



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2024 – 12:48 PM EDT

PDB ID : 7KEO
Title : Crystal structure of K29-linked di-ubiquitin in complex with synthetic antigen binding fragment
Authors : Yu, Y.; Zheng, Q.; Erramilli, S.; Pan, M.; Kossiakoff, A.; Liu, L.; Zhao, M.
Deposited on : 2020-10-11
Resolution : 2.90 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36.1
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) : 2.36.1

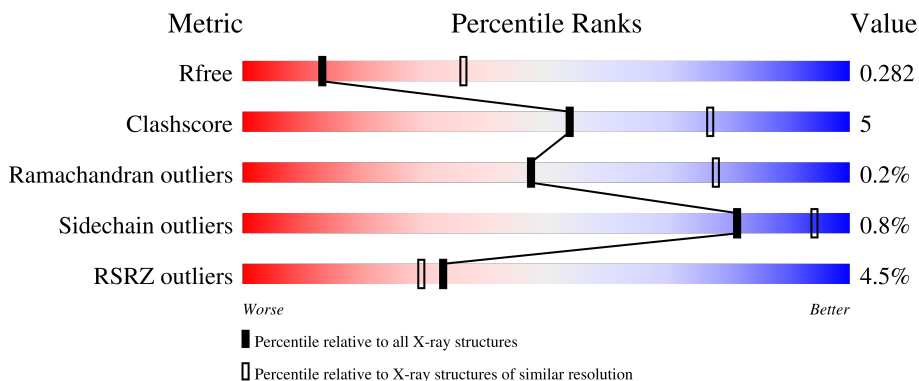
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

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






Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	235	
1	E	235	
2	B	215	
2	F	215	
3	C	76	

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Mol	Chain	Length	Quality of chain
3	D	76	 88% 12%
3	G	76	 5% 84% 11% 5%
3	H	76	 83% 17%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	PO4	A	301	-	-	-	X
4	PO4	A	302	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8889 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 Synthetic antigen binding fragment, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	227	Total	C	N	O	S	0	0	0
			1685	1068	278	332	7			
1	E	225	Total	C	N	O	S	0	0	0
			1667	1058	275	328	6			

- Molecule 2 is a protein called Synthetic antigen binding fragment, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	214	Total	C	N	O	S	1	0	0
			1612	1008	268	331	5			
2	F	214	Total	C	N	O	S	1	0	0
			1614	1009	271	328	6			

- Molecule 3 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	73	Total	C	N	O	S	0	0	0
			558	356	92	109	1			
3	D	76	Total	C	N	O	S	0	0	0
			577	364	103	109	1			
3	G	72	Total	C	N	O	S	0	0	0
			555	351	91	112	1			
3	H	76	Total	C	N	O	S	0	0	0
			578	364	102	111	1			

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 5 4 1	0	0
4	A	1	Total O P 5 4 1	0	0
4	B	1	Total O P 5 4 1	0	0
4	D	1	Total O P 5 4 1	0	0
4	E	1	Total O P 5 4 1	0	0
4	H	1	Total O P 5 4 1	0	0

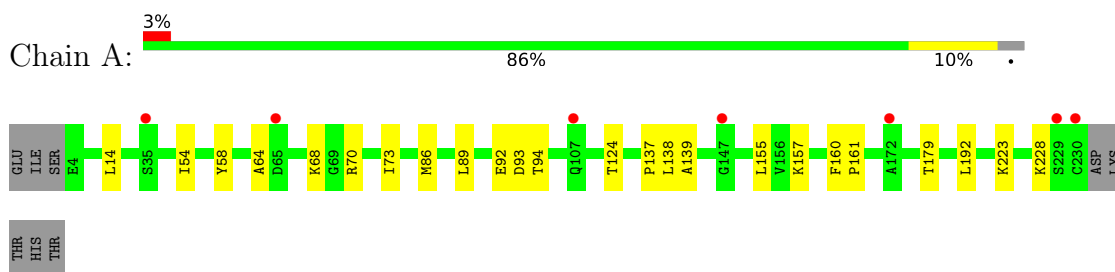
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	3	Total O 3 3	0	0
5	B	4	Total O 4 4	0	0
5	D	1	Total O 1 1	0	0
5	E	3	Total O 3 3	0	0
5	G	1	Total O 1 1	0	0
5	H	1	Total O 1 1	0	0

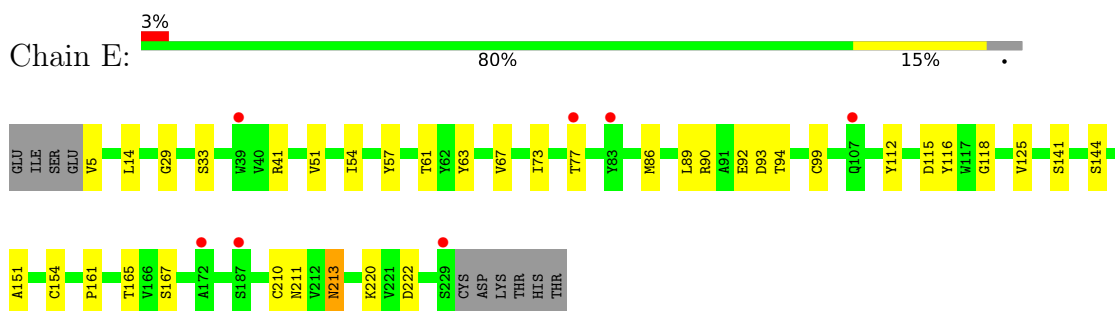
3 Residue-property plots

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.

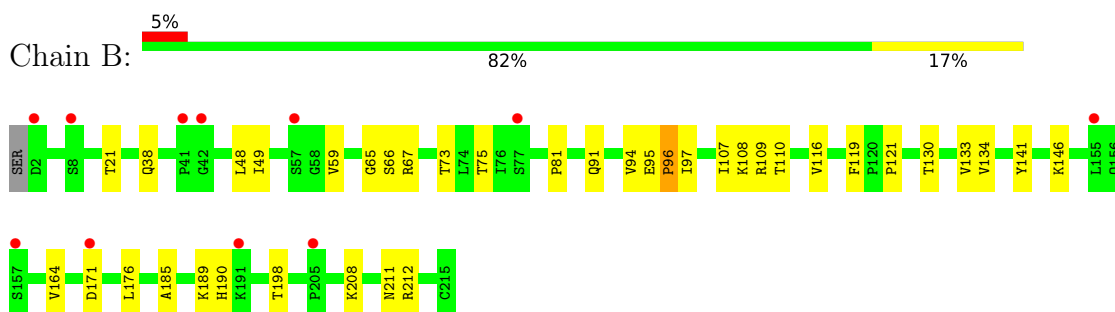
- Molecule 1: Synthetic antigen binding fragment, heavy chain



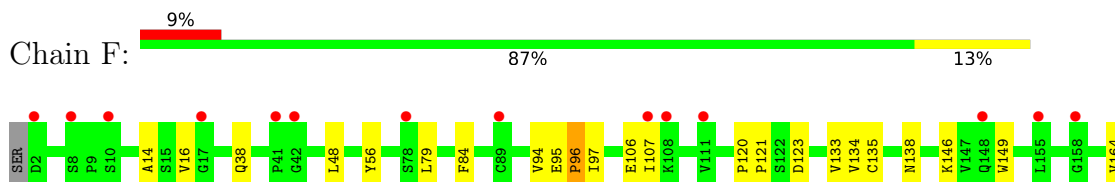
- Molecule 1: Synthetic antigen binding fragment, heavy chain

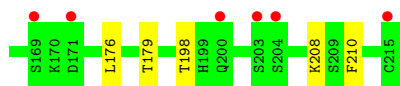


- Molecule 2: Synthetic antigen binding fragment, light chain

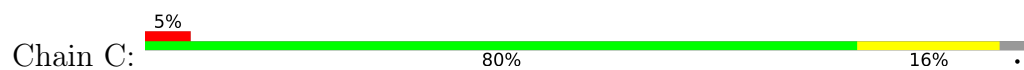


- Molecule 2: Synthetic antigen binding fragment, light chain

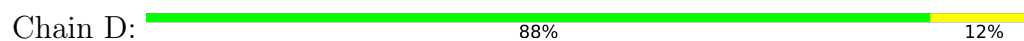




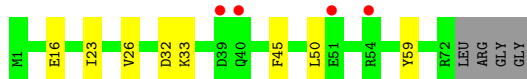
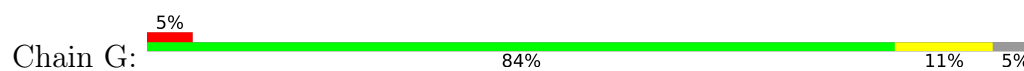
- Molecule 3: Ubiquitin



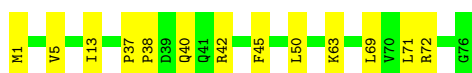
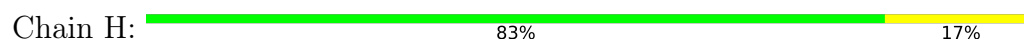
- Molecule 3: Ubiquitin



- Molecule 3: Ubiquitin



- Molecule 3: Ubiquitin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.62Å 43.74Å 174.75Å 90.00° 107.81° 90.00°	Depositor
Resolution (Å)	19.70 – 2.90 19.70 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.3 (19.70-2.90) 98.5 (19.70-2.90)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 2.88Å)	Xtrriage
Refinement program	PHENIX 1.14	Depositor
R, R_{free}	0.236 , 0.282 0.236 , 0.282	Depositor DCC
R_{free} test set	1502 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	64.2	Xtrriage
Anisotropy	0.119	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8889	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.86 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.6853e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PO4

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	0.25	0/1730	0.45	0/2363
1	E	0.25	0/1712	0.44	0/2340
2	B	0.25	0/1646	0.45	0/2239
2	F	0.25	0/1648	0.44	0/2242
3	C	0.24	0/564	0.42	0/763
3	D	0.23	0/583	0.44	0/788
3	G	0.25	0/561	0.42	0/758
3	H	0.23	0/584	0.44	0/790
All	All	0.25	0/9028	0.44	0/12283

There are no bond length outliers.

There are no bond angle outliers.

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	1685	0	1618	14	0
1	E	1667	0	1598	22	0
2	B	1612	0	1547	20	0
2	F	1614	0	1557	15	0
3	C	558	0	572	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	577	0	591	6	0
3	G	555	0	559	7	0
3	H	578	0	586	10	0
4	A	10	0	0	0	0
4	B	5	0	0	0	0
4	D	5	0	0	0	0
4	E	5	0	0	0	0
4	H	5	0	0	0	0
5	A	3	0	0	0	0
5	B	4	0	0	0	0
5	D	1	0	0	0	0
5	E	3	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
All	All	8889	0	8628	92	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 5.

All (92) 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:94:THR:HG22	1:E:125:VAL:H	1.51	0.74
3:C:23:ILE:HD13	3:C:50:LEU:HB3	1.71	0.72
3:G:23:ILE:HD13	3:G:50:LEU:HB3	1.71	0.72
3:D:13:ILE:HG22	1:E:220:LYS:HB2	1.73	0.70
2:B:121:PRO:HD3	2:B:133:VAL:HG22	1.74	0.70
1:A:70:ARG:NH2	1:A:93:ASP:OD2	2.24	0.69
2:F:121:PRO:HD3	2:F:133:VAL:HG22	1.75	0.68
2:F:16:VAL:HG22	2:F:107:ILE:HD11	1.78	0.65
3:D:7:THR:HG23	3:D:9:THR:H	1.64	0.63
3:C:11:LYS:HE3	3:D:48:LYS:HE3	1.83	0.61
3:H:5:VAL:HB	3:H:13:ILE:HG23	1.84	0.60
2:B:38:GLN:HB2	2:B:48:LEU:HD11	1.83	0.59
2:B:94:VAL:HG12	2:B:96:PRO:HD2	1.84	0.59
2:F:164:VAL:HG22	2:F:176:LEU:HD13	1.85	0.58
2:B:164:VAL:HG22	2:B:176:LEU:HD13	1.86	0.57
1:E:167:SER:HB2	1:E:211:ASN:HB2	1.85	0.57
1:E:165:THR:HB	1:E:213:ASN:HB3	1.88	0.55
1:E:115:ASP:OD1	1:E:115:ASP:N	2.37	0.55
2:B:146:LYS:HB3	2:B:198:THR:HB	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:14:LEU:HB2	1:E:161:PRO:HG3	1.87	0.55
1:E:90:ARG:NH1	1:E:92:GLU:OE1	2.40	0.54
1:E:57:TYR:OH	3:H:69:LEU:O	2.24	0.54
1:E:41:ARG:NH2	1:E:93:ASP:OD1	2.35	0.53
3:G:23:ILE:HA	3:G:26:VAL:HG12	1.90	0.53
2:F:38:GLN:HB2	2:F:48:LEU:HD11	1.91	0.53
1:A:137:PRO:HD3	1:A:223:LYS:HE3	1.91	0.52
2:B:109:ARG:NH2	2:B:171:ASP:O	2.44	0.51
3:C:23:ILE:HA	3:C:26:VAL:HG12	1.92	0.50
1:E:141:SER:H	1:E:144:SER:HB3	1.76	0.50
2:B:109:ARG:HG2	2:B:110:THR:N	2.27	0.50
2:F:94:VAL:HG12	2:F:96:PRO:HD2	1.94	0.50
1:E:57:TYR:CZ	3:H:71:LEU:HD23	2.47	0.50
1:A:92:GLU:OE1	1:A:92:GLU:N	2.43	0.49
3:H:45:PHE:HB3	3:H:50:LEU:HD21	1.94	0.49
2:B:48:LEU:HA	2:B:59:VAL:HG21	1.95	0.49
2:B:95:GLU:O	2:B:97:ILE:N	2.46	0.49
2:B:190:HIS:O	2:B:212:ARG:NH1	2.41	0.49
1:E:86:MET:HB3	1:E:89:LEU:HD21	1.94	0.48
1:A:58:TYR:OH	3:D:71:LEU:O	2.23	0.48
2:B:66:SER:OG	2:B:73:THR:OG1	2.26	0.48
1:E:51:VAL:HG13	1:E:67:VAL:HG21	1.96	0.48
1:E:63:TYR:CZ	1:E:73:ILE:HG22	2.48	0.48
1:A:86:MET:HB3	1:A:89:LEU:HD21	1.95	0.47
3:C:5:VAL:HG22	3:C:67:LEU:HB2	1.96	0.47
2:B:49:ILE:HG21	2:B:65:GLY:HA3	1.97	0.47
1:E:5:VAL:HG22	1:E:29:GLY:HA3	1.97	0.47
2:F:146:LYS:HB3	2:F:198:THR:HB	1.96	0.47
1:E:112:TYR:OH	3:H:40:GLN:HG2	2.15	0.47
3:G:33:LYS:HB3	3:H:72:ARG:HD3	1.96	0.46
2:B:81:PRO:HA	2:B:107:ILE:HG13	1.97	0.46
3:D:3:ILE:HD12	3:D:67:LEU:HD13	1.98	0.46
3:G:45:PHE:HB3	3:G:50:LEU:HD21	1.97	0.46
2:B:67:ARG:NH2	3:C:21:ASP:OD1	2.49	0.45
3:D:23:ILE:HA	3:D:26:VAL:HG12	1.99	0.45
1:A:54:ILE:HG13	1:A:73:ILE:HD13	1.98	0.45
1:E:99:CYS:O	1:E:118:GLY:N	2.47	0.45
3:H:1:MET:HB2	3:H:63:LYS:HG2	1.99	0.44
2:B:119:PHE:HB2	2:B:134:VAL:HG13	2.00	0.43
2:F:84:PHE:HD1	2:F:106:GLU:HA	1.83	0.43
1:A:14:LEU:HB2	1:A:161:PRO:HG3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:95:GLU:O	2:F:97:ILE:N	2.51	0.43
1:E:210:CYS:O	1:E:222:ASP:HA	2.17	0.43
1:A:139:ALA:HB3	1:A:228:LYS:HG2	1.99	0.43
2:B:116:VAL:HG13	2:B:208:LYS:HE3	2.00	0.43
1:A:64:ALA:O	1:A:68:LYS:HG3	2.19	0.43
1:A:138:LEU:HB3	2:B:119:PHE:CD2	2.53	0.43
1:A:94:THR:HG23	1:A:124:THR:HA	2.01	0.42
3:G:32:ASP:OD2	3:H:72:ARG:HD2	2.19	0.42
3:G:50:LEU:HD22	3:G:59:TYR:CD2	2.53	0.42
3:C:23:ILE:HD12	3:C:51:GLU:O	2.19	0.42
2:B:185:ALA:O	2:B:189:LYS:HG3	2.19	0.42
2:F:123:ASP:OD1	2:F:123:ASP:N	2.50	0.42
2:F:208:LYS:HB3	2:F:208:LYS:HE2	1.80	0.42
1:A:160:PHE:HA	1:A:161:PRO:HA	1.84	0.42
1:E:5:VAL:HB	1:E:116:TYR:HE2	1.83	0.42
3:C:50:LEU:HD22	3:C:59:TYR:CD2	2.55	0.42
2:F:14:ALA:HB3	2:F:79:LEU:HD21	2.03	0.41
3:G:33:LYS:HA	3:H:42:ARG:HD2	2.02	0.41
1:A:179:THR:HG23	1:A:192:LEU:HD21	2.02	0.41
1:A:155:LEU:HD13	1:A:157:LYS:HB2	2.02	0.41
2:F:120:PRO:HB3	2:F:210:PHE:CE2	2.55	0.41
1:E:33:SER:HB2	1:E:77:THR:HG23	2.02	0.41
2:F:134:VAL:HA	2:F:179:THR:HA	2.03	0.41
3:H:37:PRO:HA	3:H:38:PRO:HD3	1.96	0.41
2:F:135:CYS:HB2	2:F:149:TRP:CZ2	2.56	0.40
2:F:134:VAL:HB	2:F:179:THR:HG22	2.02	0.40
2:B:108:LYS:HA	2:B:141:TYR:OH	2.22	0.40
1:E:144:SER:OG	1:E:151:ALA:O	2.39	0.40
2:B:21:THR:HG22	2:B:75:THR:HG22	2.03	0.40
3:C:56:LEU:HD22	3:C:61:ILE:HD12	2.03	0.40
3:C:36:ILE:H	3:C:36:ILE:HG12	1.68	0.40
1:E:54:ILE:HD13	1:E:61:THR:HG22	2.04	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	225/235 (96%)	212 (94%)	13 (6%)	0	100	100
1	E	223/235 (95%)	216 (97%)	7 (3%)	0	100	100
2	B	212/215 (99%)	200 (94%)	11 (5%)	1 (0%)	29	61
2	F	212/215 (99%)	196 (92%)	15 (7%)	1 (0%)	29	61
3	C	71/76 (93%)	68 (96%)	3 (4%)	0	100	100
3	D	74/76 (97%)	73 (99%)	1 (1%)	0	100	100
3	G	70/76 (92%)	68 (97%)	2 (3%)	0	100	100
3	H	74/76 (97%)	72 (97%)	2 (3%)	0	100	100
All	All	1161/1204 (96%)	1105 (95%)	54 (5%)	2 (0%)	47	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	96	PRO
2	F	96	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	184/195 (94%)	184 (100%)	0	100	100
1	E	181/195 (93%)	179 (99%)	2 (1%)	73	92
2	B	182/189 (96%)	179 (98%)	3 (2%)	62	86
2	F	183/189 (97%)	181 (99%)	2 (1%)	73	92
3	C	62/68 (91%)	62 (100%)	0	100	100
3	D	62/68 (91%)	62 (100%)	0	100	100
3	G	62/68 (91%)	61 (98%)	1 (2%)	62	86

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	H	62/68 (91%)	62 (100%)	0	100	100
All	All	978/1040 (94%)	970 (99%)	8 (1%)	81	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	91	GLN
2	B	130	THR
2	B	211	ASN
1	E	154	CYS
1	E	213	ASN
2	F	56	TYR
2	F	138	ASN
3	G	16	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

6 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	$\# Z > 2$
4	PO4	E	301	-	4,4,4	0.93	0	6,6,6	0.43	0
4	PO4	A	301	-	4,4,4	0.93	0	6,6,6	0.44	0
4	PO4	B	301	-	4,4,4	0.94	0	6,6,6	0.48	0
4	PO4	A	302	-	4,4,4	0.92	0	6,6,6	0.44	0
4	PO4	D	101	-	4,4,4	0.91	0	6,6,6	0.45	0
4	PO4	H	101	-	4,4,4	0.94	0	6,6,6	0.40	0

There are no bond length outliers.

There are no bond angle outliers.

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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	227/235 (96%)	0.18	7 (3%) 49 44	40, 53, 65, 95	1 (0%)
1	E	225/235 (95%)	0.28	7 (3%) 49 44	46, 59, 72, 83	1 (0%)
2	B	214/215 (99%)	0.36	11 (5%) 28 24	45, 59, 74, 88	1 (0%)
2	F	214/215 (99%)	0.53	20 (9%) 8 6	50, 68, 89, 98	1 (0%)
3	C	73/76 (96%)	0.57	4 (5%) 25 21	54, 76, 90, 99	0
3	D	76/76 (100%)	0.02	0 100 100	45, 58, 71, 84	0
3	G	72/76 (94%)	0.51	4 (5%) 24 20	59, 73, 86, 94	0
3	H	76/76 (100%)	-0.08	0 100 100	48, 58, 73, 87	0
All	All	1177/1204 (97%)	0.31	53 (4%) 33 29	40, 60, 84, 99	4 (0%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	73	LEU	7.4
2	F	158	GLY	7.3
1	A	230	CYS	6.0
1	A	147	GLY	5.4
2	B	157	SER	5.3
1	A	229	SER	5.2
2	B	42	GLY	4.6
2	F	204	SER	4.3
2	B	155	LEU	4.1
1	E	77	THR	3.6
3	C	72	ARG	3.4
1	A	107	GLN	3.4
2	F	215	CYS	3.2
2	F	41	PRO	3.1
2	F	42	GLY	3.0
2	F	17	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
2	F	169	SER	2.9
2	B	41	PRO	2.9
2	B	171	ASP	2.9
2	F	2	ASP	2.8
2	F	148	GLN	2.8
2	B	57	SER	2.7
3	G	39	ASP	2.7
2	F	89	CYS	2.7
1	A	172	ALA	2.7
1	E	229	SER	2.6
2	F	200	GLN	2.6
2	F	171	ASP	2.6
2	F	10	SER	2.6
2	B	77	SER	2.5
3	G	40	GLN	2.5
3	G	51	GLU	2.5
1	E	107	GLN	2.5
2	B	8	SER	2.4
1	A	35	SER	2.4
2	F	203	SER	2.3
1	E	187	SER	2.3
1	E	39	TRP	2.3
2	B	205	PRO	2.3
2	F	8	SER	2.2
2	F	111	VAL	2.2
3	G	54	ARG	2.2
2	B	191	LYS	2.2
1	A	65	ASP	2.2
2	F	155	LEU	2.2
2	F	78	SER	2.2
1	E	83	TYR	2.2
3	C	35	GLY	2.1
3	C	53	GLY	2.1
2	F	108	LYS	2.1
2	B	2	ASP	2.1
2	F	107	ILE	2.0
1	E	172	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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
4	PO4	A	301	5/5	0.68	0.43	67,70,87,91	0
4	PO4	A	302	5/5	0.78	0.52	50,52,59,62	5
4	PO4	E	301	5/5	0.89	0.28	72,77,83,104	0
4	PO4	B	301	5/5	0.90	0.27	69,79,88,99	0
4	PO4	D	101	5/5	0.95	0.14	48,53,58,65	0
4	PO4	H	101	5/5	0.95	0.15	60,64,73,73	0

6.5 Other polymers [i](#)

There are no such residues in this entry.