



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:57 AM GMT

PDB ID : 2JIU
Title : CRYSTAL STRUCTURE OF EGFR KINASE DOMAIN T790M MUTATION
IN COMPLEX WITH AEE788
Authors : Yun, C.-H.; Mengwasser, K.E.; Toms, A.V.; Woo, M.S.; Greulich, H.; Wong,
K.-K.; Meyerson, M.; Eck, M.J.
Deposited on : 2007-07-01
Resolution : 3.05 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

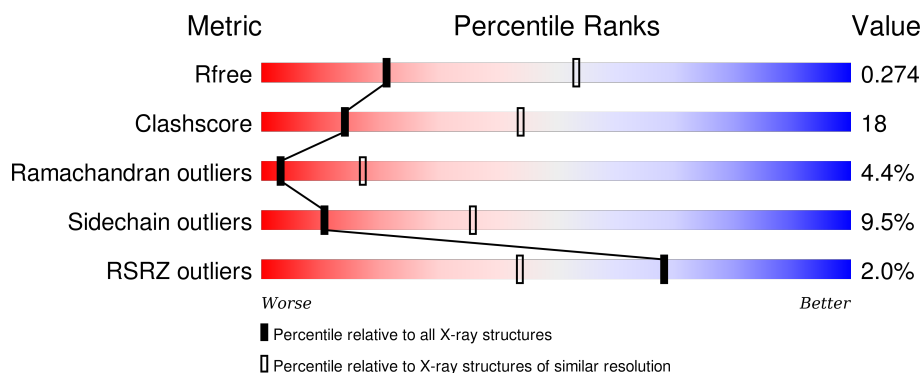
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

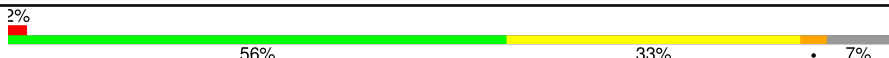
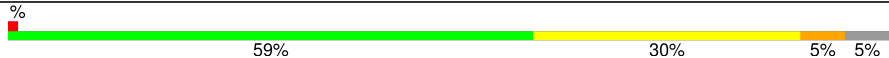
The reported resolution of this entry is 3.05 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1191 (3.10-3.02)
Clashscore	102246	1303 (3.10-3.02)
Ramachandran outliers	100387	1254 (3.10-3.02)
Sidechain outliers	100360	1254 (3.10-3.02)
RSRZ outliers	91569	1197 (3.10-3.02)

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

Mol	Chain	Length	Quality of chain
1	A	328	 2% 56% 33% 7%
1	B	328	 % 59% 30% 5% 5%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4880 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.

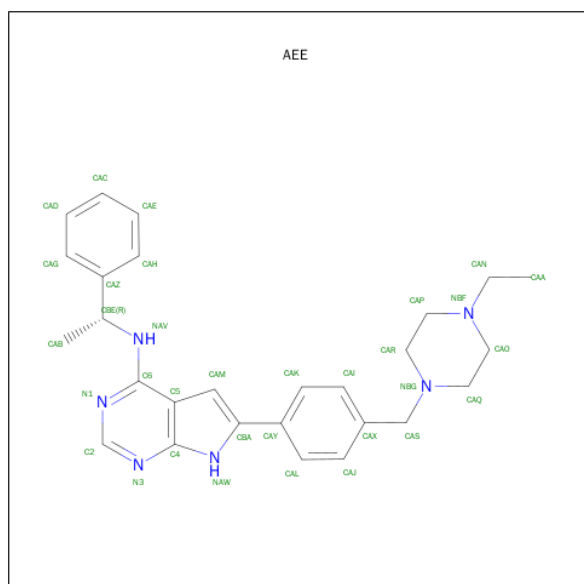
- Molecule 1 is a protein called EPIDERMAL GROWTH FACTOR RECEPTOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	1
			2380	1531	406	426	17			
1	B	310	Total	C	N	O	S	0	0	1
			2407	1544	415	430	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	790	MET	THR	ENGINEERED MUTATION	UNP P00533
B	790	MET	THR	ENGINEERED MUTATION	UNP P00533

- Molecule 2 is 6-{4-[(4-ETHYLPIPERAZIN-1-YL)METHYL]PHENYL}-N-[(1R)-1-PHENYLETHYL]-7H-PYRROLO[2,3-D]PYRIMIDIN-4-AMINE (three-letter code: AEE) (formula: C₂₇H₃₂N₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			33	27	6		

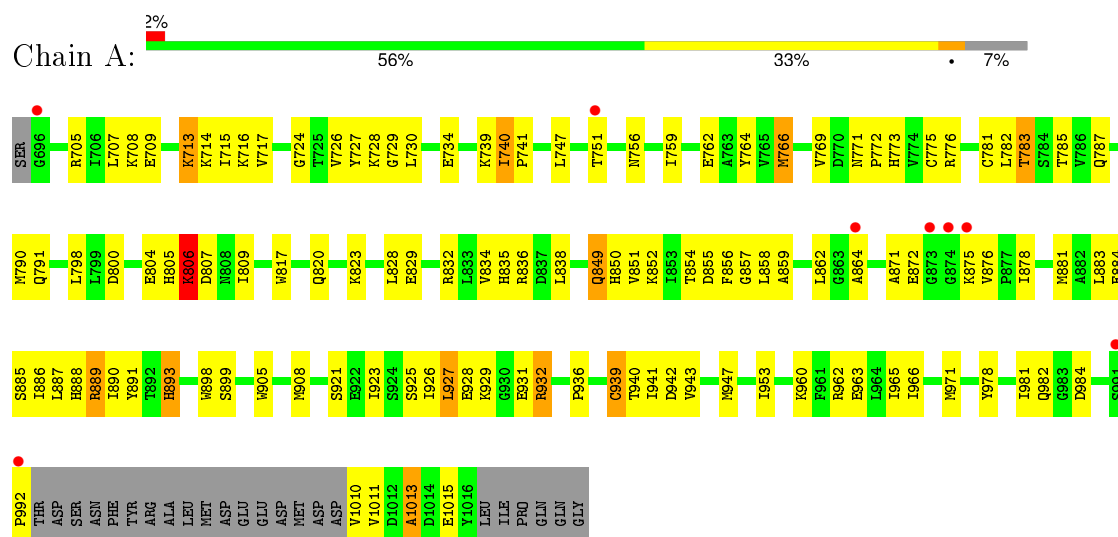
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	38	Total	O	0	0
			38	38		
3	B	22	Total	O	0	0
			22	22		

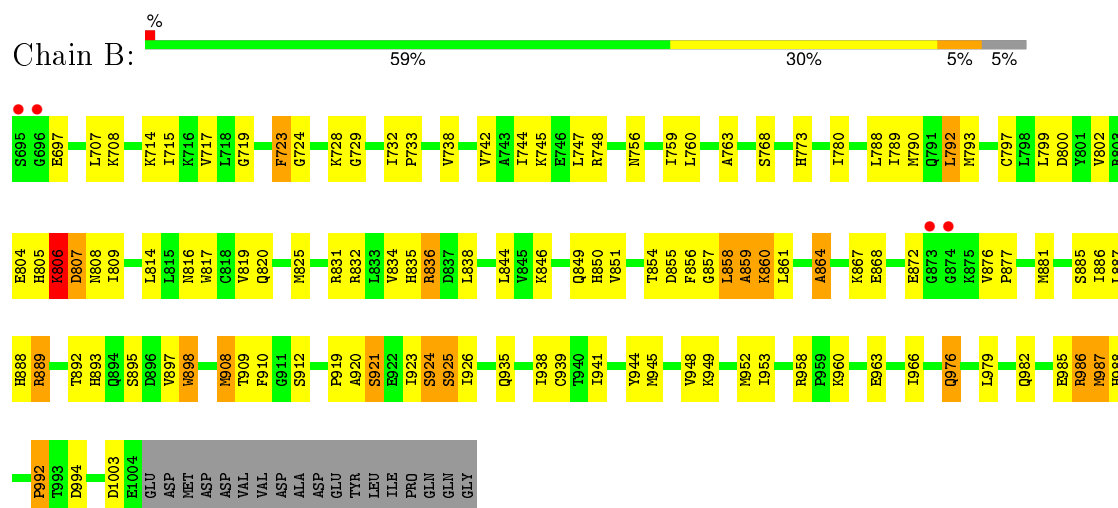
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: EPIDERMAL GROWTH FACTOR RECEPTOR



• Molecule 1: EPIDERMAL GROWTH FACTOR RECEPTOR



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	48.41Å 88.64Å 164.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.12 – 3.05 24.12 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.8 (24.12-3.05) 99.8 (24.12-3.05)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 3.05Å)	Xtriage
Refinement program	REFMAC 5.3.0026	Depositor
R, R_{free}	0.211 , 0.277 0.206 , 0.274	Depositor DCC
R_{free} test set	716 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	72.6	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 33.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 14082 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4880	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.67	0/2431	0.67	1/3293 (0.0%)
1	B	0.68	0/2459	0.70	1/3333 (0.0%)
All	All	0.67	0/4890	0.69	2/6626 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	992	PRO	N-CA-CB	5.91	110.39	103.30
1	B	992	PRO	N-CA-CB	5.88	110.36	103.30

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	2380	0	2398	85	0
1	B	2407	0	2403	91	0
2	A	33	0	32	2	0
3	A	38	0	0	4	0
3	B	22	0	0	4	0
All	All	4880	0	4833	176	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 18.

All (176) 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:836:ARG:NH1	1:B:859:ALA:CB	1.86	1.37
1:B:836:ARG:NH1	1:B:859:ALA:HB3	0.97	1.28
1:B:834:VAL:H	1:B:860:LYS:CB	1.54	1.19
1:B:836:ARG:HH12	1:B:859:ALA:CB	1.54	1.01
1:B:835:HIS:O	1:B:836:ARG:HB2	1.58	1.01
1:B:808:ASN:HD22	1:B:988:HIS:HA	1.22	1.01
1:A:791:GLN:HE21	1:A:1013:ALA:HB3	1.28	0.99
1:A:790:MET:HE1	3:A:3016:HOH:O	1.69	0.91
1:A:730:LEU:HD23	1:A:739:LYS:HB3	1.53	0.89
1:B:807:ASP:O	1:B:986:ARG:CZ	2.22	0.88
1:B:836:ARG:HH11	1:B:859:ALA:HB3	1.09	0.87
1:A:716:LYS:HE2	1:A:728:LYS:HD3	1.55	0.87
1:B:834:VAL:N	1:B:860:LYS:CB	2.36	0.87
1:A:881:MET:HB2	1:A:886:ILE:HD11	1.58	0.86
1:A:940:THR:HB	1:A:978:TYR:O	1.75	0.85
1:B:808:ASN:ND2	1:B:988:HIS:HA	1.95	0.81
1:B:836:ARG:HH12	1:B:859:ALA:HB3	1.07	0.80
1:A:724:GLY:HA3	1:A:747:LEU:HA	1.62	0.80
1:A:856:PHE:HB3	1:A:859:ALA:HB2	1.64	0.79
1:B:808:ASN:HD22	1:B:988:HIS:CA	1.99	0.76
1:B:825:MET:CE	1:B:838:LEU:HD22	2.16	0.76
1:B:836:ARG:CZ	1:B:859:ALA:HB3	2.07	0.75
2:A:2017:AEE:HAO1	3:A:3038:HOH:O	1.86	0.74
1:B:715:ILE:HG12	1:B:729:GLY:HA2	1.70	0.74
1:B:790:MET:HE2	1:B:855:ASP:H	1.53	0.73
1:B:836:ARG:HH12	1:B:859:ALA:HB1	1.54	0.73
1:B:835:HIS:O	1:B:836:ARG:CB	2.39	0.71
1:B:707:LEU:HD12	1:B:789:ILE:HG13	1.74	0.69
1:B:800:ASP:O	1:B:804:GLU:HG3	1.94	0.68
1:A:856:PHE:HB3	1:A:859:ALA:CB	2.23	0.68
1:A:716:LYS:HE2	1:A:728:LYS:CD	2.22	0.67
1:B:953:ILE:HD11	3:B:3014:HOH:O	1.93	0.67
1:A:836:ARG:HE	1:A:858:LEU:HB3	1.60	0.67
1:A:923:ILE:HG22	1:A:927:LEU:HD22	1.75	0.67
1:A:747:LEU:HD11	1:A:759:ILE:HD12	1.76	0.66
1:A:820:GLN:NE2	1:A:851:VAL:HG22	2.11	0.66
1:B:825:MET:HE1	1:B:838:LEU:HD22	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:790:MET:HE2	1:B:855:ASP:OD2	1.97	0.64
1:B:806:LYS:HE2	1:B:910:PHE:HB2	1.78	0.64
1:A:800:ASP:O	1:A:804:GLU:HG2	1.97	0.64
1:A:835:HIS:O	1:A:836:ARG:HB2	1.97	0.64
1:B:908:MET:HG2	1:B:939:CYS:SG	2.39	0.63
1:A:936:PRO:HB2	1:A:939:CYS:HB2	1.80	0.63
1:A:769:VAL:O	1:A:776:ARG:HD3	1.99	0.62
1:A:791:GLN:NE2	1:A:1013:ALA:HB3	2.09	0.61
1:B:805:HIS:O	1:B:807:ASP:N	2.33	0.61
1:B:846:LYS:HD3	1:B:850:HIS:CD2	2.34	0.61
1:A:883:LEU:HD22	1:A:953:ILE:HD12	1.82	0.61
1:A:971:MET:HG2	1:A:978:TYR:CD2	2.35	0.61
1:A:905:TRP:HD1	1:A:947:MET:CE	2.15	0.60
1:A:740:ILE:HD12	1:A:741:PRO:HD2	1.83	0.60
1:A:762:GLU:HG2	1:A:766:MET:HE2	1.83	0.59
1:A:836:ARG:HG2	1:A:891:TYR:CD1	2.38	0.58
1:A:829:GLU:HG3	1:A:893:HIS:HD2	1.68	0.58
1:B:857:GLY:C	1:B:859:ALA:H	2.07	0.58
1:A:963:GLU:O	1:A:966:ILE:HG13	2.03	0.58
1:A:740:ILE:HD13	1:A:1013:ALA:N	2.19	0.58
1:A:836:ARG:HH22	1:A:871:ALA:HB2	1.69	0.58
1:A:715:ILE:HG12	1:A:729:GLY:HA2	1.86	0.57
1:B:976:GLN:HE21	1:B:976:GLN:H	1.52	0.57
1:B:858:LEU:C	1:B:860:LYS:H	2.08	0.57
1:B:763:ALA:HB2	1:B:788:LEU:HD21	1.86	0.57
1:A:881:MET:HB2	1:A:886:ILE:CD1	2.32	0.57
1:B:898:TRP:HD1	1:B:958:ARG:CZ	2.17	0.56
1:B:923:ILE:O	1:B:925:SER:N	2.38	0.56
1:B:938:ILE:HD12	1:B:979:LEU:HD22	1.87	0.56
1:B:806:LYS:HE3	1:B:910:PHE:O	2.05	0.56
1:A:931:GLU:O	1:A:932:ARG:HD3	2.06	0.56
1:A:782:LEU:HD11	1:B:948:VAL:HG13	1.87	0.56
1:B:825:MET:HE2	1:B:838:LEU:HD22	1.88	0.56
1:B:976:GLN:H	1:B:976:GLN:NE2	2.05	0.55
1:A:940:THR:HG22	1:A:942:ASP:H	1.70	0.55
1:B:858:LEU:O	1:B:860:LYS:N	2.40	0.54
1:A:926:ILE:HG13	1:A:927:LEU:N	2.21	0.54
1:A:828:LEU:HD11	1:A:856:PHE:HE1	1.72	0.54
1:A:836:ARG:HD3	1:A:858:LEU:O	2.08	0.54
1:A:829:GLU:CG	1:A:893:HIS:HD2	2.21	0.54
1:B:909:THR:HB	1:B:912:SER:HB2	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:885:SER:O	1:B:889:ARG:HA	2.08	0.53
1:A:714:LYS:HD3	1:A:727:TYR:CE2	2.44	0.53
1:A:838:LEU:HB3	1:A:899:SER:HB3	1.90	0.53
1:A:883:LEU:HD21	1:A:928:GLU:HG3	1.90	0.53
1:A:714:LYS:HD3	1:A:727:TYR:CD2	2.43	0.53
1:B:742:VAL:O	1:B:792:LEU:HD22	2.09	0.53
1:B:728:LYS:NZ	3:B:3002:HOH:O	2.42	0.52
1:A:940:THR:HG22	1:A:942:ASP:N	2.24	0.52
1:A:971:MET:HG2	1:A:978:TYR:CG	2.45	0.52
1:A:805:HIS:O	1:A:809:ILE:HG12	2.09	0.52
1:A:740:ILE:HD12	1:A:741:PRO:CD	2.39	0.52
1:B:908:MET:CG	1:B:939:CYS:SG	2.98	0.51
1:A:805:HIS:O	1:A:807:ASP:N	2.41	0.51
1:B:834:VAL:HB	1:B:860:LYS:CB	2.40	0.51
1:B:855:ASP:HB2	1:B:858:LEU:HD11	1.93	0.51
1:A:713:LYS:N	1:A:713:LYS:HD2	2.26	0.51
1:A:878:ILE:HG23	1:A:886:ILE:HD12	1.93	0.50
1:A:925:SER:O	1:A:929:LYS:HG2	2.12	0.50
1:B:986:ARG:O	1:B:988:HIS:N	2.42	0.50
1:A:790:MET:CE	3:A:3016:HOH:O	2.43	0.50
1:B:923:ILE:HA	1:B:926:ILE:HG12	1.93	0.49
1:B:876:VAL:HG21	1:B:889:ARG:NH2	2.27	0.49
1:A:805:HIS:C	1:A:807:ASP:H	2.15	0.49
1:A:960:LYS:HD2	1:A:962:ARG:HH11	1.77	0.49
1:B:893:HIS:O	1:B:897:VAL:HG23	2.12	0.49
1:A:806:LYS:HD3	1:A:806:LYS:O	2.12	0.49
1:A:1013:ALA:C	1:A:1015:GLU:H	2.15	0.49
1:B:945:MET:HG3	1:B:949:LYS:HE3	1.95	0.48
1:A:943:VAL:HG22	1:A:971:MET:SD	2.53	0.48
1:A:849:GLN:H	1:A:849:GLN:NE2	2.10	0.48
1:B:773:HIS:CE1	1:B:820:GLN:NE2	2.81	0.48
1:B:857:GLY:C	1:B:859:ALA:N	2.66	0.48
1:B:790:MET:CE	1:B:855:ASP:H	2.24	0.48
1:B:858:LEU:C	1:B:860:LYS:N	2.67	0.48
1:A:905:TRP:CD1	1:A:947:MET:CE	2.96	0.48
1:A:705:ARG:NE	3:A:3002:HOH:O	2.41	0.48
1:B:748:ARG:HG2	3:B:3004:HOH:O	2.12	0.48
1:A:829:GLU:HG3	1:A:893:HIS:CD2	2.49	0.48
1:B:724:GLY:HA2	3:B:3004:HOH:O	2.14	0.48
1:B:747:LEU:HG	1:B:759:ILE:HD13	1.95	0.48
1:B:804:GLU:O	1:B:805:HIS:ND1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:775:CYS:HB3	1:A:790:MET:CE	2.43	0.47
1:A:856:PHE:O	1:A:858:LEU:N	2.48	0.47
1:A:923:ILE:O	1:A:926:ILE:HG12	2.14	0.47
1:B:877:PRO:O	1:B:881:MET:HG3	2.15	0.47
1:A:888:HIS:O	1:A:890:ILE:HG12	2.14	0.47
1:B:935:GLN:HA	1:B:944:TYR:CE2	2.50	0.47
1:A:771:ASN:OD1	1:A:772:PRO:HD2	2.15	0.46
1:A:881:MET:HB3	1:A:885:SER:HB2	1.98	0.46
1:A:809:ILE:HD12	1:A:817:TRP:HH2	1.79	0.46
1:B:793:MET:HE2	1:B:793:MET:HA	1.97	0.46
1:A:834:VAL:HG12	1:A:836:ARG:HG3	1.98	0.46
1:A:850:HIS:CE1	1:A:852:LYS:HE3	2.50	0.46
1:B:889:ARG:HA	1:B:889:ARG:HD3	1.84	0.45
1:B:886:ILE:HG21	1:B:924:SER:HB3	1.98	0.45
1:B:780:ILE:HG22	1:B:788:LEU:HD23	1.98	0.45
2:A:2017:AEE:HAN2	2:A:2017:AEE:HAQ2	1.79	0.45
1:A:941:ILE:HG23	1:A:942:ASP:N	2.33	0.44
1:A:832:ARG:HB3	1:A:862:LEU:HB2	2.00	0.44
1:B:814:LEU:HA	1:B:817:TRP:CE3	2.53	0.44
1:B:723:PHE:O	1:B:745:LYS:HE3	2.18	0.43
1:A:713:LYS:H	1:A:713:LYS:HD2	1.84	0.43
1:B:816:ASN:O	1:B:819:VAL:HB	2.19	0.43
1:B:835:HIS:NE2	1:B:854:THR:O	2.44	0.43
1:B:820:GLN:NE2	1:B:851:VAL:H	2.16	0.43
1:B:963:GLU:HA	1:B:966:ILE:HG13	2.00	0.43
1:B:806:LYS:HG3	1:B:910:PHE:HB3	2.00	0.43
1:A:1010:VAL:HG23	1:A:1011:VAL:HG23	2.01	0.43
1:B:909:THR:HB	1:B:912:SER:CB	2.49	0.43
1:B:717:VAL:HG22	1:B:719:GLY:H	1.83	0.43
1:B:732:ILE:HG23	1:B:738:VAL:O	2.19	0.43
1:B:919:PRO:C	1:B:921:SER:H	2.22	0.42
1:B:806:LYS:HG3	1:B:910:PHE:CB	2.49	0.42
1:A:764:TYR:HE2	1:B:941:ILE:HG23	1.85	0.42
1:A:790:MET:HE2	1:A:790:MET:HB3	1.57	0.42
1:B:892:THR:HG23	1:B:895:SER:H	1.84	0.42
1:B:876:VAL:HG21	1:B:889:ARG:HH22	1.85	0.41
1:B:756:ASN:O	1:B:759:ILE:HG22	2.21	0.41
1:A:936:PRO:O	1:A:939:CYS:HB2	2.19	0.41
1:B:836:ARG:HH11	1:B:859:ALA:CB	1.90	0.41
1:B:857:GLY:O	1:B:859:ALA:N	2.45	0.41
1:A:823:LYS:HA	1:A:965:ILE:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:781:CYS:HB3	1:A:787:GLN:HB2	2.03	0.41
1:B:799:LEU:O	1:B:802:VAL:HG12	2.21	0.41
1:B:832:ARG:HH12	1:B:864:ALA:HA	1.86	0.41
1:A:905:TRP:HD1	1:A:947:MET:HE3	1.84	0.41
1:A:854:THR:HG23	1:A:855:ASP:HB2	2.01	0.41
1:A:771:ASN:ND2	1:A:773:HIS:H	2.18	0.41
1:B:768:SER:C	1:B:831:ARG:HH12	2.24	0.41
1:A:709:GLU:OE1	1:A:783:THR:HG21	2.20	0.41
1:B:952:MET:HB2	1:B:958:ARG:HG2	2.02	0.40
1:B:820:GLN:HB3	1:B:820:GLN:HE21	1.61	0.40
1:A:889:ARG:N	1:A:889:ARG:HD3	2.36	0.40
1:A:940:THR:HG22	1:A:942:ASP:HB2	2.04	0.40
1:B:797:CYS:HA	1:B:844:LEU:HA	2.04	0.40
1:A:898:TRP:C	1:A:898:TRP:CD1	2.94	0.40
1:B:744:ILE:H	1:B:744:ILE:HD12	1.86	0.40
1:B:986:ARG:C	1:B:988:HIS:H	2.24	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	300/328 (92%)	254 (85%)	40 (13%)	6 (2%)	9	36
1	B	308/328 (94%)	256 (83%)	31 (10%)	21 (7%)	1	8
All	All	608/656 (93%)	510 (84%)	71 (12%)	27 (4%)	3	17

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	806	LYS
1	A	1013	ALA

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Mol	Chain	Res	Type
1	B	733	PRO
1	B	806	LYS
1	B	887	LEU
1	B	986	ARG
1	B	987	MET
1	B	992	PRO
1	A	857	GLY
1	B	723	PHE
1	B	858	LEU
1	B	859	ALA
1	B	860	LYS
1	B	864	ALA
1	B	872	GLU
1	B	925	SER
1	B	982	GLN
1	A	864	ALA
1	A	872	GLU
1	B	836	ARG
1	B	861	LEU
1	B	888	HIS
1	B	920	ALA
1	A	734	GLU
1	B	924	SER
1	B	994	ASP
1	B	1003	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	254/288 (88%)	226 (89%)	28 (11%)	8	29
1	B	252/288 (88%)	232 (92%)	20 (8%)	15	46
All	All	506/576 (88%)	458 (90%)	48 (10%)	11	36

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	707	LEU
1	A	708	LYS
1	A	713	LYS
1	A	717	VAL
1	A	726	VAL
1	A	740	ILE
1	A	751	THR
1	A	756	ASN
1	A	766	MET
1	A	783	THR
1	A	785	THR
1	A	798	LEU
1	A	806	LYS
1	A	849	GLN
1	A	875	LYS
1	A	876	VAL
1	A	884	GLU
1	A	887	LEU
1	A	889	ARG
1	A	893	HIS
1	A	908	MET
1	A	921	SER
1	A	927	LEU
1	A	932	ARG
1	A	939	CYS
1	A	981	ILE
1	A	982	GLN
1	A	984	ASP
1	B	697	GLU
1	B	708	LYS
1	B	714	LYS
1	B	760	LEU
1	B	792	LEU
1	B	806	LYS
1	B	807	ASP
1	B	809	ILE
1	B	849	GLN
1	B	856	PHE
1	B	867	LYS
1	B	868	GLU
1	B	889	ARG
1	B	898	TRP
1	B	908	MET

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Mol	Chain	Res	Type
1	B	921	SER
1	B	960	LYS
1	B	976	GLN
1	B	985	GLU
1	B	987	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	756	ASN
1	A	771	ASN
1	A	787	GLN
1	A	791	GLN
1	A	826	ASN
1	A	849	GLN
1	A	893	HIS
1	A	935	GLN
1	B	756	ASN
1	B	808	ASN
1	B	820	GLN
1	B	842	ASN
1	B	849	GLN
1	B	850	HIS
1	B	976	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 ⓘ

1 ligand is 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	AEE	A	2017	-	35,37,37	1.57	3 (8%)	41,51,51	2.11	5 (12%)

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	AEE	A	2017	-	-	0/18/28/28	0/5/5/5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2017	AEE	CAY-CBA	-6.23	1.39	1.48
2	A	2017	AEE	C6-C5	-4.86	1.39	1.44
2	A	2017	AEE	CAM-CBA	-3.38	1.33	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2017	AEE	N3-C2-N1	-9.41	121.69	128.89
2	A	2017	AEE	C5-C6-N1	-3.55	118.93	121.46
2	A	2017	AEE	CAM-CBA-CAY	-2.83	125.46	129.40
2	A	2017	AEE	CAY-CBA-NAW	3.53	125.45	120.62
2	A	2017	AEE	C2-N1-C6	6.00	120.80	116.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2017	AEE	2	0

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	304/328 (92%)	-0.15	8 (2%) 59 33	41, 62, 99, 113	0
1	B	310/328 (94%)	-0.25	4 (1%) 79 59	37, 60, 98, 109	0
All	All	614/656 (93%)	-0.20	12 (1%) 68 44	37, 61, 99, 113	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	874	GLY	3.6
1	A	696	GLY	3.3
1	A	873	GLY	3.2
1	B	873	GLY	3.1
1	B	874	GLY	2.9
1	A	875	LYS	2.7
1	A	992	PRO	2.6
1	A	751	THR	2.6
1	B	696	GLY	2.5
1	B	695	SER	2.4
1	A	991	SER	2.3
1	A	864	ALA	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	AEE	A	2017	33/33	0.84	0.27	1.25	68,73,91,91	0

6.5 Other polymers [i](#)

There are no such residues in this entry.