



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:54 PM UTC

PDB ID : 1UWJ / pdb_00001uwj
Title : The complex of mutant V599E B-RAF and BAY439006.
Authors : Barford, D.; Roe, S.M.; Wan, P.T.C.; Cancer Genome Project
Deposited on : 2004-02-05
Resolution : 3.50 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

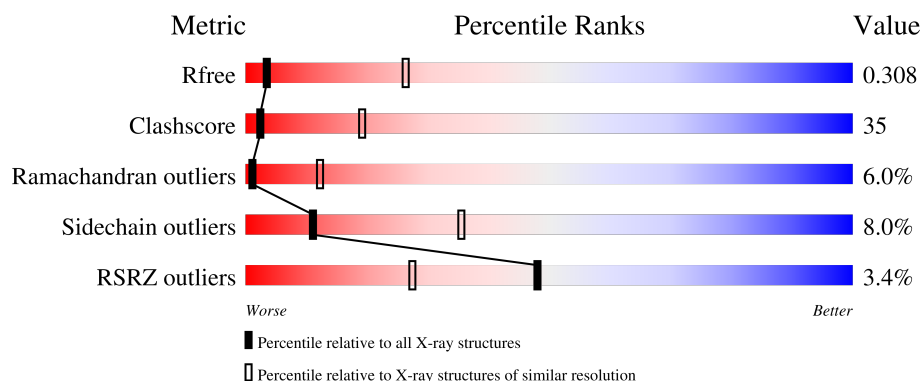
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

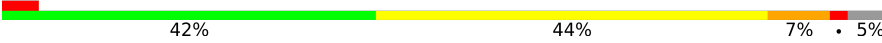
The reported resolution of this entry is 3.50 Å.

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}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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 of 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	276	
1	B	276	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4281 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.

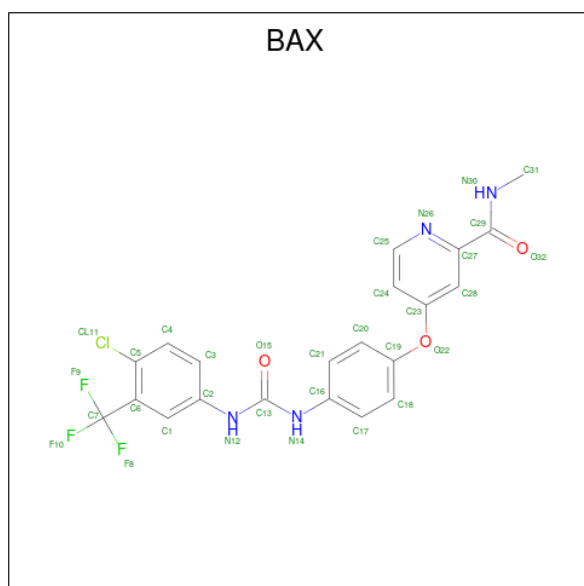
- Molecule 1 is a protein called B-RAF PROTO-ONCOGENE SERINE/THREONINE-PROTEIN KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2111	1352	367	379	13			
1	B	263	Total	C	N	O	S	0	0	0
			2105	1349	366	377	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	599	GLU	VAL	engineered mutation	UNP P15056
B	599	GLU	VAL	engineered mutation	UNP P15056

- Molecule 2 is 4-{4-[(4-CHLORO-3-(TRIFLUOROMETHYL)PHENYL)AMINO]CARBOXYL)AMINO]PHENOXY}-N-METHYLPYRIDINE-2-CARBOXAMIDE (CCD ID: BAX) (formula: $C_{21}H_{16}ClF_3N_4O_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Cl	F	N	O	0	0
			32	21	1	3	4	3		
2	B	1	Total	C	Cl	F	N	O	0	0
			32	21	1	3	4	3		

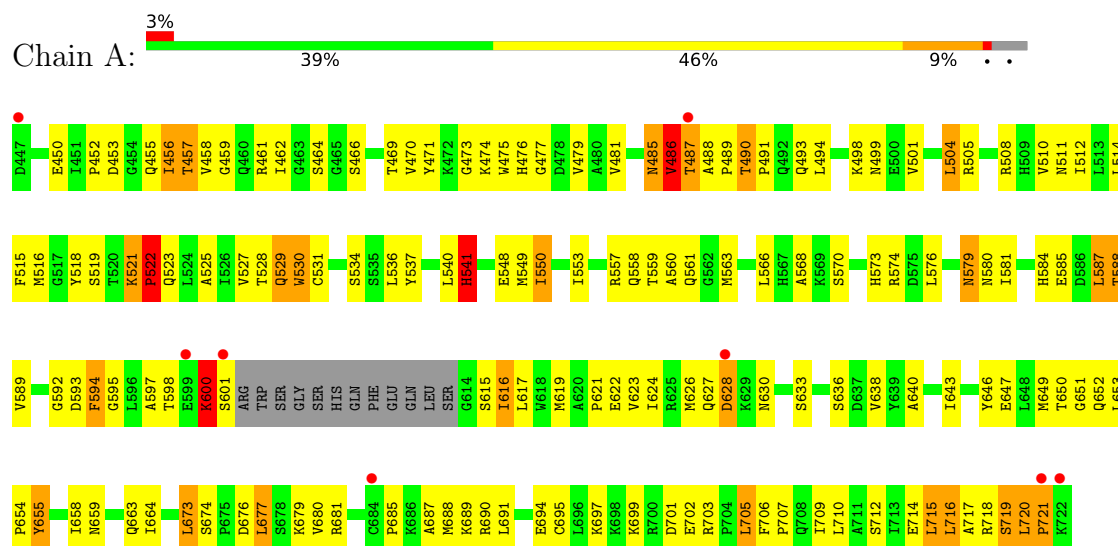
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		

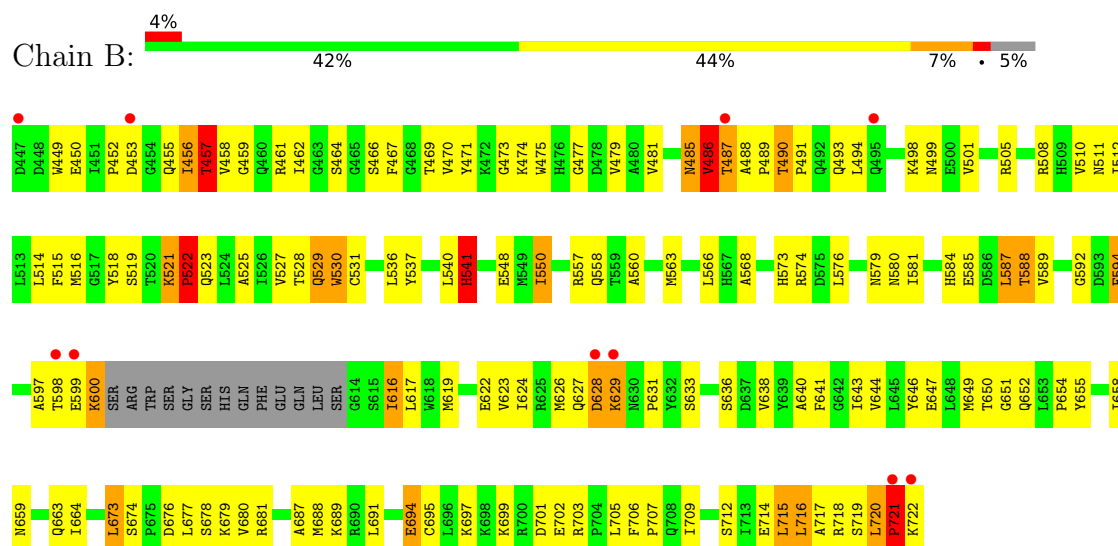
3 Residue-property plots [i](#)

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: B-RAF PROTO-ONCOGENE SERINE/THREONINE-PROTEIN KINASE



• Molecule 1: B-RAF PROTO-ONCOGENE SERINE/THREONINE-PROTEIN KINASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	97.66Å 97.66Å 168.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.39 – 3.50 29.39 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.39-3.50) 99.6 (29.39-3.50)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.68 (at 3.47Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.275 , 0.357 0.246 , 0.308	Depositor DCC
R_{free} test set	566 reflections (5.25%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.615	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 15.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	4281	wwPDB-VP
Average B, all atoms (Å ²)	5.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.21% 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: BAX

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	0.63	0/2156	1.03	8/2907 (0.3%)
1	B	0.61	0/2150	1.00	5/2899 (0.2%)
All	All	0.62	0/4306	1.01	13/5806 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	630	ASN	CA-C-N	10.13	129.92	119.19
1	A	630	ASN	C-N-CA	10.13	129.92	119.19
1	A	548	GLU	N-CA-C	-7.39	100.75	110.43
1	B	548	GLU	N-CA-C	-7.20	101.00	110.43
1	A	541	HIS	N-CA-C	6.60	121.49	113.17
1	B	541	HIS	N-CA-C	6.35	121.17	113.17
1	A	485	ASN	N-CA-C	6.06	117.71	110.19
1	B	485	ASN	N-CA-C	5.94	117.55	110.19
1	A	530	TRP	N-CA-C	-5.59	102.20	110.48
1	B	616	ILE	N-CA-C	5.52	120.82	109.34
1	A	719	SER	N-CA-C	5.29	117.17	109.71
1	B	530	TRP	N-CA-C	-5.16	102.11	110.32
1	A	705	LEU	N-CA-C	-5.06	103.91	110.19

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	2111	0	2150	155	0
1	B	2105	0	2145	155	0
2	A	32	0	16	0	0
2	B	32	0	16	0	0
3	A	1	0	0	0	0
All	All	4281	0	4327	303	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 35.

All (303) 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:563:MET:HE3	1:A:576:LEU:HD22	1.40	1.04
1:B:563:MET:HE3	1:B:576:LEU:HD22	1.42	0.99
1:B:626:MET:SD	1:B:631:PRO:HG2	2.10	0.91
1:A:550:ILE:HD12	1:A:550:ILE:H	1.38	0.89
1:B:550:ILE:HD12	1:B:550:ILE:H	1.37	0.89
1:B:598:THR:HG22	1:B:599:GLU:H	1.39	0.88
1:B:521:LYS:HB3	1:B:522:PRO:HD2	1.59	0.85
1:B:464:SER:HB3	1:B:469:THR:HG23	1.59	0.84
1:A:521:LYS:HB3	1:A:522:PRO:HD2	1.59	0.82
1:B:490:THR:HG22	1:B:493:GLN:H	1.46	0.81
1:A:479:VAL:HG12	1:A:516:MET:HE2	1.63	0.81
1:A:490:THR:HG22	1:A:493:GLN:H	1.48	0.79
1:A:464:SER:HB3	1:A:469:THR:HG23	1.65	0.78
1:B:479:VAL:HG12	1:B:516:MET:HE2	1.65	0.77
1:A:453:ASP:HB2	1:A:521:LYS:HB3	1.68	0.75
1:A:490:THR:HG23	1:A:491:PRO:HD2	1.69	0.74
1:B:453:ASP:HB2	1:B:521:LYS:HB3	1.69	0.74
1:A:452:PRO:HB2	1:A:455:GLN:HE21	1.53	0.72
1:A:550:ILE:H	1:A:550:ILE:CD1	2.03	0.72
1:A:563:MET:CE	1:A:576:LEU:HD22	2.16	0.72
1:B:563:MET:CE	1:B:576:LEU:HD22	2.20	0.72
1:B:550:ILE:H	1:B:550:ILE:CD1	2.01	0.71
1:B:680:VAL:HG21	1:B:689:LYS:HD2	1.71	0.71
1:A:680:VAL:HG21	1:A:689:LYS:HD2	1.72	0.71
1:B:719:SER:C	1:B:720:LEU:HD22	2.15	0.71
1:A:600:LYS:HG3	1:A:601:SER:H	1.56	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:452:PRO:HB2	1:B:455:GLN:HE21	1.57	0.70
1:B:687:ALA:HB1	1:B:716:LEU:HD22	1.74	0.70
1:B:453:ASP:HB2	1:B:521:LYS:CB	2.22	0.69
1:B:490:THR:HG23	1:B:491:PRO:HD2	1.73	0.69
1:A:479:VAL:CG1	1:A:516:MET:HE2	2.23	0.69
1:B:550:ILE:HD12	1:B:550:ILE:N	2.07	0.68
1:A:456:ILE:HG22	1:A:475:TRP:HB2	1.75	0.68
1:B:616:ILE:HD12	1:B:617:LEU:N	2.08	0.68
1:B:453:ASP:H	1:B:521:LYS:HE3	1.59	0.68
1:B:646:TYR:CE1	1:B:650:THR:HG21	2.29	0.68
1:A:453:ASP:HB2	1:A:521:LYS:CB	2.23	0.67
1:B:456:ILE:HG22	1:B:475:TRP:HB2	1.76	0.67
1:A:616:ILE:HD13	1:A:616:ILE:H	1.60	0.67
1:A:521:LYS:CB	1:A:522:PRO:HD2	2.24	0.67
1:A:687:ALA:HB1	1:A:716:LEU:HD22	1.76	0.67
1:B:479:VAL:CG1	1:B:516:MET:HE2	2.25	0.67
1:A:481:VAL:HG22	1:A:527:VAL:HG22	1.75	0.66
1:A:453:ASP:CG	1:A:522:PRO:HD2	2.20	0.66
1:B:521:LYS:CB	1:B:522:PRO:HD2	2.23	0.66
1:A:646:TYR:CE1	1:A:650:THR:HG21	2.30	0.66
1:A:550:ILE:HD12	1:A:550:ILE:N	2.08	0.66
1:A:452:PRO:HD2	1:A:455:GLN:NE2	2.11	0.66
1:B:453:ASP:CG	1:B:522:PRO:HD2	2.21	0.66
1:B:585:GLU:O	1:B:587:LEU:HD23	1.97	0.65
1:A:453:ASP:H	1:A:521:LYS:HE3	1.62	0.65
1:B:485:ASN:ND2	1:B:489:PRO:HG3	2.12	0.65
1:A:453:ASP:OD2	1:A:522:PRO:HG2	1.97	0.64
1:A:580:ASN:HB3	1:A:592:GLY:O	1.96	0.64
1:B:452:PRO:HD2	1:B:455:GLN:NE2	2.12	0.64
1:A:490:THR:HG23	1:A:491:PRO:CD	2.27	0.64
1:A:720:LEU:H	1:A:721:PRO:CD	2.12	0.63
1:A:485:ASN:ND2	1:A:489:PRO:HG3	2.14	0.63
1:B:481:VAL:HG22	1:B:527:VAL:HG22	1.81	0.63
1:B:599:GLU:C	1:B:600:LYS:HD3	2.23	0.63
1:B:717:ALA:HA	1:B:720:LEU:HD23	1.79	0.63
1:B:580:ASN:HB3	1:B:592:GLY:O	1.98	0.62
1:B:490:THR:HG23	1:B:491:PRO:CD	2.30	0.62
1:B:453:ASP:OD2	1:B:522:PRO:HG2	2.00	0.62
1:B:514:LEU:HD12	1:B:515:PHE:H	1.65	0.61
1:A:466:SER:HB2	1:A:597:ALA:HA	1.82	0.61
1:A:587:LEU:O	1:A:588:THR:CB	2.48	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:616:ILE:HG22	1:B:619:MET:CE	2.30	0.61
1:A:673:LEU:HD23	1:A:674:SER:H	1.66	0.60
1:A:537:TYR:HB3	1:A:579:ASN:ND2	2.17	0.60
1:B:466:SER:HB2	1:B:597:ALA:HA	1.82	0.60
1:A:453:ASP:OD1	1:A:521:LYS:HB2	2.01	0.60
1:B:587:LEU:O	1:B:588:THR:CB	2.50	0.60
1:B:537:TYR:HB3	1:B:579:ASN:ND2	2.18	0.59
1:B:453:ASP:OD1	1:B:521:LYS:HB2	2.03	0.59
1:A:581:ILE:HG23	1:A:589:VAL:CG1	2.33	0.59
1:B:646:TYR:O	1:B:650:THR:HG23	2.03	0.59
1:A:585:GLU:HG2	1:B:587:LEU:HD11	1.84	0.58
1:A:646:TYR:O	1:A:650:THR:HG23	2.04	0.58
1:B:587:LEU:O	1:B:587:LEU:HG	2.03	0.58
1:B:494:LEU:O	1:B:498:LYS:HG3	2.03	0.58
1:A:587:LEU:O	1:A:587:LEU:HG	2.03	0.57
1:A:658:ILE:HD12	1:A:673:LEU:HD11	1.86	0.57
1:B:461:ARG:HE	1:B:471:TYR:HE2	1.52	0.57
1:A:633:SER:O	1:A:636:SER:HB3	2.05	0.57
1:A:716:LEU:HD12	1:A:716:LEU:O	2.05	0.57
1:A:522:PRO:O	1:A:523:GLN:HB2	2.05	0.57
1:B:581:ILE:HG23	1:B:589:VAL:CG1	2.35	0.57
1:A:514:LEU:HD12	1:A:515:PHE:H	1.69	0.56
1:B:514:LEU:HB3	1:B:529:GLN:HG2	1.87	0.56
1:B:633:SER:O	1:B:636:SER:HB3	2.04	0.56
1:B:673:LEU:HD23	1:B:674:SER:H	1.70	0.56
1:A:514:LEU:HB3	1:A:529:GLN:HG2	1.86	0.56
1:A:581:ILE:HG23	1:A:589:VAL:HG13	1.88	0.56
1:A:658:ILE:CD1	1:A:673:LEU:HD11	2.35	0.56
1:B:716:LEU:O	1:B:716:LEU:HD12	2.06	0.55
1:A:587:LEU:O	1:A:588:THR:HB	2.05	0.55
1:A:622:GLU:HG2	1:A:623:VAL:N	2.20	0.55
1:B:490:THR:HB	1:B:493:GLN:CD	2.31	0.55
1:B:587:LEU:O	1:B:588:THR:HB	2.06	0.55
1:B:658:ILE:HD12	1:B:673:LEU:HD11	1.88	0.55
1:A:585:GLU:O	1:A:587:LEU:HD23	2.06	0.55
1:B:659:ASN:OD1	1:B:659:ASN:O	2.25	0.55
1:B:522:PRO:O	1:B:523:GLN:HB2	2.05	0.55
1:B:658:ILE:CD1	1:B:673:LEU:HD11	2.37	0.55
1:B:581:ILE:HG23	1:B:589:VAL:HG13	1.89	0.54
1:A:494:LEU:O	1:A:498:LYS:HG3	2.08	0.54
1:B:714:GLU:O	1:B:717:ALA:HB3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:THR:HB	1:A:493:GLN:CD	2.33	0.53
1:B:691:LEU:HD12	1:B:691:LEU:O	2.08	0.53
1:A:691:LEU:HD12	1:A:691:LEU:O	2.09	0.53
1:A:695:CYS:O	1:A:703:ARG:HD3	2.09	0.53
1:B:622:GLU:HG2	1:B:623:VAL:N	2.23	0.53
1:A:600:LYS:CG	1:A:601:SER:H	2.22	0.52
1:B:486:VAL:O	1:B:488:ALA:N	2.43	0.52
1:A:461:ARG:HE	1:A:471:TYR:HE2	1.56	0.52
1:A:553:ILE:HD11	1:A:685:PRO:HG2	1.92	0.52
1:B:490:THR:HB	1:B:493:GLN:NE2	2.25	0.52
1:B:633:SER:H	1:B:636:SER:HB3	1.74	0.52
1:B:699:LYS:HD2	1:B:702:GLU:OE2	2.09	0.52
1:A:519:SER:HB3	1:A:525:ALA:HB3	1.92	0.52
1:A:587:LEU:HD11	1:B:585:GLU:HG2	1.92	0.51
1:A:462:ILE:HB	1:A:470:VAL:HG12	1.93	0.51
1:B:453:ASP:HB2	1:B:521:LYS:CG	2.41	0.51
1:B:514:LEU:HD12	1:B:515:PHE:N	2.25	0.51
1:A:699:LYS:HD2	1:A:702:GLU:OE2	2.11	0.50
1:A:715:LEU:HG	1:A:716:LEU:N	2.26	0.50
1:B:563:MET:HE3	1:B:576:LEU:CD2	2.28	0.50
1:B:654:PRO:HB2	1:B:655:TYR:CD2	2.45	0.50
1:B:462:ILE:HB	1:B:470:VAL:HG12	1.93	0.50
1:B:462:ILE:HD11	1:B:530:TRP:HH2	1.76	0.50
1:A:462:ILE:CD1	1:A:530:TRP:HH2	2.25	0.50
1:B:521:LYS:O	1:B:522:PRO:C	2.55	0.50
1:B:519:SER:HB3	1:B:525:ALA:HB3	1.93	0.50
1:B:695:CYS:O	1:B:703:ARG:HD3	2.12	0.50
1:B:715:LEU:C	1:B:717:ALA:H	2.20	0.50
1:A:486:VAL:O	1:A:487:THR:C	2.54	0.50
1:A:521:LYS:O	1:A:523:GLN:O	2.30	0.50
1:A:557:ARG:O	1:A:560:ALA:HB3	2.12	0.50
1:A:452:PRO:HA	1:A:521:LYS:HE3	1.92	0.50
1:A:633:SER:H	1:A:636:SER:HB3	1.77	0.49
1:A:462:ILE:HD11	1:A:530:TRP:HH2	1.77	0.49
1:A:619:MET:HB3	1:A:624:ILE:HD11	1.94	0.49
1:B:462:ILE:CD1	1:B:530:TRP:HH2	2.25	0.49
1:B:655:TYR:CE1	1:B:673:LEU:HD22	2.47	0.49
1:A:490:THR:HB	1:A:493:GLN:NE2	2.27	0.49
1:B:536:LEU:HG	1:B:540:LEU:HD13	1.95	0.49
1:A:511:ASN:ND2	1:A:558:GLN:HB3	2.27	0.49
1:A:616:ILE:H	1:A:616:ILE:CD1	2.24	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:655:TYR:CE1	1:A:673:LEU:HD22	2.48	0.49
1:A:453:ASP:HB2	1:A:521:LYS:CG	2.42	0.49
1:B:511:ASN:ND2	1:B:558:GLN:HB3	2.27	0.49
1:A:715:LEU:HD13	1:A:718:ARG:HH21	1.77	0.49
1:B:521:LYS:O	1:B:523:GLN:O	2.30	0.49
1:A:486:VAL:O	1:A:488:ALA:N	2.45	0.49
1:B:719:SER:O	1:B:720:LEU:HD13	2.12	0.49
1:A:659:ASN:OD1	1:A:659:ASN:O	2.31	0.49
1:B:452:PRO:HA	1:B:521:LYS:HE3	1.95	0.48
1:B:486:VAL:O	1:B:487:THR:C	2.55	0.48
1:B:541:HIS:ND1	1:B:541:HIS:N	2.60	0.48
1:B:616:ILE:HD12	1:B:616:ILE:C	2.38	0.48
1:B:640:ALA:O	1:B:643:ILE:N	2.46	0.48
1:B:717:ALA:CA	1:B:720:LEU:HD23	2.42	0.48
1:B:499:ASN:C	1:B:501:VAL:N	2.71	0.48
1:B:499:ASN:C	1:B:501:VAL:H	2.20	0.48
1:B:627:GLN:O	1:B:628:ASP:HB2	2.14	0.48
1:A:563:MET:HE3	1:A:576:LEU:CD2	2.27	0.48
1:A:573:HIS:O	1:A:574:ARG:HB2	2.12	0.48
1:A:714:GLU:O	1:A:717:ALA:HB3	2.14	0.48
1:A:508:ARG:NH2	1:B:505:ARG:HA	2.28	0.48
1:A:521:LYS:O	1:A:522:PRO:C	2.56	0.48
1:A:537:TYR:CB	1:A:579:ASN:ND2	2.76	0.48
1:B:654:PRO:HB2	1:B:655:TYR:CE2	2.49	0.48
1:A:499:ASN:C	1:A:501:VAL:H	2.21	0.48
1:A:499:ASN:C	1:A:501:VAL:N	2.71	0.48
1:B:619:MET:HB3	1:B:624:ILE:HD11	1.95	0.48
1:B:697:LYS:HB2	1:B:703:ARG:HG2	1.96	0.48
1:A:541:HIS:ND1	1:A:541:HIS:N	2.61	0.47
1:A:536:LEU:HG	1:A:540:LEU:HD13	1.96	0.47
1:B:644:VAL:O	1:B:647:GLU:HB2	2.14	0.47
1:A:490:THR:HB	1:A:493:GLN:CG	2.44	0.47
1:A:593:ASP:CG	1:A:600:LYS:HZ3	2.22	0.47
1:B:461:ARG:HA	1:B:471:TYR:HD2	1.79	0.47
1:B:557:ARG:O	1:B:560:ALA:HB3	2.15	0.47
1:A:508:ARG:HB2	1:B:449:TRP:CH2	2.49	0.47
1:A:697:LYS:O	1:A:703:ARG:NH1	2.48	0.47
1:B:521:LYS:HB3	1:B:522:PRO:CD	2.37	0.47
1:A:705:LEU:HB2	1:A:707:PRO:HD2	1.97	0.47
1:B:537:TYR:CB	1:B:579:ASN:ND2	2.78	0.47
1:B:715:LEU:HD13	1:B:718:ARG:HH21	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:450:GLU:HG3	1:B:518:TYR:CE1	2.51	0.46
1:B:705:LEU:HB2	1:B:707:PRO:HD2	1.97	0.46
1:A:715:LEU:C	1:A:717:ALA:H	2.23	0.46
1:A:720:LEU:N	1:A:721:PRO:CD	2.76	0.46
1:A:505:ARG:HA	1:B:508:ARG:NH2	2.31	0.46
1:A:521:LYS:HB3	1:A:522:PRO:CD	2.39	0.46
1:A:501:VAL:O	1:A:505:ARG:HG2	2.16	0.46
1:B:467:PHE:CD1	1:B:598:THR:O	2.69	0.46
1:A:654:PRO:HB2	1:A:655:TYR:CD2	2.50	0.46
1:B:720:LEU:O	1:B:722:LYS:N	2.49	0.46
1:A:461:ARG:HA	1:A:471:TYR:HD2	1.80	0.46
1:A:647:GLU:O	1:A:651:GLY:N	2.46	0.45
1:B:715:LEU:HG	1:B:716:LEU:N	2.31	0.45
1:B:490:THR:HB	1:B:493:GLN:CG	2.47	0.45
1:A:459:GLY:H	1:A:473:GLY:HA2	1.81	0.45
1:A:453:ASP:N	1:A:521:LYS:HG3	2.32	0.45
1:A:719:SER:O	1:A:720:LEU:HB2	2.16	0.45
1:A:512:ILE:CD1	1:A:566:LEU:HG	2.47	0.45
1:B:719:SER:O	1:B:720:LEU:HD22	2.17	0.45
1:A:504:LEU:HD22	1:A:515:PHE:HB2	1.98	0.45
1:A:663:GLN:O	1:A:664:ILE:C	2.60	0.45
1:B:706:PHE:N	1:B:707:PRO:CD	2.80	0.45
1:A:452:PRO:HD2	1:A:455:GLN:HE22	1.82	0.45
1:A:688:MET:O	1:A:691:LEU:N	2.50	0.44
1:B:475:TRP:C	1:B:477:GLY:H	2.25	0.44
1:B:521:LYS:CB	1:B:522:PRO:CD	2.92	0.44
1:A:654:PRO:HB2	1:A:655:TYR:CE2	2.52	0.44
1:A:640:ALA:O	1:A:643:ILE:N	2.51	0.44
1:B:616:ILE:HG22	1:B:619:MET:HE3	1.98	0.44
1:B:647:GLU:O	1:B:651:GLY:N	2.44	0.44
1:A:594:PHE:CD1	1:A:594:PHE:N	2.86	0.44
1:B:629:LYS:C	1:B:631:PRO:HD3	2.43	0.44
1:B:514:LEU:O	1:B:528:THR:HB	2.18	0.44
1:A:514:LEU:HD12	1:A:515:PHE:N	2.30	0.43
1:A:712:SER:O	1:A:716:LEU:HG	2.18	0.43
1:A:514:LEU:O	1:A:528:THR:HB	2.17	0.43
1:B:457:THR:HB	1:B:474:LYS:HB2	1.99	0.43
1:B:617:LEU:HD13	1:B:654:PRO:HD2	2.00	0.43
1:B:573:HIS:O	1:B:574:ARG:HB2	2.18	0.43
1:B:501:VAL:O	1:B:505:ARG:HG2	2.18	0.43
1:A:697:LYS:HB2	1:A:703:ARG:HG2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:594:PHE:CD1	1:B:594:PHE:N	2.86	0.43
1:B:453:ASP:CB	1:B:521:LYS:HB3	2.46	0.43
1:B:720:LEU:HB2	1:B:721:PRO:HD3	2.00	0.43
1:A:508:ARG:HB2	1:B:449:TRP:CZ3	2.54	0.43
1:B:529:GLN:HE21	1:B:529:GLN:HB2	1.67	0.43
1:A:521:LYS:CB	1:A:522:PRO:CD	2.94	0.43
1:B:655:TYR:CD1	1:B:673:LEU:HD22	2.54	0.43
1:B:584:HIS:O	1:B:587:LEU:O	2.37	0.43
1:B:691:LEU:HA	1:B:694:GLU:HB2	2.00	0.42
1:A:688:MET:O	1:A:689:LYS:C	2.62	0.42
1:A:490:THR:CG2	1:A:491:PRO:N	2.82	0.42
1:A:695:CYS:O	1:A:703:ARG:CD	2.68	0.42
1:B:712:SER:O	1:B:716:LEU:HG	2.20	0.42
1:A:649:MET:O	1:A:681:ARG:HG3	2.19	0.42
1:B:676:ASP:O	1:B:678:SER:N	2.53	0.42
1:A:687:ALA:CB	1:A:716:LEU:HD22	2.46	0.42
1:B:695:CYS:C	1:B:697:LYS:H	2.26	0.42
1:B:697:LYS:O	1:B:703:ARG:NH1	2.52	0.42
1:B:453:ASP:N	1:B:521:LYS:HG3	2.34	0.42
1:A:450:GLU:HG3	1:A:518:TYR:CE1	2.54	0.42
1:A:522:PRO:O	1:A:523:GLN:CB	2.68	0.42
1:A:643:ILE:O	1:A:647:GLU:HG3	2.20	0.42
1:A:695:CYS:C	1:A:697:LYS:H	2.26	0.42
1:A:706:PHE:N	1:A:707:PRO:CD	2.83	0.42
1:B:459:GLY:H	1:B:473:GLY:HA2	1.84	0.42
1:B:522:PRO:O	1:B:523:GLN:CB	2.68	0.42
1:A:559:THR:HG23	1:A:581:ILE:HD11	2.01	0.42
1:A:617:LEU:CD1	1:A:653:LEU:HD22	2.49	0.42
1:A:655:TYR:CD1	1:A:673:LEU:HD22	2.55	0.42
1:A:584:HIS:O	1:A:587:LEU:O	2.38	0.42
1:A:650:THR:O	1:A:652:GLN:HG3	2.19	0.42
1:A:715:LEU:HD13	1:A:718:ARG:NH2	2.35	0.42
1:A:619:MET:CB	1:A:624:ILE:HD11	2.50	0.41
1:B:598:THR:HG22	1:B:599:GLU:N	2.19	0.41
1:B:687:ALA:CB	1:B:716:LEU:HD22	2.45	0.41
1:A:457:THR:HB	1:A:474:LYS:HB2	2.02	0.41
1:A:688:MET:C	1:A:690:ARG:N	2.75	0.41
1:B:638:VAL:HG13	1:B:709:ILE:HD11	2.02	0.41
1:B:650:THR:O	1:B:652:GLN:HG3	2.20	0.41
1:B:585:GLU:O	1:B:587:LEU:CD2	2.66	0.41
1:A:622:GLU:OE1	1:A:703:ARG:NH2	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:720:LEU:N	1:A:721:PRO:HD3	2.36	0.41
1:B:563:MET:CE	1:B:576:LEU:HB2	2.50	0.41
1:B:640:ALA:O	1:B:641:PHE:C	2.63	0.41
1:B:649:MET:O	1:B:681:ARG:HG3	2.20	0.41
1:A:529:GLN:HE21	1:A:529:GLN:HB2	1.70	0.41
1:A:616:ILE:HG22	1:A:619:MET:CE	2.51	0.41
1:A:563:MET:CE	1:A:576:LEU:HB2	2.51	0.41
1:B:531:CYS:HA	1:B:584:HIS:CE1	2.55	0.41
1:B:688:MET:O	1:B:689:LYS:C	2.63	0.41
1:A:549:MET:O	1:A:550:ILE:C	2.64	0.41
1:A:638:VAL:HG13	1:A:709:ILE:HD11	2.01	0.41
1:B:599:GLU:O	1:B:600:LYS:HD3	2.20	0.41
1:A:676:ASP:O	1:A:679:LYS:HG3	2.20	0.41
1:A:691:LEU:HA	1:A:694:GLU:HB2	2.02	0.41
1:B:490:THR:CG2	1:B:491:PRO:N	2.84	0.41
1:B:512:ILE:CD1	1:B:566:LEU:HG	2.51	0.41
1:B:715:LEU:O	1:B:717:ALA:N	2.45	0.41
1:A:531:CYS:HA	1:A:584:HIS:CE1	2.56	0.41
1:B:663:GLN:O	1:B:664:ILE:C	2.64	0.40
1:A:561:GLN:HG3	1:A:710:LEU:HD11	2.03	0.40
1:A:593:ASP:C	1:A:595:GLY:H	2.29	0.40
1:A:677:LEU:O	1:A:689:LYS:HE3	2.22	0.40
1:A:475:TRP:C	1:A:477:GLY:H	2.29	0.40
1:A:476:HIS:CE1	1:B:568:ALA:HB2	2.55	0.40
1:A:638:VAL:CG1	1:A:709:ILE:HD11	2.51	0.40
1:B:688:MET:O	1:B:691:LEU:N	2.55	0.40
1:A:568:ALA:C	1:A:570:SER:H	2.28	0.40
1:A:617:LEU:HD13	1:A:654:PRO:HD2	2.03	0.40
1:B:622:GLU:OE1	1:B:703:ARG:NH2	2.41	0.40
1:B:519:SER:N	1:B:525:ALA:O	2.55	0.40
1:B:676:ASP:O	1:B:679:LYS:HG3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	260/276 (94%)	204 (78%)	40 (15%)	16 (6%)	1	12
1	B	259/276 (94%)	203 (78%)	41 (16%)	15 (6%)	1	13
All	All	519/552 (94%)	407 (78%)	81 (16%)	31 (6%)	1	12

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	522	PRO
1	A	628	ASP
1	A	720	LEU
1	B	486	VAL
1	B	487	THR
1	B	522	PRO
1	B	628	ASP
1	A	486	VAL
1	A	487	THR
1	A	588	THR
1	A	600	LYS
1	B	588	THR
1	B	677	LEU
1	B	721	PRO
1	A	677	LEU
1	B	594	PHE
1	B	629	LYS
1	B	720	LEU
1	A	521	LYS
1	A	594	PHE
1	A	615	SER
1	A	655	TYR
1	B	521	LYS
1	A	587	LEU
1	B	587	LEU
1	B	716	LEU
1	A	716	LEU
1	B	457	THR
1	A	721	PRO
1	A	458	VAL
1	B	458	VAL

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	232/243 (96%)	210 (90%)	22 (10%)	8	31
1	B	231/243 (95%)	216 (94%)	15 (6%)	15	42
All	All	463/486 (95%)	426 (92%)	37 (8%)	11	35

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	456	ILE
1	A	457	THR
1	A	486	VAL
1	A	490	THR
1	A	504	LEU
1	A	510	VAL
1	A	522	PRO
1	A	529	GLN
1	A	534	SER
1	A	541	HIS
1	A	550	ILE
1	A	579	ASN
1	A	598	THR
1	A	600	LYS
1	A	616	ILE
1	A	621	PRO
1	A	626	MET
1	A	627	GLN
1	A	628	ASP
1	A	673	LEU
1	A	701	ASP
1	A	715	LEU
1	B	456	ILE
1	B	457	THR
1	B	486	VAL
1	B	490	THR
1	B	510	VAL

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Mol	Chain	Res	Type
1	B	522	PRO
1	B	529	GLN
1	B	541	HIS
1	B	550	ILE
1	B	600	LYS
1	B	673	LEU
1	B	694	GLU
1	B	701	ASP
1	B	715	LEU
1	B	721	PRO

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

Mol	Chain	Res	Type
1	A	455	GLN
1	A	476	HIS
1	A	492	GLN
1	A	493	GLN
1	A	499	ASN
1	A	509	HIS
1	A	511	ASN
1	A	529	GLN
1	A	538	HIS
1	A	558	GLN
1	A	579	ASN
1	A	659	ASN
1	B	455	GLN
1	B	476	HIS
1	B	493	GLN
1	B	499	ASN
1	B	509	HIS
1	B	511	ASN
1	B	529	GLN
1	B	558	GLN
1	B	579	ASN
1	B	659	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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	BAX	B	1723	-	34,34,34	3.83	13 (38%)	48,48,48	4.21	17 (35%)
2	BAX	A	1723	-	34,34,34	3.62	13 (38%)	48,48,48	4.09	16 (33%)

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	BAX	B	1723	-	-	11/24/24/24	0/3/3/3
2	BAX	A	1723	-	-	12/24/24/24	0/3/3/3

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1723	BAX	C29-N30	12.54	1.50	1.33
2	A	1723	BAX	C29-N30	11.93	1.49	1.33
2	B	1723	BAX	C16-N14	8.70	1.59	1.41
2	A	1723	BAX	C16-N14	8.55	1.58	1.41
2	B	1723	BAX	C13-N12	8.04	1.55	1.37
2	A	1723	BAX	C13-N12	7.73	1.54	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1723	BAX	C25-N26	5.70	1.46	1.34
2	A	1723	BAX	C25-N26	5.38	1.45	1.34
2	B	1723	BAX	C24-C23	5.10	1.48	1.38
2	B	1723	BAX	C7-C6	-4.93	1.39	1.50
2	A	1723	BAX	C24-C23	4.91	1.47	1.38
2	B	1723	BAX	C17-C16	4.83	1.47	1.39
2	B	1723	BAX	C5-C6	4.81	1.47	1.39
2	B	1723	BAX	C27-N26	4.68	1.42	1.34
2	A	1723	BAX	C7-C6	-4.34	1.41	1.50
2	A	1723	BAX	C17-C16	4.31	1.46	1.39
2	A	1723	BAX	C27-N26	4.25	1.42	1.34
2	A	1723	BAX	C28-C23	-3.95	1.32	1.39
2	A	1723	BAX	C5-C6	3.51	1.45	1.39
2	A	1723	BAX	C28-C27	-2.94	1.32	1.40
2	B	1723	BAX	C28-C23	-2.94	1.34	1.39
2	B	1723	BAX	C28-C27	-2.87	1.32	1.40
2	B	1723	BAX	C5-CL11	2.57	1.79	1.73
2	B	1723	BAX	C27-C29	2.26	1.56	1.50
2	A	1723	BAX	C27-C29	2.16	1.56	1.50
2	A	1723	BAX	C13-N14	2.03	1.41	1.37

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1723	BAX	C27-C29-N30	20.39	129.53	115.56
2	A	1723	BAX	C27-C29-N30	19.40	128.85	115.56
2	B	1723	BAX	C2-N12-C13	9.45	145.89	126.61
2	A	1723	BAX	C2-N12-C13	8.21	143.37	126.61
2	A	1723	BAX	C29-C27-N26	7.93	126.69	117.42
2	B	1723	BAX	C29-C27-N26	7.82	126.56	117.42
2	A	1723	BAX	O15-C13-N12	-7.60	110.13	123.64
2	B	1723	BAX	O15-C13-N12	-7.59	110.16	123.64
2	A	1723	BAX	N14-C13-N12	7.06	126.10	112.44
2	B	1723	BAX	O32-C29-N30	-6.38	113.24	122.67
2	A	1723	BAX	O32-C29-N30	-6.12	113.61	122.67
2	B	1723	BAX	N14-C13-N12	5.90	123.86	112.44
2	A	1723	BAX	C31-N30-C29	-5.47	111.39	121.80
2	B	1723	BAX	C31-N30-C29	-5.19	111.92	121.80
2	A	1723	BAX	C24-C25-N26	-5.00	117.85	123.97
2	B	1723	BAX	C24-C25-N26	-4.71	118.20	123.97
2	B	1723	BAX	C7-C6-C5	4.02	124.72	121.86
2	B	1723	BAX	O15-C13-N14	-3.52	117.39	123.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1723	BAX	C7-C6-C5	3.47	124.33	121.86
2	B	1723	BAX	O32-C29-C27	-3.37	113.64	121.08
2	A	1723	BAX	O15-C13-N14	-3.17	118.00	123.64
2	B	1723	BAX	C28-C27-N26	-2.95	119.73	123.27
2	A	1723	BAX	C28-C27-N26	-2.82	119.89	123.27
2	A	1723	BAX	O32-C29-C27	-2.78	114.94	121.08
2	A	1723	BAX	O22-C23-C28	-2.70	110.69	119.16
2	B	1723	BAX	O22-C23-C28	-2.64	110.89	119.16
2	A	1723	BAX	O22-C23-C24	2.63	128.53	119.36
2	A	1723	BAX	C16-N14-C13	-2.59	121.33	126.61
2	B	1723	BAX	O22-C23-C24	2.56	128.26	119.36
2	B	1723	BAX	C25-N26-C27	2.35	119.96	116.93
2	A	1723	BAX	C25-N26-C27	2.32	119.92	116.93
2	B	1723	BAX	F8-C7-C6	-2.30	108.57	112.65
2	B	1723	BAX	C6-C5-CL11	2.10	124.24	121.70

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1723	BAX	C27-C29-N30-C31
2	A	1723	BAX	O32-C29-N30-C31
2	A	1723	BAX	C5-C6-C7-F8
2	A	1723	BAX	C5-C6-C7-F9
2	A	1723	BAX	C5-C6-C7-F10
2	B	1723	BAX	C27-C29-N30-C31
2	B	1723	BAX	O32-C29-N30-C31
2	B	1723	BAX	C28-C27-C29-O32
2	A	1723	BAX	C28-C27-C29-N30
2	A	1723	BAX	N26-C27-C29-N30
2	A	1723	BAX	C28-C27-C29-O32
2	B	1723	BAX	C28-C27-C29-N30
2	B	1723	BAX	N26-C27-C29-O32
2	B	1723	BAX	N26-C27-C29-N30
2	A	1723	BAX	N26-C27-C29-O32
2	B	1723	BAX	O15-C13-N12-C2
2	A	1723	BAX	O15-C13-N12-C2
2	B	1723	BAX	O15-C13-N14-C16
2	B	1723	BAX	C3-C2-N12-C13
2	B	1723	BAX	C1-C2-N12-C13
2	B	1723	BAX	C5-C6-C7-F8
2	A	1723	BAX	O15-C13-N14-C16

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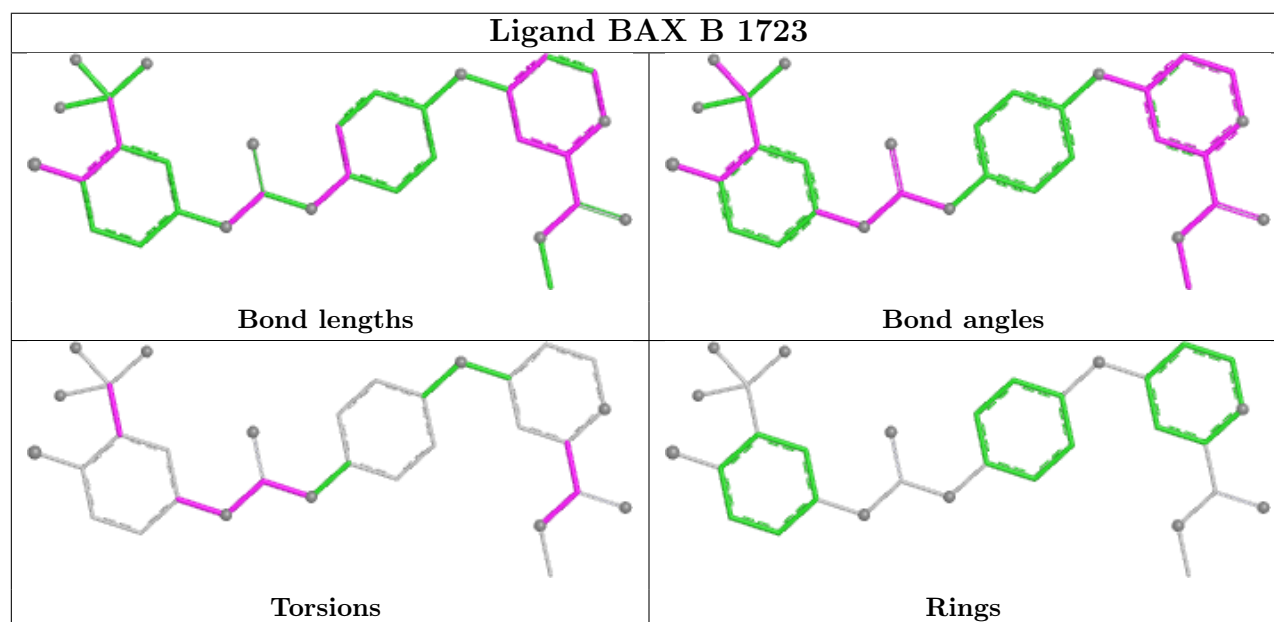
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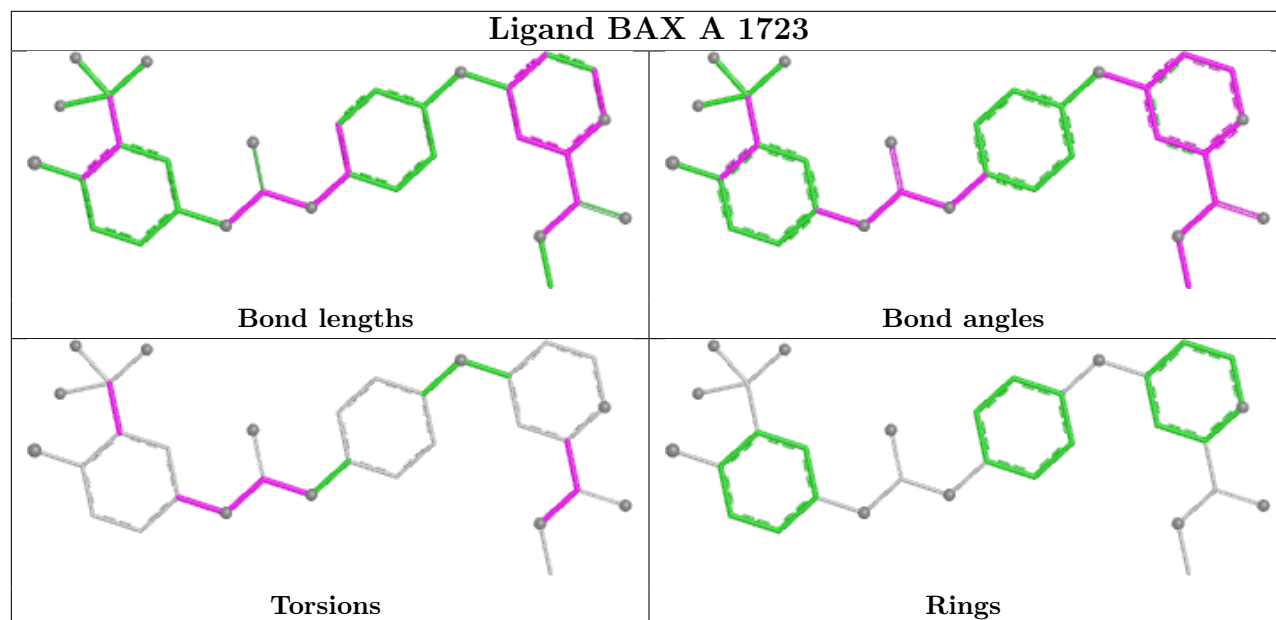
Mol	Chain	Res	Type	Atoms
2	A	1723	BAX	C3-C2-N12-C13

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	264/276 (95%)	0.31	8 (3%) 52 30	4, 4, 4, 4	0
1	B	263/276 (95%)	0.37	10 (3%) 44 24	7, 7, 7, 7	0
All	All	527/552 (95%)	0.34	18 (3%) 48 27	4, 4, 7, 7	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	721	PRO	4.0
1	B	721	PRO	3.8
1	B	453	ASP	3.7
1	B	447	ASP	3.4
1	A	628	ASP	3.2
1	B	628	ASP	2.5
1	B	598	THR	2.5
1	A	601	SER	2.5
1	B	629	LYS	2.4
1	A	487	THR	2.4
1	A	722	LYS	2.4
1	B	599	GLU	2.4
1	B	722	LYS	2.3
1	A	447	ASP	2.1
1	B	495	GLN	2.1
1	A	684	CYS	2.1
1	B	487	THR	2.0
1	A	599	GLU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

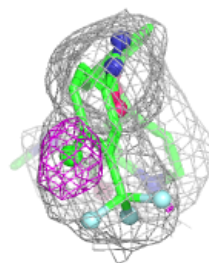
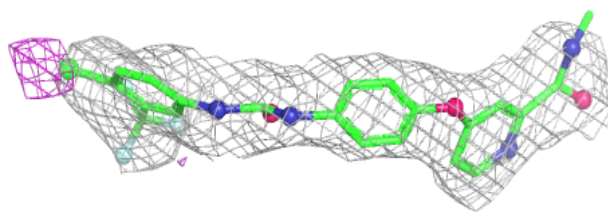
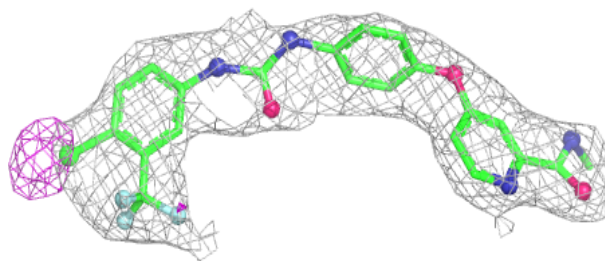
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

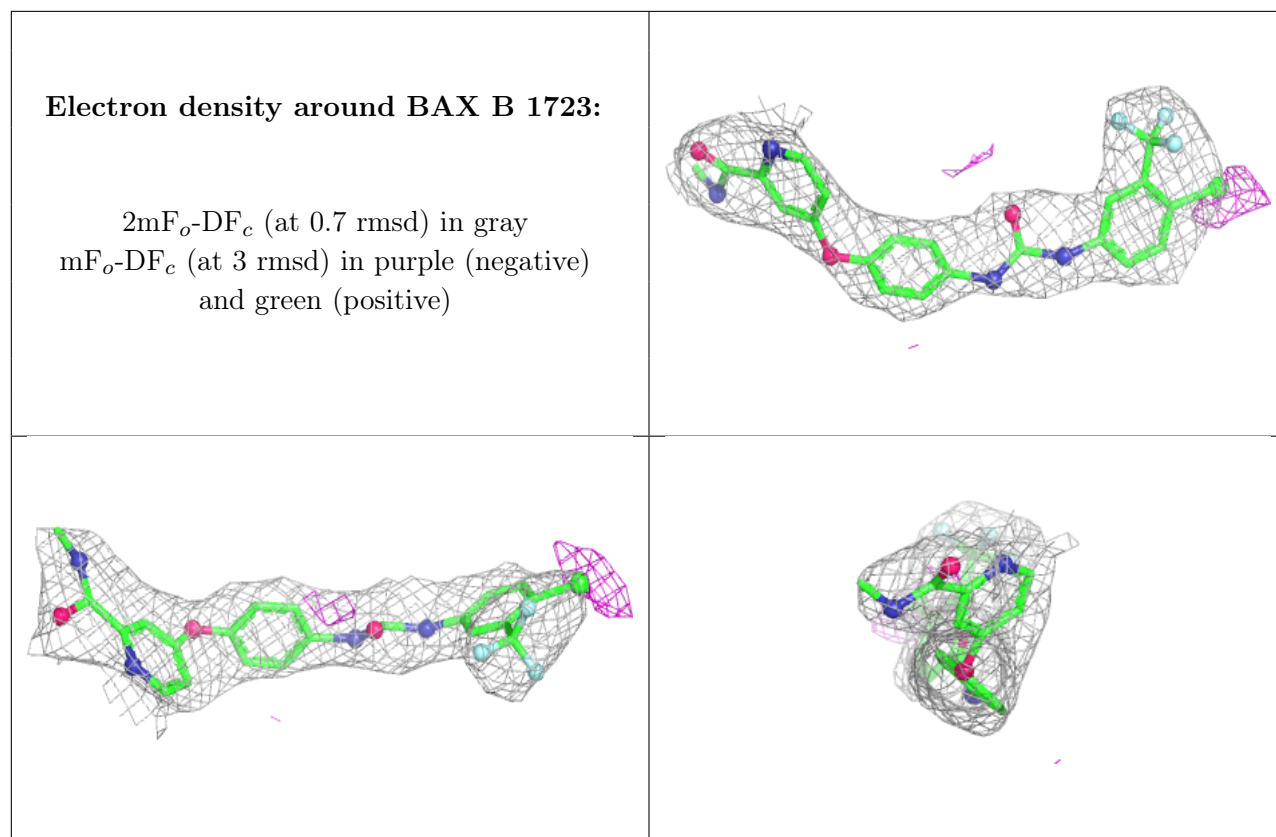
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BAX	A	1723	32/32	0.90	0.14	2,2,2,2	0
2	BAX	B	1723	32/32	0.90	0.13	2,2,2,2	0

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 BAX A 1723:

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.