



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:32 PM GMT

PDB ID : 1XMI
Title : Crystal structure of human F508A NBD1 domain with ATP
Authors : Lewis, H.A.; Zhao, X.; Wang, C.; Sauder, J.M.; Rooney, I.; Noland, B.W.; Lorimer, D.; Kearins, M.C.; Conners, K.; Condon, B.; Maloney, P.C.; Guggino, W.B.; Hunt, J.F.; Emtage, S.; Structural GenomiX
Deposited on : 2004-10-02
Resolution : 2.25 Å(reported)

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

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

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

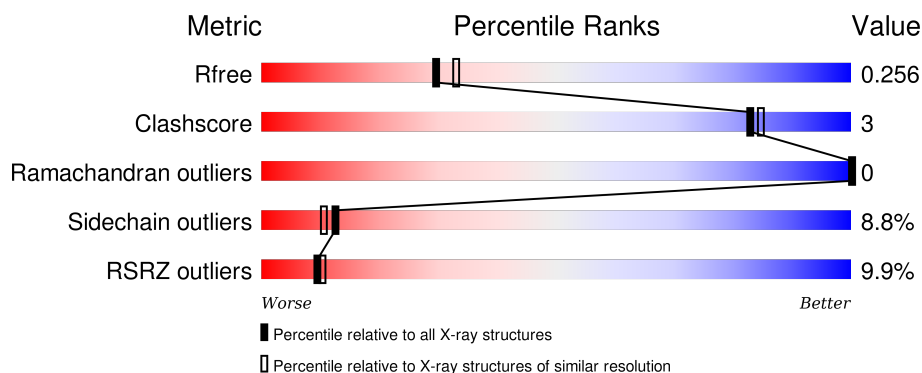
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1640 (2.28-2.24)
Clashscore	102246	1095 (2.26-2.26)
Ramachandran outliers	100387	1063 (2.26-2.26)
Sidechain outliers	100360	1063 (2.26-2.26)
RSRZ outliers	91569	1647 (2.28-2.24)

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	291	<div> <div>9%</div> <div>77%</div> <div>14%</div> <div>8%</div> </div>
1	B	291	<div> <div>4%</div> <div>72%</div> <div>15%</div> <div>11%</div> </div>
1	C	291	<div> <div>8%</div> <div>78%</div> <div>14%</div> <div>7%</div> </div>
1	D	291	<div> <div>13%</div> <div>77%</div> <div>13%</div> <div>8%</div> </div>
1	E	291	<div> <div>11%</div> <div>73%</div> <div>17%</div> <div>9%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11228 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 Cystic fibrosis transmembrane conductance regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2056	1308	342	394	12			
1	B	258	Total	C	N	O	S	0	0	0
			2027	1294	334	386	13			
1	C	270	Total	C	N	O	S	0	0	0
			2117	1347	349	408	13			
1	D	268	Total	C	N	O	S	0	0	0
			2085	1325	344	403	13			
1	E	266	Total	C	N	O	S	0	0	0
			2079	1325	343	398	13			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	388	SER	THR	CLONING ARTIFACT	UNP P13569
A	429	SER	PHE	ENGINEERED	UNP P13569
A	508	ALA	PHE	ENGINEERED	UNP P13569
A	667	ARG	HIS	ENGINEERED	UNP P13569
B	388	SER	THR	CLONING ARTIFACT	UNP P13569
B	429	SER	PHE	ENGINEERED	UNP P13569
B	508	ALA	PHE	ENGINEERED	UNP P13569
B	667	ARG	HIS	ENGINEERED	UNP P13569
C	388	SER	THR	CLONING ARTIFACT	UNP P13569
C	429	SER	PHE	ENGINEERED	UNP P13569
C	508	ALA	PHE	ENGINEERED	UNP P13569
C	667	ARG	HIS	ENGINEERED	UNP P13569
D	388	SER	THR	CLONING ARTIFACT	UNP P13569
D	429	SER	PHE	ENGINEERED	UNP P13569
D	508	ALA	PHE	ENGINEERED	UNP P13569
D	667	ARG	HIS	ENGINEERED	UNP P13569
E	388	SER	THR	CLONING ARTIFACT	UNP P13569
E	429	SER	PHE	ENGINEERED	UNP P13569
E	508	ALA	PHE	ENGINEERED	UNP P13569

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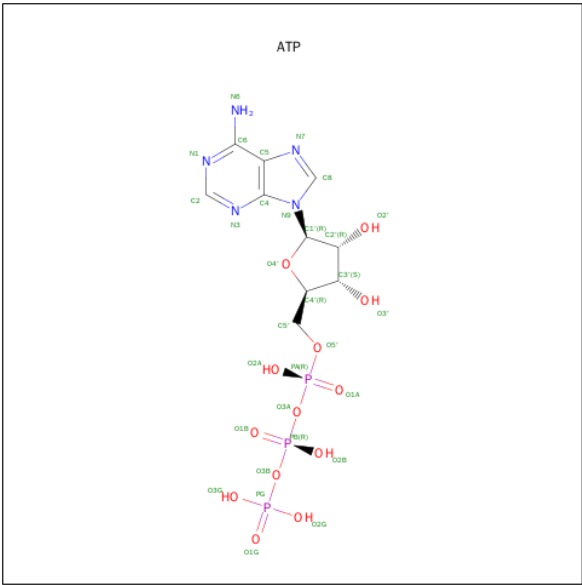
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Chain	Residue	Modelled	Actual	Comment	Reference
E	667	ARG	HIS	ENGINEERED	UNP P13569

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		
2	D	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		
2	E	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	E	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

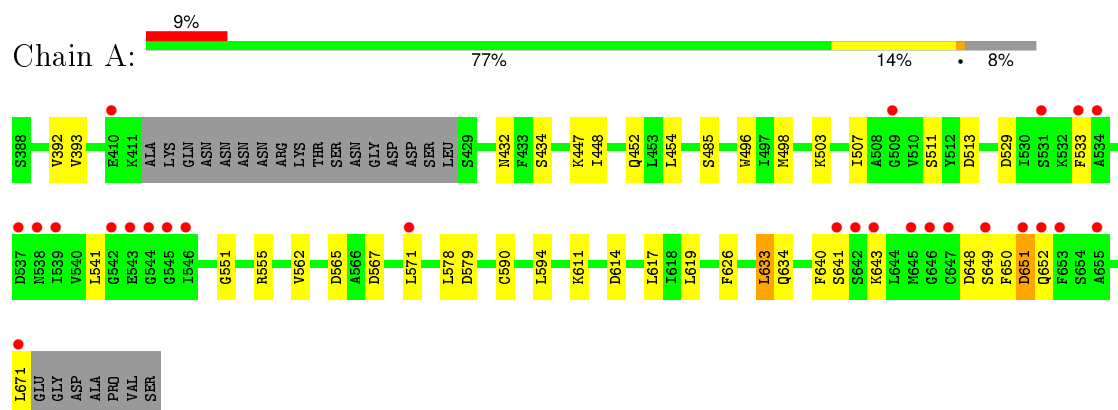
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	155	Total	O	0	0
			155	155		
4	B	163	Total	O	0	0
			163	163		
4	C	146	Total	O	0	0
			146	146		
4	D	132	Total	O	0	0
			132	132		
4	E	108	Total	O	0	0
			108	108		

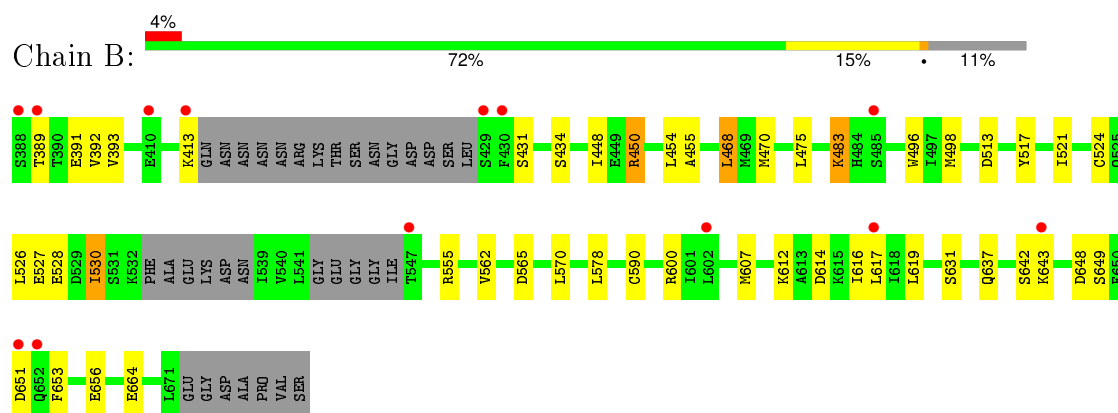
3 Residue-property plots [i](#)

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.

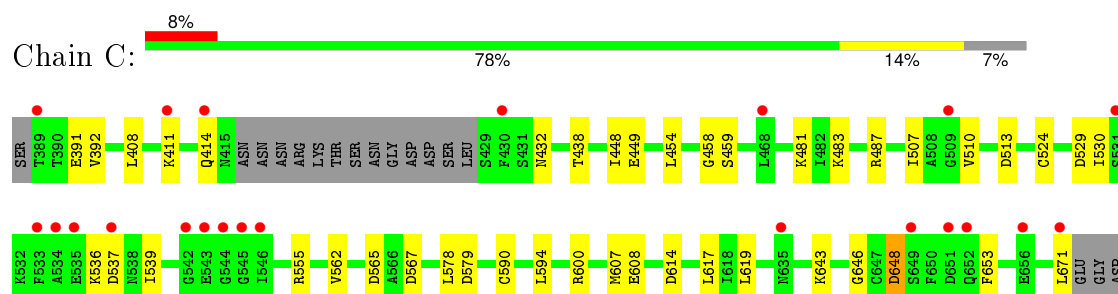
- Molecule 1: Cystic fibrosis transmembrane conductance regulator



- Molecule 1: Cystic fibrosis transmembrane conductance regulator




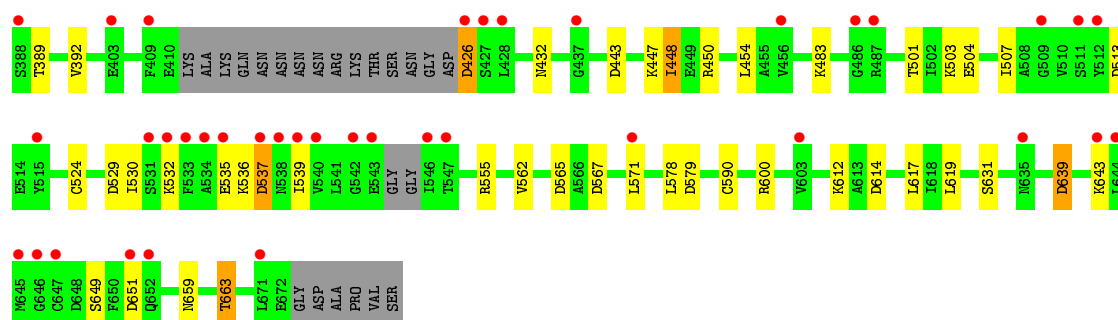
- Molecule 1: Cystic fibrosis transmembrane conductance regulator



ALA
PRO
VAL
SER

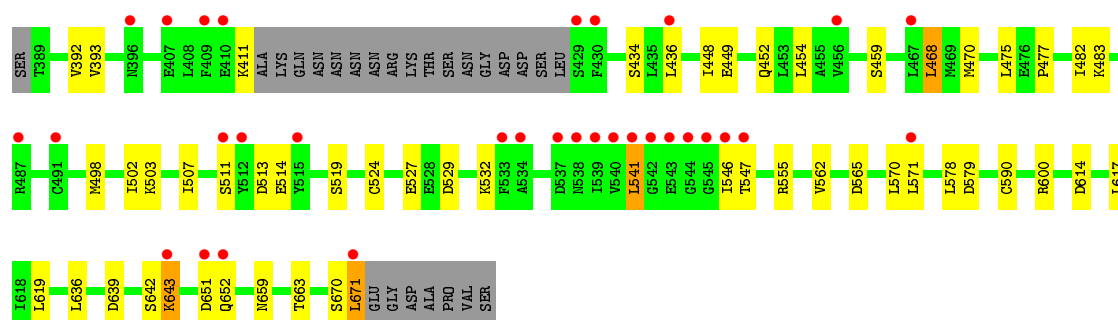
• Molecule 1: Cystic fibrosis transmembrane conductance regulator

Chain D: 



• Molecule 1: Cystic fibrosis transmembrane conductance regulator

Chain E: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	144.46Å 154.02Å 136.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.20 – 2.25 27.20 – 2.25	Depositor EDS
% Data completeness (in resolution range)	(Not available) (27.20-2.25) 99.9 (27.20-2.25)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.13 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.212 , 0.265 0.207 , 0.256	Depositor DCC
R_{free} test set	3582 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	29.7	Xtriage
Anisotropy	0.089	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.4	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 71898 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11228	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.67% 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

5.1 Standard geometry

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

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	0.33	0/2091	0.65	8/2813 (0.3%)
1	B	0.34	0/2059	0.65	4/2762 (0.1%)
1	C	0.34	0/2152	0.65	7/2890 (0.2%)
1	D	0.34	0/2118	0.67	11/2845 (0.4%)
1	E	0.34	0/2114	0.66	7/2839 (0.2%)
All	All	0.34	0/10534	0.66	37/14149 (0.3%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	565	ASP	CB-CG-OD2	6.61	124.25	118.30
1	D	639	ASP	CB-CG-OD2	6.42	124.08	118.30
1	B	651	ASP	CB-CG-OD2	6.13	123.82	118.30
1	D	651	ASP	CB-CG-OD2	5.97	123.67	118.30
1	D	565	ASP	CB-CG-OD2	5.88	123.59	118.30
1	E	513	ASP	CB-CG-OD2	5.86	123.58	118.30
1	C	614	ASP	CB-CG-OD2	5.84	123.56	118.30
1	C	565	ASP	CB-CG-OD2	5.67	123.41	118.30
1	D	426	ASP	CB-CG-OD2	5.50	123.25	118.30
1	C	513	ASP	CB-CG-OD2	5.48	123.23	118.30
1	C	579	ASP	CB-CG-OD2	5.47	123.22	118.30
1	A	579	ASP	CB-CG-OD2	5.44	123.20	118.30
1	E	639	ASP	CB-CG-OD2	5.41	123.17	118.30
1	D	537	ASP	CB-CG-OD2	5.40	123.16	118.30
1	B	614	ASP	CB-CG-OD2	5.36	123.12	118.30
1	D	567	ASP	CB-CG-OD2	5.36	123.12	118.30
1	A	529	ASP	CB-CG-OD2	5.33	123.10	118.30
1	D	513	ASP	CB-CG-OD2	5.32	123.09	118.30
1	A	567	ASP	CB-CG-OD2	5.32	123.08	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	537	ASP	CB-CG-OD2	5.31	123.08	118.30
1	E	614	ASP	CB-CG-OD2	5.29	123.06	118.30
1	A	614	ASP	CB-CG-OD2	5.27	123.05	118.30
1	D	614	ASP	CB-CG-OD2	5.26	123.03	118.30
1	E	579	ASP	CB-CG-OD2	5.26	123.04	118.30
1	B	513	ASP	CB-CG-OD2	5.25	123.03	118.30
1	C	567	ASP	CB-CG-OD2	5.24	123.02	118.30
1	A	513	ASP	CB-CG-OD2	5.21	122.99	118.30
1	E	565	ASP	CB-CG-OD2	5.19	122.97	118.30
1	D	579	ASP	CB-CG-OD2	5.17	122.95	118.30
1	E	651	ASP	CB-CG-OD2	5.14	122.93	118.30
1	A	565	ASP	CB-CG-OD2	5.10	122.89	118.30
1	D	529	ASP	CB-CG-OD2	5.08	122.87	118.30
1	C	529	ASP	CB-CG-OD2	5.08	122.87	118.30
1	E	529	ASP	CB-CG-OD2	5.02	122.82	118.30
1	A	648	ASP	CB-CG-OD2	5.01	122.81	118.30
1	D	443	ASP	CB-CG-OD2	5.00	122.81	118.30
1	A	651	ASP	CB-CG-OD2	5.00	122.80	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	2056	0	2014	12	0
1	B	2027	0	2033	15	0
1	C	2117	0	2106	12	0
1	D	2085	0	2056	12	0
1	E	2079	0	2068	14	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
3	A	31	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	31	0	12	0	0
3	C	31	0	12	0	0
3	D	31	0	12	0	0
3	E	31	0	12	0	0
4	A	155	0	0	0	0
4	B	163	0	0	0	0
4	C	146	0	0	0	0
4	D	132	0	0	1	0
4	E	108	0	0	0	0
All	All	11228	0	10337	64	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 3.

All (64) 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:E:670:SER:O	1:E:671:LEU:HB2	1.85	0.77
1:D:530:ILE:HG12	1:D:537:ASP:HB3	1.77	0.66
1:E:524:CYS:O	1:E:555:ARG:HD2	2.03	0.59
1:B:392:VAL:HB	1:B:448:ILE:HG22	1.83	0.59
1:D:524:CYS:O	1:D:555:ARG:HD2	2.03	0.58
1:E:659:ASN:O	1:E:663:THR:HG23	2.04	0.57
1:D:392:VAL:HB	1:D:448:ILE:HG23	1.87	0.56
1:C:459:SER:HB3	1:C:643:LYS:O	2.05	0.56
1:D:659:ASN:O	1:D:663:THR:CG2	2.55	0.55
1:E:546:ILE:O	1:E:547:THR:HG22	2.06	0.55
1:C:524:CYS:O	1:C:555:ARG:HD2	2.07	0.55
1:E:393:VAL:HB	1:E:483:LYS:HB2	1.89	0.55
1:A:393:VAL:HG22	1:A:447:LYS:HG3	1.90	0.54
1:D:501:THR:HG23	1:D:504:GLU:H	1.71	0.54
1:D:659:ASN:O	1:D:663:THR:HG22	2.08	0.53
1:B:524:CYS:O	1:B:555:ARG:HD2	2.08	0.53
1:D:503:LYS:O	1:D:507:ILE:HG12	2.09	0.53
1:A:448:ILE:HG13	1:A:452:GLN:HB2	1.91	0.52
1:E:470:MET:HB2	1:E:475:LEU:HD23	1.90	0.52
1:B:526:LEU:HB3	1:B:530:ILE:HD12	1.92	0.51
1:B:562:VAL:HG21	1:B:590:CYS:SG	2.51	0.51
1:A:485:SER:HB2	1:B:527:GLU:HG2	1.93	0.50
1:A:562:VAL:HG11	1:A:594:LEU:HD23	1.92	0.50
1:D:389:THR:HG23	1:D:450:ARG:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:455:ALA:HB3	1:B:616:ILE:HG22	1.94	0.48
1:E:477:PRO:HG3	1:E:482:ILE:HD12	1.94	0.48
1:C:562:VAL:HG11	1:C:594:LEU:HD23	1.94	0.48
1:D:562:VAL:HG21	1:D:590:CYS:SG	2.53	0.48
1:C:562:VAL:HG21	1:C:590:CYS:SG	2.54	0.48
1:B:517:TYR:O	1:B:521:ILE:HG12	2.13	0.48
1:A:633:LEU:HD21	1:A:640:PHE:CD2	2.49	0.47
1:A:392:VAL:HB	1:A:448:ILE:HG22	1.95	0.47
1:E:459:SER:HB3	1:E:643:LYS:O	2.15	0.47
1:E:468:LEU:HG	1:E:570:LEU:HB3	1.96	0.47
1:E:562:VAL:HG21	1:E:590:CYS:SG	2.55	0.46
1:B:468:LEU:HG	1:B:570:LEU:HB3	1.97	0.45
1:B:393:VAL:HB	1:B:483:LYS:HB2	1.98	0.45
1:C:392:VAL:HB	1:C:448:ILE:HG22	1.99	0.45
1:C:507:ILE:HG23	1:C:510:VAL:O	2.16	0.45
1:E:502:ILE:HG13	1:E:541:LEU:HD11	1.98	0.45
1:A:626:PHE:CE2	1:A:633:LEU:HB2	2.52	0.45
1:C:530:ILE:O	1:C:536:LYS:HA	2.17	0.45
1:C:648:ASP:HB3	1:C:653:PHE:CE1	2.52	0.44
1:B:555:ARG:NH1	1:B:664:GLU:OE2	2.49	0.44
1:A:562:VAL:HG21	1:A:590:CYS:SG	2.58	0.44
1:C:458:GLY:O	1:C:646:GLY:HA3	2.18	0.43
1:E:503:LYS:O	1:E:507:ILE:HG12	2.18	0.43
1:D:447:LYS:HB3	4:D:777:HOH:O	2.19	0.43
1:C:648:ASP:HB3	1:C:653:PHE:HE1	1.85	0.42
1:B:470:MET:HB2	1:B:475:LEU:HD23	2.00	0.42
1:B:496:TRP:CZ2	1:B:498:MET:HG3	2.55	0.41
1:A:650:PHE:C	1:A:652:GLN:H	2.24	0.41
1:D:392:VAL:HB	1:D:448:ILE:CG2	2.50	0.41
1:E:392:VAL:HB	1:E:448:ILE:HG22	2.02	0.41
1:A:551:GLY:O	1:A:555:ARG:HG3	2.20	0.41
1:B:648:ASP:HB3	1:B:653:PHE:CZ	2.56	0.41
1:C:391:GLU:HG3	1:C:449:GLU:HA	2.02	0.41
1:D:535:GLU:HB2	1:D:539:ILE:HB	2.03	0.41
1:A:503:LYS:O	1:A:507:ILE:HG12	2.21	0.41
1:C:459:SER:HA	1:C:646:GLY:HA2	2.02	0.40
1:B:521:ILE:HD12	1:B:530:ILE:CD1	2.52	0.40
1:B:389:THR:HG23	1:B:450:ARG:HB3	2.02	0.40
1:E:448:ILE:HG13	1:E:452:GLN:HB2	2.04	0.40
1:A:496:TRP:CZ2	1:A:498:MET:HG3	2.57	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	263/291 (90%)	252 (96%)	11 (4%)	0	100	100
1	B	250/291 (86%)	246 (98%)	4 (2%)	0	100	100
1	C	266/291 (91%)	258 (97%)	8 (3%)	0	100	100
1	D	262/291 (90%)	253 (97%)	9 (3%)	0	100	100
1	E	262/291 (90%)	256 (98%)	6 (2%)	0	100	100
All	All	1303/1455 (90%)	1265 (97%)	38 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	220/252 (87%)	202 (92%)	18 (8%)	14	13
1	B	224/252 (89%)	202 (90%)	22 (10%)	10	7
1	C	232/252 (92%)	214 (92%)	18 (8%)	16	14
1	D	227/252 (90%)	209 (92%)	18 (8%)	15	13
1	E	227/252 (90%)	204 (90%)	23 (10%)	9	7
All	All	1130/1260 (90%)	1031 (91%)	99 (9%)	12	10

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	432	ASN
1	A	434	SER
1	A	454	LEU
1	A	511	SER
1	A	533	PHE
1	A	541	LEU
1	A	571	LEU
1	A	578	LEU
1	A	611	LYS
1	A	617	LEU
1	A	619	LEU
1	A	633	LEU
1	A	634	GLN
1	A	641	SER
1	A	643	LYS
1	A	649	SER
1	A	651	ASP
1	A	671	LEU
1	B	391	GLU
1	B	413	LYS
1	B	431	SER
1	B	434	SER
1	B	450	ARG
1	B	454	LEU
1	B	468	LEU
1	B	483	LYS
1	B	528	GLU
1	B	530	ILE
1	B	578	LEU
1	B	600	ARG
1	B	607	MET
1	B	612	LYS
1	B	617	LEU
1	B	619	LEU
1	B	631	SER
1	B	637	GLN
1	B	642	SER
1	B	643	LYS
1	B	649	SER
1	B	656	GLU
1	C	408	LEU
1	C	411	LYS
1	C	414	GLN

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Mol	Chain	Res	Type
1	C	432	ASN
1	C	438	THR
1	C	454	LEU
1	C	481	LYS
1	C	483	LYS
1	C	487	ARG
1	C	539	ILE
1	C	578	LEU
1	C	600	ARG
1	C	607	MET
1	C	608	GLU
1	C	617	LEU
1	C	619	LEU
1	C	648	ASP
1	C	671	LEU
1	D	426	ASP
1	D	432	ASN
1	D	448	ILE
1	D	454	LEU
1	D	483	LYS
1	D	532	LYS
1	D	536	LYS
1	D	571	LEU
1	D	578	LEU
1	D	600	ARG
1	D	612	LYS
1	D	617	LEU
1	D	619	LEU
1	D	631	SER
1	D	639	ASP
1	D	643	LYS
1	D	649	SER
1	D	663	THR
1	E	411	LYS
1	E	434	SER
1	E	436	LEU
1	E	449	GLU
1	E	454	LEU
1	E	468	LEU
1	E	498	MET
1	E	511	SER
1	E	514	GLU

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Mol	Chain	Res	Type
1	E	519	SER
1	E	527	GLU
1	E	532	LYS
1	E	541	LEU
1	E	571	LEU
1	E	578	LEU
1	E	600	ARG
1	E	617	LEU
1	E	619	LEU
1	E	636	LEU
1	E	642	SER
1	E	643	LYS
1	E	652	GLN
1	E	671	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	432	ASN
1	A	484	HIS
1	A	637	GLN
1	B	637	GLN
1	C	432	ASN
1	C	484	HIS
1	C	637	GLN
1	D	432	ASN
1	D	538	ASN
1	D	637	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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$
3	ATP	A	2	2	24,33,33	1.10	2 (8%)	31,52,52	2.37	4 (12%)
3	ATP	B	1	2	24,33,33	1.12	2 (8%)	31,52,52	2.39	3 (9%)
3	ATP	C	3	2	24,33,33	1.12	2 (8%)	31,52,52	2.47	3 (9%)
3	ATP	D	4	2	24,33,33	1.16	2 (8%)	31,52,52	2.46	4 (12%)
3	ATP	E	5	2	24,33,33	1.11	2 (8%)	31,52,52	2.52	3 (9%)

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
3	ATP	A	2	2	-	0/18/38/38	0/3/3/3
3	ATP	B	1	2	-	0/18/38/38	0/3/3/3
3	ATP	C	3	2	-	0/18/38/38	0/3/3/3
3	ATP	D	4	2	-	0/18/38/38	0/3/3/3
3	ATP	E	5	2	-	0/18/38/38	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2	ATP	C2-N1	2.41	1.38	1.33
3	E	5	ATP	C2-N1	2.53	1.38	1.33
3	C	3	ATP	C2-N1	2.55	1.38	1.33
3	B	1	ATP	C2-N1	2.55	1.38	1.33
3	D	4	ATP	C2-N1	2.71	1.39	1.33
3	B	1	ATP	C2-N3	3.69	1.38	1.32
3	A	2	ATP	C2-N3	3.73	1.38	1.32
3	E	5	ATP	C2-N3	3.74	1.38	1.32
3	C	3	ATP	C2-N3	3.75	1.38	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	4	ATP	C2-N3	3.93	1.39	1.32

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	5	ATP	N3-C2-N1	-12.47	119.35	128.89
3	C	3	ATP	N3-C2-N1	-12.37	119.42	128.89
3	D	4	ATP	N3-C2-N1	-12.09	119.64	128.89
3	B	1	ATP	N3-C2-N1	-11.95	119.75	128.89
3	A	2	ATP	N3-C2-N1	-11.86	119.81	128.89
3	D	4	ATP	PA-O3A-PB	-3.22	123.68	132.73
3	E	5	ATP	PA-O3A-PB	-2.69	125.16	132.73
3	C	3	ATP	PA-O3A-PB	-2.69	125.18	132.73
3	B	1	ATP	PB-O3B-PG	-2.65	123.78	132.67
3	B	1	ATP	PA-O3A-PB	-2.53	125.62	132.73
3	E	5	ATP	PB-O3B-PG	-2.49	124.33	132.67
3	D	4	ATP	PB-O3B-PG	-2.44	124.49	132.67
3	C	3	ATP	PB-O3B-PG	-2.36	124.76	132.67
3	A	2	ATP	PA-O3A-PB	-2.21	126.52	132.73
3	A	2	ATP	PB-O3B-PG	-2.15	125.46	132.67
3	A	2	ATP	C4-C5-N7	-2.02	107.62	109.48
3	D	4	ATP	O2B-PB-O3A	2.00	114.17	105.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	267/291 (91%)	0.50	26 (9%) 10 11	10, 28, 62, 73	0
1	B	258/291 (88%)	0.31	13 (5%) 32 36	12, 27, 47, 76	0
1	C	270/291 (92%)	0.41	22 (8%) 15 16	9, 26, 56, 72	0
1	D	268/291 (92%)	0.67	38 (14%) 4 4	11, 33, 59, 69	0
1	E	266/291 (91%)	0.60	32 (12%) 6 6	14, 32, 58, 68	0
All	All	1329/1455 (91%)	0.50	131 (9%) 9 10	9, 29, 59, 76	0

All (131) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	545	GLY	11.9
1	A	544	GLY	8.8
1	E	546	ILE	7.7
1	D	426	ASP	6.6
1	C	546	ILE	6.4
1	A	546	ILE	6.4
1	E	542	GLY	6.4
1	A	533	PHE	6.0
1	A	534	ALA	5.9
1	C	544	GLY	5.7
1	A	647	CYS	5.6
1	D	509	GLY	5.6
1	A	531	SER	5.5
1	C	649	SER	5.4
1	B	430	PHE	5.2
1	E	544	GLY	5.0
1	D	647	CYS	5.0
1	D	645	MET	4.8
1	A	646	GLY	4.7
1	A	543	GLU	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	547	THR	4.5
1	B	429	SER	4.4
1	E	533	PHE	4.4
1	E	511	SER	4.4
1	A	671	LEU	4.3
1	A	542	GLY	4.3
1	D	388	SER	4.1
1	E	539	ILE	4.0
1	D	646	GLY	4.0
1	C	534	ALA	4.0
1	A	651	ASP	3.8
1	A	641	SER	3.7
1	E	429	SER	3.7
1	A	539	ILE	3.5
1	D	543	GLU	3.5
1	E	651	ASP	3.5
1	C	543	GLU	3.5
1	E	538	ASN	3.4
1	C	652	GLN	3.4
1	D	540	VAL	3.4
1	E	515	TYR	3.4
1	B	388	SER	3.4
1	B	413	LYS	3.4
1	E	652	GLN	3.3
1	B	652	GLN	3.3
1	C	430	PHE	3.2
1	C	535	GLU	3.2
1	C	651	ASP	3.2
1	D	428	LEU	3.2
1	E	436	LEU	3.2
1	E	671	LEU	3.2
1	A	642	SER	3.2
1	E	409	PHE	3.0
1	D	537	ASP	2.9
1	D	603	VAL	2.9
1	C	531	SER	2.9
1	D	547	THR	2.9
1	D	546	ILE	2.9
1	D	534	ALA	2.8
1	A	410	GLU	2.8
1	C	533	PHE	2.8
1	D	515	TYR	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	410	GLU	2.8
1	D	456	VAL	2.8
1	E	643	LYS	2.8
1	A	545	GLY	2.8
1	A	643	LYS	2.8
1	C	411	LYS	2.8
1	C	509	GLY	2.8
1	D	651	ASP	2.7
1	C	414	GLN	2.7
1	E	430	PHE	2.7
1	D	635	ASN	2.7
1	E	407	GLU	2.7
1	E	543	GLU	2.7
1	D	427	SER	2.7
1	A	652	GLN	2.7
1	E	547	THR	2.7
1	E	540	VAL	2.7
1	D	644	LEU	2.6
1	D	542	GLY	2.6
1	D	643	LYS	2.6
1	E	534	ALA	2.6
1	E	512	TYR	2.6
1	D	533	PHE	2.6
1	D	532	LYS	2.6
1	E	545	GLY	2.5
1	B	651	ASP	2.5
1	D	512	TYR	2.5
1	B	643	LYS	2.5
1	D	487	ARG	2.5
1	D	437	GLY	2.4
1	E	571	LEU	2.4
1	E	491	CYS	2.4
1	A	645	MET	2.4
1	C	537	ASP	2.4
1	B	485	SER	2.4
1	D	535	GLU	2.4
1	A	537	ASP	2.3
1	A	649	SER	2.3
1	D	531	SER	2.3
1	E	541	LEU	2.3
1	D	409	PHE	2.3
1	D	571	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	511	SER	2.2
1	C	389	THR	2.2
1	A	538	ASN	2.2
1	E	456	VAL	2.2
1	B	602	LEU	2.2
1	D	539	ILE	2.2
1	A	653	PHE	2.2
1	D	538	ASN	2.2
1	D	671	LEU	2.2
1	A	509	GLY	2.1
1	C	635	ASN	2.1
1	E	487	ARG	2.1
1	A	571	LEU	2.1
1	C	542	GLY	2.1
1	C	656	GLU	2.1
1	D	652	GLN	2.1
1	E	467	LEU	2.1
1	D	486	GLY	2.1
1	D	403	GLU	2.1
1	E	410	GLU	2.1
1	B	389	THR	2.0
1	E	396	ASN	2.0
1	C	468	LEU	2.0
1	C	671	LEU	2.0
1	A	655	ALA	2.0
1	E	537	ASP	2.0
1	B	617	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	ATP	D	4	31/31	0.98	0.10	-0.67	8,17,24,29	0
3	ATP	C	3	31/31	0.98	0.10	-0.70	9,21,25,29	0
3	ATP	B	1	31/31	0.98	0.10	-0.76	13,20,26,33	0
3	ATP	E	5	31/31	0.97	0.10	-0.78	17,23,30,32	0
3	ATP	A	2	31/31	0.98	0.10	-0.82	12,17,25,26	0
2	MG	E	10	1/1	0.84	0.10	-	20,20,20,20	0
2	MG	D	9	1/1	0.93	0.08	-	13,13,13,13	0
2	MG	B	6	1/1	0.95	0.06	-	18,18,18,18	0
2	MG	C	8	1/1	0.92	0.10	-	16,16,16,16	0
2	MG	A	7	1/1	0.91	0.07	-	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.