Res	Type
1	A	186	CYS
1	A	190	ILE
1	A	191	ARG
1	A	199	TYR
1	A	230	MET
1	A	231	GLN
1	A	232	GLU
1	A	251	ILE
1	A	252	LYS
1	B	304	ARG
1	B	307	LEU
1	B	308	ARG
1	B	316	GLU
1	B	319	LYS
1	B	332	ASP
1	B	335	MET
1	B	343	LYS
1	B	347	ARG
1	B	357	LEU
1	B	361	ARG
1	B	372	LEU
1	B	376	ASP
1	B	381	LEU
1	B	398	SER
1	B	403	PHE
1	B	406	SER
1	B	408	LYS
1	B	410	LYS
1	B	420	LEU
1	B	431	SER
1	B	435	ARG
1	B	438	GLU
1	B	439	ARG
1	B	446	GLU
1	B	451	ASN
1	B	454	ILE
1	B	459	ARG
1	B	461	PHE
1	B	465	ARG
1	B	466	MET
1	B	488	LEU
1	B	494	LYS

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Mol	Chain	Res	Type
1	B	500	ASP
1	B	514	LYS
1	B	516	THR
1	B	518	LEU
1	B	521	GLU
1	B	534	LEU
1	B	536	ILE
1	B	542	LYS
1	B	543	LEU
1	B	551	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	235	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FBP	A	295	3	18,20,20	1.38	3 (16%)	23,32,32	1.77	6 (26%)
2	FBP	B	595	3	18,20,20	1.37	3 (16%)	23,32,32	1.91	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	FBP	A	295	3	-	6/13/32/32	0/1/1/1
2	FBP	B	595	3	-	7/13/32/32	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	295	FBP	O5-C2	-3.36	1.38	1.43
2	B	595	FBP	O5-C2	-3.06	1.38	1.43
2	B	595	FBP	P1-O1	-2.71	1.51	1.60
2	B	595	FBP	P1-O3P	-2.68	1.44	1.54
2	A	295	FBP	P1-O3P	-2.66	1.44	1.54
2	A	295	FBP	P1-O1	-2.40	1.52	1.60

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	295	FBP	O4-C4-C5	4.56	124.24	111.05
2	B	595	FBP	O3P-P1-O1	4.26	118.08	106.73
2	B	595	FBP	O4-C4-C5	4.07	122.82	111.05
2	A	295	FBP	O4-C4-C3	-3.47	101.77	112.15
2	B	595	FBP	O4-C4-C3	-3.35	102.11	112.15
2	A	295	FBP	O3P-P1-O1	2.99	114.69	106.73
2	B	595	FBP	O6P-P2-O6	2.80	114.19	106.73
2	B	595	FBP	P1-O1-C1	2.77	125.93	118.30
2	B	595	FBP	O2-C2-O5	-2.52	104.64	109.50
2	A	295	FBP	O6P-P2-O6	2.50	113.39	106.73
2	A	295	FBP	O2-C2-O5	-2.02	105.60	109.50
2	A	295	FBP	P2-O6-C6	2.02	123.85	118.30

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	595	FBP	C1-O1-P1-O1P
2	B	595	FBP	C1-O1-P1-O2P
2	B	595	FBP	C1-O1-P1-O3P
2	B	595	FBP	C6-O6-P2-O4P
2	B	595	FBP	C6-O6-P2-O5P
2	B	595	FBP	C6-O6-P2-O6P
2	A	295	FBP	C4-C5-C6-O6
2	A	295	FBP	C6-O6-P2-O6P
2	A	295	FBP	O5-C5-C6-O6
2	A	295	FBP	C6-O6-P2-O4P
2	A	295	FBP	C5-C6-O6-P2
2	B	595	FBP	C5-C6-O6-P2
2	A	295	FBP	C6-O6-P2-O5P

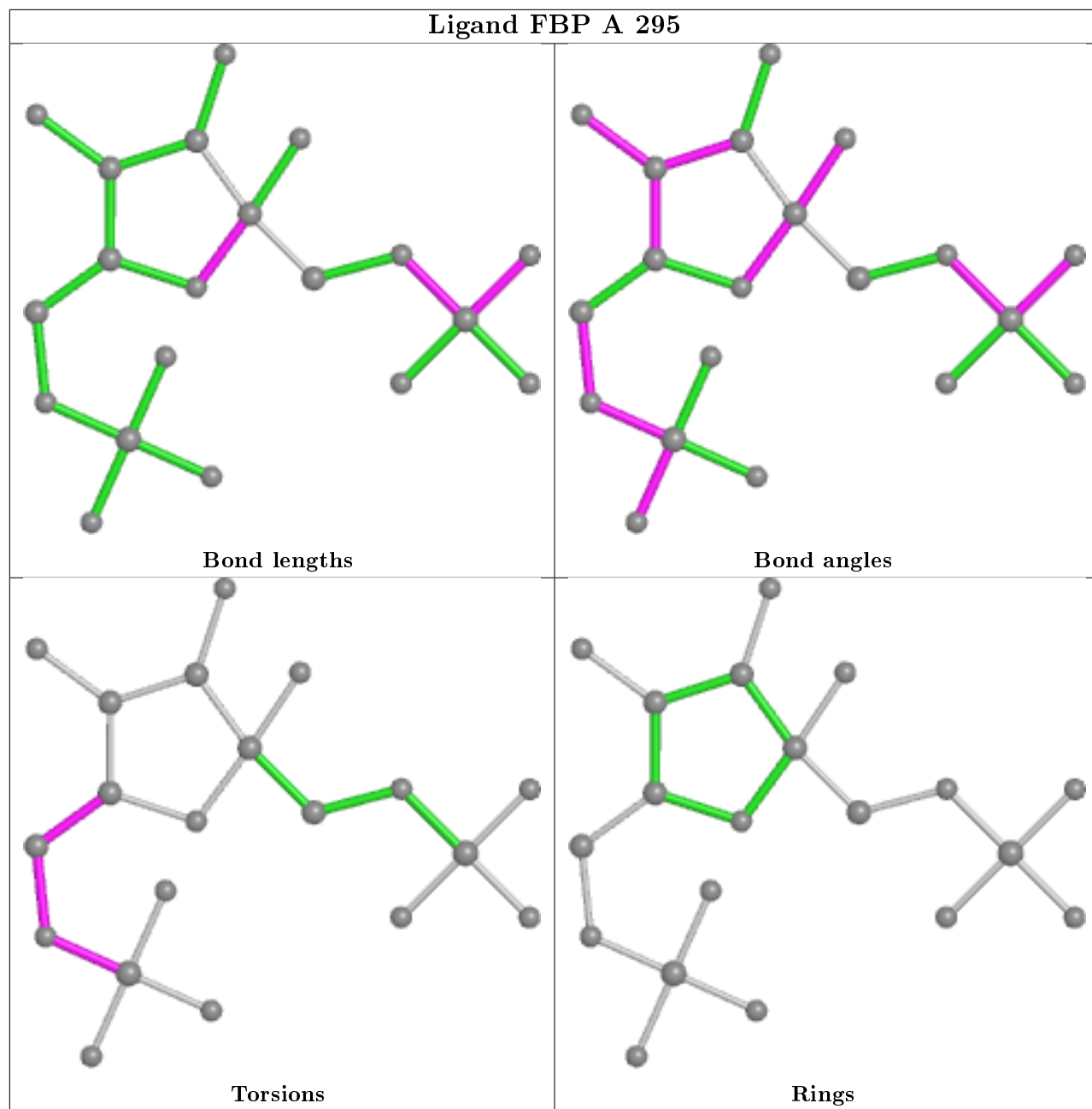
There are no ring outliers.

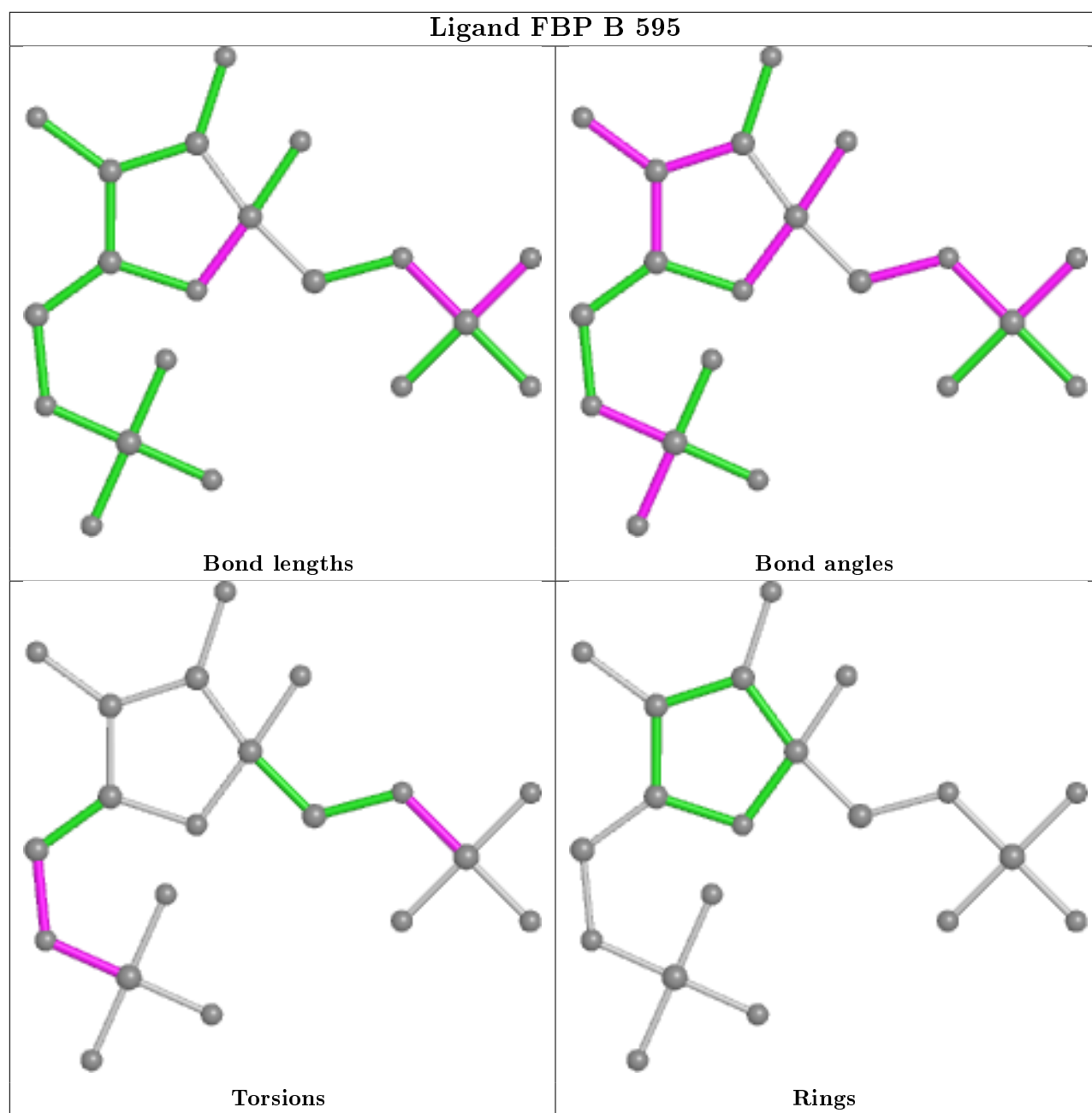
2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	295	FBP	4	0
2	B	595	FBP	2	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.

Ligand FBP A 295





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	252/252 (100%)	-0.54	0 100 100	20, 38, 63, 83	0
1	B	252/252 (100%)	-0.55	0 100 100	16, 37, 63, 84	0
All	All	504/504 (100%)	-0.54	0 100 100	16, 37, 63, 84	0

There are no RSRZ outliers to report.

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 monosaccharides 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(Å ²)	Q<0.9
2	FBP	B	595	20/20	0.95	0.13	12,58,97,98	0
3	CA	B	591	1/1	0.98	0.04	11,11,11,11	0
2	FBP	A	295	20/20	0.98	0.11	18,32,54,76	0
3	CA	B	590	1/1	0.99	0.07	10,10,10,10	0
3	CA	B	592	1/1	1.00	0.03	13,13,13,13	0
3	CA	A	290	1/1	1.00	0.02	15,15,15,15	0
3	CA	A	291	1/1	1.00	0.03	10,10,10,10	0

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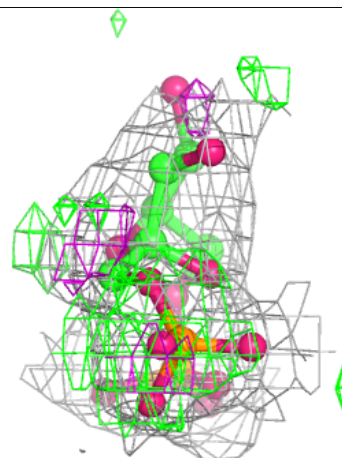
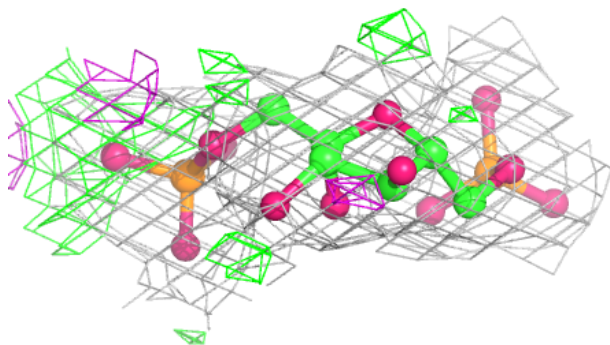
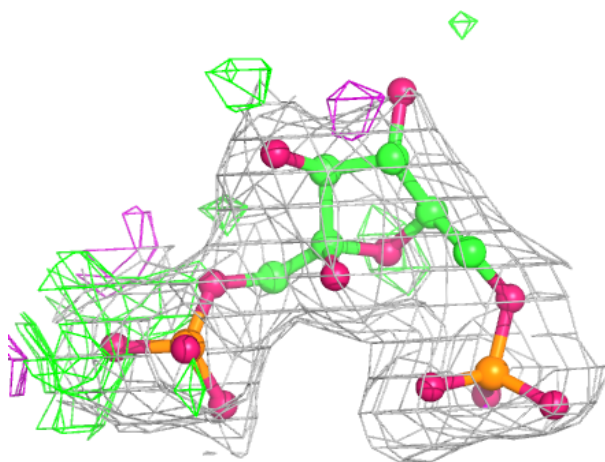
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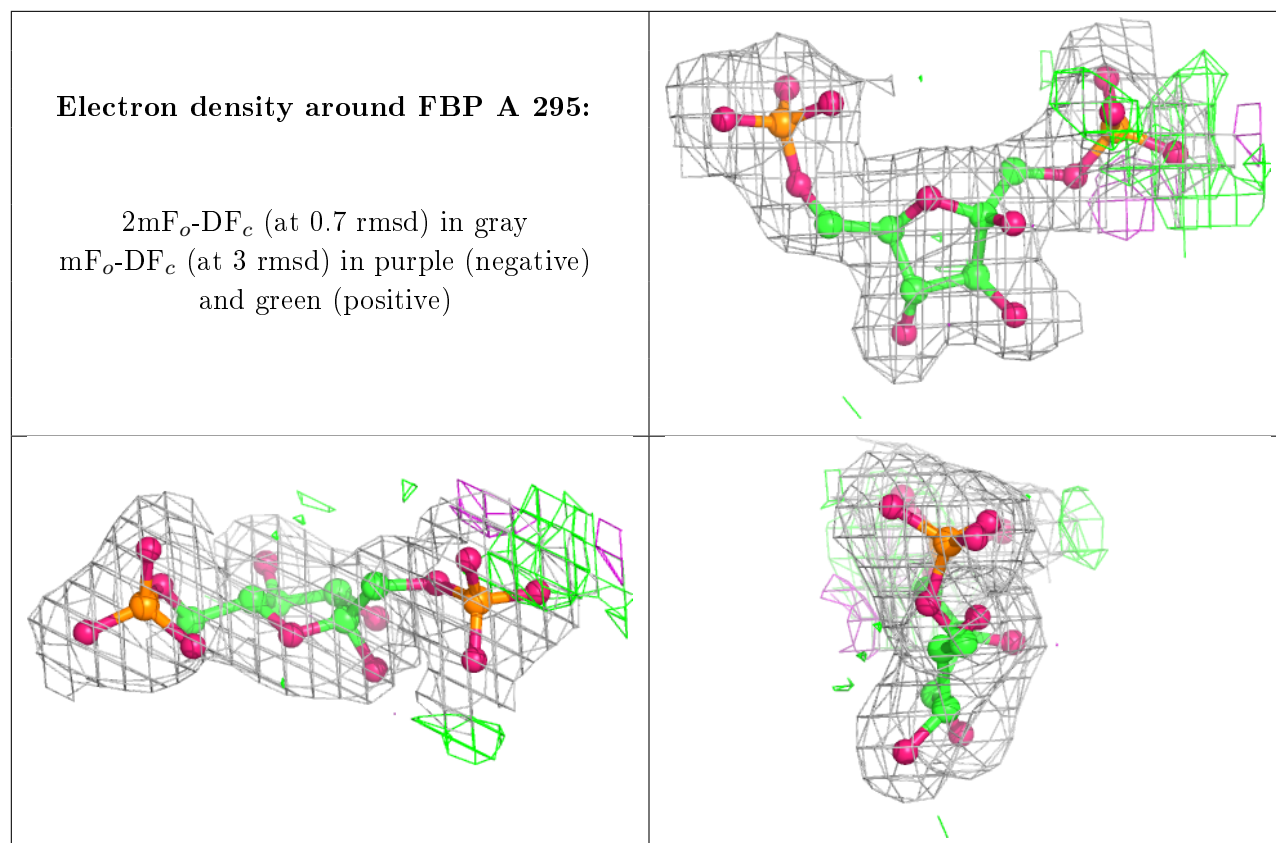
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	CA	A	292	1/1	1.00	0.03	10,10,10,10	0

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 FBP B 595:

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.