



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 01:33 am GMT

PDB ID : 4X6C  
Title : CD1a ternary complex with lysophosphatidylcholine and BK6 TCR  
Authors : Birkinshaw, R.W.; Rossjohn, J.  
Deposited on : 2014-12-08  
Resolution : 2.80 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](mailto: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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

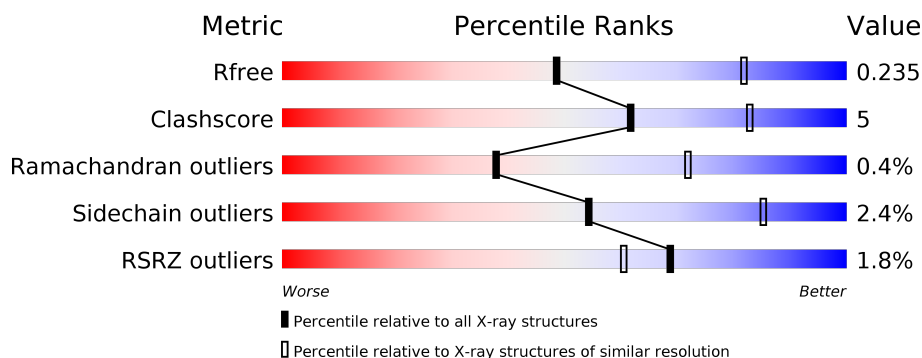
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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.



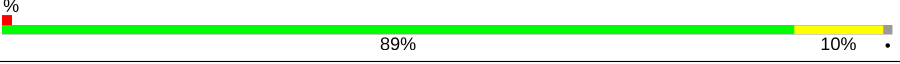

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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  $\geq 3$ , 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	281	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>13%</div> <div>6%</div> </div> </div>
1	C	281	<div> <div>4%</div> <div> <div></div> <div>70%</div> <div>16%</div> <div>11%</div> </div> </div>
2	B	105	<div> <div>%</div> <div> <div></div> <div>72%</div> <div>20%</div> <div>7%</div> </div> </div>
2	D	105	<div> <div>3%</div> <div> <div></div> <div>69%</div> <div>16%</div> <div>14%</div> </div> </div>
3	E	207	<div> <div></div> <div> <div></div> <div>86%</div> <div>10%</div> <div>.</div> </div> </div>
3	G	207	<div> <div></div> <div> <div></div> <div>84%</div> <div>13%</div> <div>.</div> </div> </div>

Continued on next page...

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
4	F	245	 89% 10%
4	H	245	 91% 8%

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
5	NAG	A	501	-	-	-	X
6	42H	A	503	-	-	-	X
6	42H	C	302	-	-	-	X

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 12865 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 T-cell surface glycoprotein CD1a.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2122	1359	370	385	8			
1	C	249	Total	C	N	O	S	0	0	0
			2054	1324	357	365	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	ILE	THR	variant	UNP P06126
A	51	TRP	CYS	variant	UNP P06126
A	279	GLY	-	expression tag	UNP P06126
A	280	SER	-	expression tag	UNP P06126
A	281	LEU	-	expression tag	UNP P06126
A	282	VAL	-	expression tag	UNP P06126
A	283	PRO	-	expression tag	UNP P06126
A	284	ARG	-	expression tag	UNP P06126
C	13	ILE	THR	variant	UNP P06126
C	51	TRP	CYS	variant	UNP P06126
C	279	GLY	-	expression tag	UNP P06126
C	280	SER	-	expression tag	UNP P06126
C	281	LEU	-	expression tag	UNP P06126
C	282	VAL	-	expression tag	UNP P06126
C	283	PRO	-	expression tag	UNP P06126
C	284	ARG	-	expression tag	UNP P06126

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	0	0
			820	523	139	156	2			
2	D	90	Total	C	N	O	S	0	0	0
			762	490	128	142	2			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	100	GLY	-	expression tag	UNP P61769
B	101	SER	-	expression tag	UNP P61769
B	102	LEU	-	expression tag	UNP P61769
B	103	VAL	-	expression tag	UNP P61769
B	104	PRO	-	expression tag	UNP P61769
B	105	ARG	-	expression tag	UNP P61769
D	100	GLY	-	expression tag	UNP P61769
D	101	SER	-	expression tag	UNP P61769
D	102	LEU	-	expression tag	UNP P61769
D	103	VAL	-	expression tag	UNP P61769
D	104	PRO	-	expression tag	UNP P61769
D	105	ARG	-	expression tag	UNP P61769

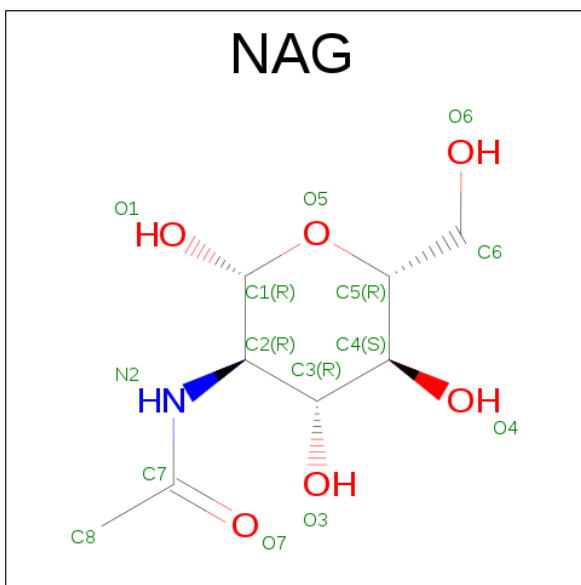
- Molecule 3 is a protein called TCR alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	198	Total	C	N	O	S	0	0	0
			1548	969	251	318	10			
3	G	201	Total	C	N	O	S	0	0	0
			1569	981	255	323	10			

- Molecule 4 is a protein called TCR beta.

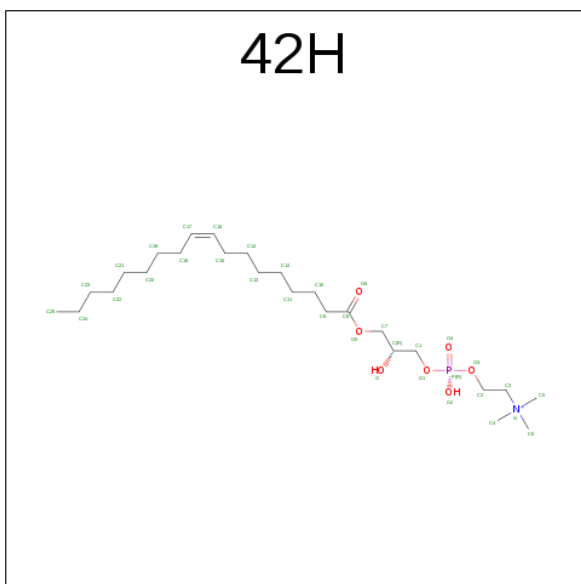
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	243	Total	C	N	O	S	0	0	0
			1930	1223	336	361	10			
4	H	243	Total	C	N	O	S	0	0	0
			1927	1222	335	360	10			

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is (4R,7R,18Z)-4,7-dihydroxy-N,N,N-trimethyl-10-oxo-3,5,9-trioxa-4-phosphaheptacos-18-en-1-aminium 4-oxide (three-letter code: 42H) (formula: C<sub>26</sub>H<sub>53</sub>NO<sub>7</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			35	26	1	7	1		
6	C	1	Total	C	N	O	P	0	0
			35	26	1	7	1		

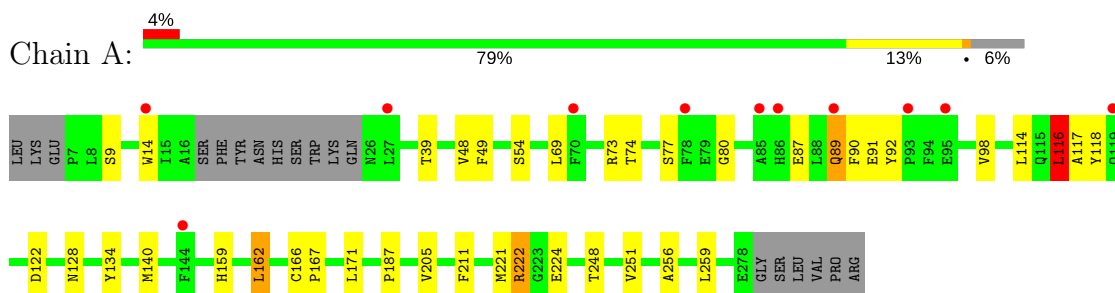
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	E	3	Total	O	0	0
			3	3		
7	F	7	Total	O	0	0
			7	7		
7	G	4	Total	O	0	0
			4	4		
7	H	7	Total	O	0	0
			7	7		

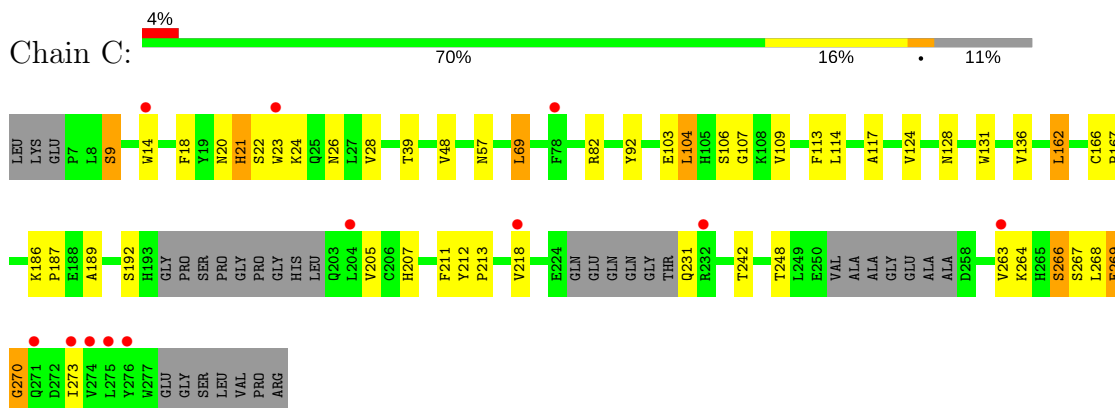
### 3 Residue-property plots

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

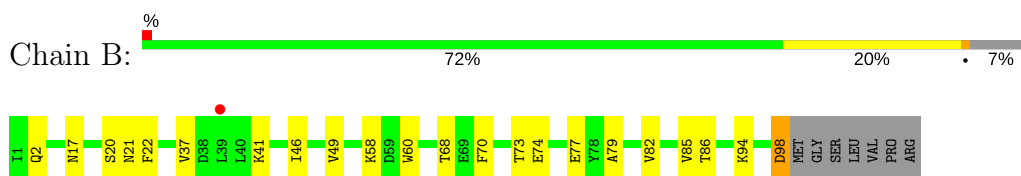
#### • Molecule 1: T-cell surface glycoprotein CD1a



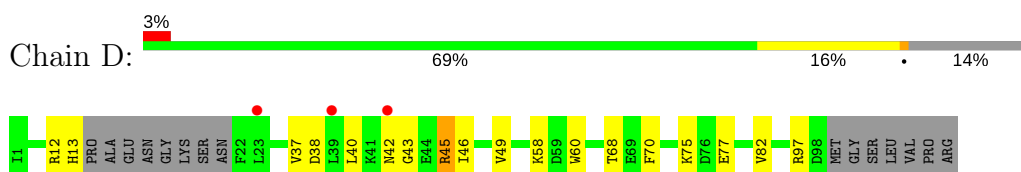
#### • Molecule 1: T-cell surface glycoprotein CD1a



#### • Molecule 2: Beta-2-microglobulin

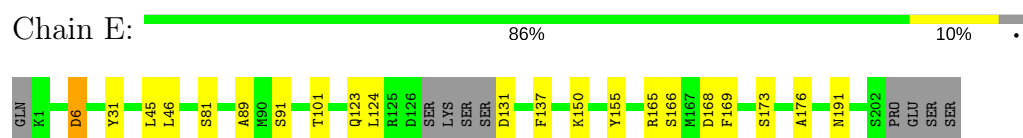


#### • Molecule 2: Beta-2-microglobulin

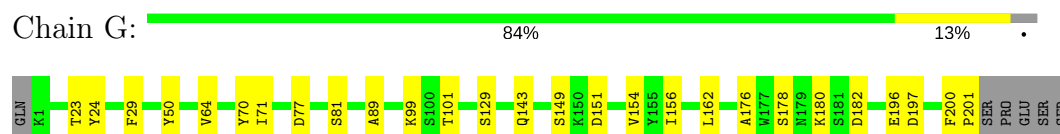




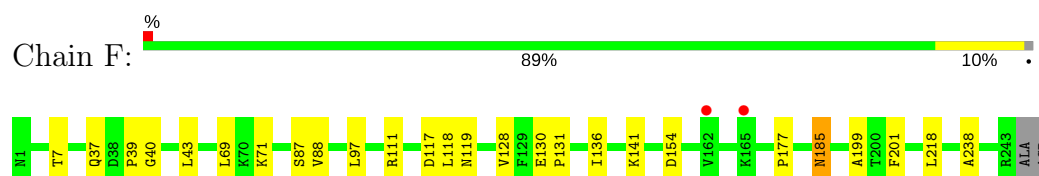
## ● Molecule 3: TCR alpha



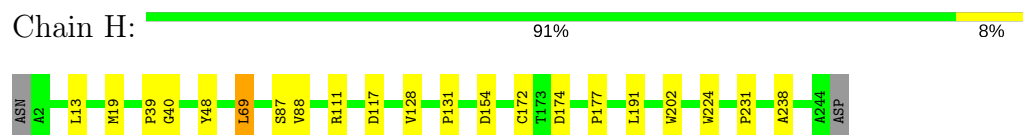
## ● Molecule 3: TCR alpha



## ● Molecule 4: TCR beta



## ● Molecule 4: TCR beta



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.27Å 142.14Å 175.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	74.75 – 2.80 79.59 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (74.75-2.80) 99.9 (79.59-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.04 (at 2.82Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.176 , 0.236 0.176 , 0.235	Depositor DCC
$R_{free}$ test set	2648 reflections (4.75%)	DCC
Wilson B-factor (Å <sup>2</sup> )	58.0	Xtriage
Anisotropy	0.441	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 58.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	12865	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 42H, NAG

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	0.44	0/2189	0.62	1/2978 (0.0%)
1	C	0.43	0/2120	0.59	0/2879
2	B	0.35	0/843	0.52	0/1142
2	D	0.38	0/783	0.54	0/1059
3	E	0.53	0/1583	0.65	0/2146
3	G	0.50	0/1605	0.61	0/2176
4	F	0.52	0/1983	0.67	0/2696
4	H	0.54	0/1980	0.64	1/2692 (0.0%)
All	All	0.48	0/13086	0.62	2/17768 (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	116	LEU	CA-CB-CG	7.10	131.62	115.30
4	H	69	LEU	CA-CB-CG	-5.04	103.71	115.30

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	2122	0	2013	24	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2054	0	1948	32	0
2	B	820	0	785	12	0
2	D	762	0	728	11	0
3	E	1548	0	1459	11	0
3	G	1569	0	1485	16	0
4	F	1930	0	1854	15	0
4	H	1927	0	1850	14	0
5	A	28	0	26	0	0
5	C	14	0	13	1	0
6	A	35	0	52	6	0
6	C	35	0	52	0	0
7	E	3	0	0	0	0
7	F	7	0	0	0	0
7	G	4	0	0	0	0
7	H	7	0	0	0	0
All	All	12865	0	12265	128	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 (128) 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:80:GLY:HA3	6:A:503:42H:H13	1.45	0.98
4:F:185:ASN:OD1	4:F:185:ASN:N	2.12	0.82
6:A:503:42H:O2	6:A:503:42H:H11	1.80	0.81
1:C:109:VAL:HG21	1:C:167:PRO:HB3	1.63	0.78
1:A:80:GLY:CA	6:A:503:42H:H13	2.15	0.77
1:C:9:SER:HB2	1:C:103:GLU:HG2	1.67	0.76
4:H:88:VAL:HG22	4:H:111:ARG:HG2	1.68	0.76
1:A:187:PRO:HB3	1:A:211:PHE:HB3	1.68	0.74
3:E:6:ASP:OD2	3:E:6:ASP:N	2.17	0.72
1:C:189:ALA:HB3	1:C:273:ILE:HG21	1.72	0.70
3:G:23:THR:HG22	3:G:70:TYR:HD1	1.57	0.68
2:D:42:ASN:ND2	2:D:77:GLU:OE2	2.28	0.67
3:G:149:SER:HB2	3:G:154:VAL:HG23	1.75	0.67
2:B:98:ASP:N	2:B:98:ASP:OD2	2.31	0.63
1:A:39:THR:HG22	1:A:48:VAL:HB	1.80	0.63
3:G:154:VAL:HG12	3:G:178:SER:HB2	1.79	0.63
2:D:77:GLU:OE1	2:D:77:GLU:N	2.31	0.62
1:C:187:PRO:HB3	1:C:211:PHE:HB3	1.80	0.61

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:ARG:HB2	1:A:259:LEU:HD23	1.82	0.61
1:C:39:THR:HG22	1:C:48:VAL:HB	1.83	0.61
2:D:37:VAL:HB	2:D:82:VAL:HG22	1.82	0.61
4:F:136:ILE:HG23	4:F:199:ALA:HB1	1.82	0.60
1:C:266:SER:O	1:C:269:GLU:HG3	2.02	0.60
1:C:186:LYS:HD3	1:C:268:LEU:HD13	1.83	0.60
3:G:151:ASP:HB3	3:G:154:VAL:HG22	1.83	0.60
3:E:150:LYS:O	4:H:117:ASP:HB2	2.02	0.60
1:A:159:HIS:ND1	3:G:50:TYR:OH	2.25	0.60
3:G:180:LYS:HD3	3:G:182:ASP:OD2	2.02	0.59
1:A:91:GLU:HG3	1:A:92:TYR:H	1.68	0.59
1:A:87:GLU:OE1	1:A:87:GLU:N	2.36	0.59
1:C:166:CYS:HB3	1:C:167:PRO:HD3	1.87	0.57
1:C:114:LEU:HB2	1:C:162:LEU:HD21	1.87	0.55
1:C:106:SER:OG	1:C:107:GLY:N	2.39	0.55
2:B:41:LYS:HB2	2:B:46:ILE:HD11	1.88	0.55
1:C:124:VAL:HG22	1:C:131:TRP:HE3	1.71	0.54
4:H:128:VAL:HG23	4:H:238:ALA:HB3	1.90	0.54
1:A:122:ASP:HB3	1:A:134:TYR:CD1	2.42	0.54
1:C:242:THR:OG1	2:D:12:ARG:NH1	2.40	0.54
1:A:221:MET:HE3	1:A:224:GLU:HA	1.90	0.54
2:D:75:LYS:H	2:D:97:ARG:NH2	2.06	0.53
3:G:156:ILE:HG12	3:G:176:ALA:HB2	1.91	0.53
1:C:192:SER:HB3	1:C:207:HIS:NE2	2.24	0.53
1:C:104:LEU:HD22	1:C:109:VAL:HA	1.90	0.53
1:C:218:VAL:HG22	1:C:263:VAL:HG13	1.91	0.52
2:D:38:ASP:HB3	2:D:45:ARG:HG3	1.91	0.52
4:H:87:SER:OG	4:H:88:VAL:N	2.43	0.52
1:C:268:LEU:O	1:C:270:GLY:N	2.39	0.52
2:B:37:VAL:HG22	2:B:82:VAL:HG22	1.91	0.52
1:A:251:VAL:HG21	1:A:256:ALA:HB2	1.92	0.52
1:C:20:ASN:HB3	1:C:24:LYS:HE2	1.92	0.51
3:G:99:LYS:HE2	4:H:48:TYR:CE1	2.46	0.51
1:C:186:LYS:HA	1:C:267:SER:OG	2.11	0.51
2:D:46:ILE:HD12	2:D:68:THR:HG21	1.93	0.51
1:A:73:ARG:CZ	6:A:503:42H:H3	2.42	0.50
1:C:103:GLU:HG3	1:C:113:PHE:HE2	1.77	0.50
4:F:88:VAL:HG22	4:F:111:ARG:HG2	1.94	0.49
1:C:57:ASN:OD1	5:C:301:NAG:N2	2.46	0.49
1:A:166:CYS:HB3	1:A:167:PRO:HD3	1.95	0.48
3:E:89:ALA:HA	3:E:101:THR:O	2.13	0.48

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:154:ASP:CG	4:H:177:PRO:HG3	2.33	0.48
1:A:49:PHE:HB3	1:A:54:SER:HB2	1.96	0.48
1:C:23:TRP:H	1:C:82:ARG:HD2	1.79	0.48
2:B:2:GLN:HB3	2:B:86:THR:HG22	1.95	0.48
4:F:69:LEU:HA	4:F:69:LEU:HD23	1.71	0.47
2:B:49:VAL:HG12	2:B:68:THR:HB	1.96	0.47
4:F:39:PRO:HA	4:F:40:GLY:HA2	1.74	0.47
3:G:196:GLU:O	3:G:196:GLU:HG2	2.14	0.47
4:H:13:LEU:HD11	4:H:19:MET:HG2	1.95	0.47
1:A:89:GLN:HG3	1:A:118:TYR:CE1	2.50	0.46
3:E:165:ARG:O	3:E:166:SER:OG	2.25	0.46
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.50	0.46
2:B:17:ASN:OD1	2:B:74:GLU:N	2.48	0.46
1:C:231:GLN:N	1:C:248:THR:O	2.48	0.46
3:G:162:LEU:HB3	4:H:172:CYS:HB2	1.97	0.46
3:E:155:TYR:O	3:E:176:ALA:HA	2.16	0.46
3:E:150:LYS:HB3	3:E:191:ASN:ND2	2.31	0.46
4:H:39:PRO:HA	4:H:40:GLY:HA2	1.63	0.46
3:E:31:TYR:HB2	3:E:91:SER:HB2	1.97	0.46
2:D:49:VAL:HG12	2:D:68:THR:HB	1.99	0.45
4:F:154:ASP:CG	4:F:177:PRO:HG3	2.35	0.45
1:C:205:VAL:HA	1:C:248:THR:HG22	1.97	0.45
3:E:123:GLN:C	3:E:124:LEU:HD12	2.35	0.45
1:C:117:ALA:HB2	2:D:60:TRP:CE2	2.51	0.45
4:F:37:GLN:HB2	4:F:43:LEU:HD23	1.98	0.45
3:G:89:ALA:HA	3:G:101:THR:O	2.17	0.45
4:F:87:SER:OG	4:F:88:VAL:N	2.50	0.44
2:B:21:ASN:OD1	2:B:22:PHE:N	2.40	0.44
4:H:131:PRO:HD2	4:H:202:TRP:CZ2	2.53	0.44
4:H:69:LEU:HA	4:H:69:LEU:HD23	1.64	0.44
3:E:137:PHE:O	3:E:173:SER:HA	2.18	0.43
4:F:118:LEU:HD22	4:F:218:LEU:HD21	2.00	0.43
1:A:73:ARG:NH2	6:A:503:42H:H3	2.33	0.43
1:C:212:TYR:CG	1:C:213:PRO:HA	2.54	0.43
1:C:104:LEU:HA	1:C:104:LEU:HD22	1.87	0.43
4:H:224:TRP:CZ2	4:H:231:PRO:HD3	2.54	0.43
4:F:128:VAL:HG23	4:F:238:ALA:HB3	2.00	0.43
1:C:109:VAL:CG2	1:C:167:PRO:HB3	2.42	0.43
1:A:114:LEU:HB2	1:A:162:LEU:HD21	2.01	0.42
1:A:171:LEU:HD23	1:A:171:LEU:HA	1.71	0.42
1:A:14:TRP:HB2	1:A:98:VAL:HB	2.01	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:20:SER:O	2:B:20:SER:OG	2.36	0.42
3:G:64:VAL:HG13	3:G:71:ILE:HG22	2.01	0.42
1:A:205:VAL:HG22	1:A:248:THR:HG22	2.01	0.42
1:C:20:ASN:OD1	1:C:21:HIS:N	2.51	0.42
2:B:79:ALA:HA	2:B:94:LYS:HA	2.02	0.42
2:D:40:LEU:HD13	2:D:43:GLY:HA2	2.01	0.42
1:A:98:VAL:HG22	1:A:116:LEU:HD22	2.02	0.42
1:C:20:ASN:O	1:C:22:SER:N	2.48	0.42
2:D:75:LYS:HA	2:D:75:LYS:HD2	1.70	0.42
4:F:154:ASP:OD1	4:F:177:PRO:HG3	2.20	0.42
1:A:91:GLU:CG	1:A:92:TYR:H	2.30	0.41
4:F:117:ASP:OD2	4:F:119:ASN:HB2	2.20	0.41
1:A:74:THR:O	1:A:77:SER:OG	2.38	0.41
1:C:18:PHE:HB3	1:C:92:TYR:HE1	1.86	0.41
3:E:169:PHE:CE2	4:F:141:LYS:HE2	2.55	0.41
3:G:200:PHE:HA	3:G:201:PRO:HD2	1.93	0.41
4:H:154:ASP:OD1	4:H:177:PRO:HG3	2.20	0.41
4:H:174:ASP:HB2	4:H:191:LEU:HD12	2.03	0.41
4:F:130:GLU:HA	4:F:131:PRO:HD3	1.92	0.41
6:A:503:42H:H7	6:A:503:42H:H15	1.78	0.40
1:C:69:LEU:HD21	4:F:97:LEU:HD13	2.03	0.40
3:G:197:ASP:OD1	3:G:197:ASP:N	2.41	0.40
1:A:117:ALA:HB2	2:B:60:TRP:CD2	2.57	0.40
1:C:14:TRP:CD1	1:C:28:VAL:HG22	2.57	0.40
3:G:143:GLN:N	3:G:143:GLN:OE1	2.55	0.40
3:G:24:TYR:CD2	3:G:29:PHE:HD2	2.40	0.40
2:B:77:GLU:N	2:B:77:GLU:OE1	2.54	0.40
3:E:45:LEU:HG	3:E:46:LEU:N	2.35	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	259/281 (92%)	249 (96%)	8 (3%)	2 (1%)	22	55
1	C	241/281 (86%)	228 (95%)	9 (4%)	4 (2%)	11	34
2	B	96/105 (91%)	94 (98%)	2 (2%)	0	100	100
2	D	86/105 (82%)	83 (96%)	3 (4%)	0	100	100
3	E	194/207 (94%)	188 (97%)	6 (3%)	0	100	100
3	G	199/207 (96%)	190 (96%)	9 (4%)	0	100	100
4	F	241/245 (98%)	235 (98%)	6 (2%)	0	100	100
4	H	241/245 (98%)	232 (96%)	9 (4%)	0	100	100
All	All	1557/1676 (93%)	1499 (96%)	52 (3%)	6 (0%)	38	72

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	PHE
1	C	269	GLU
1	A	128	ASN
1	C	128	ASN
1	C	21	HIS
1	C	270	GLY

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	228/245 (93%)	221 (97%)	7 (3%)	45	79
1	C	223/245 (91%)	215 (96%)	8 (4%)	40	74
2	B	93/99 (94%)	89 (96%)	4 (4%)	33	67
2	D	86/99 (87%)	82 (95%)	4 (5%)	30	64
3	E	178/187 (95%)	174 (98%)	4 (2%)	57	87
3	G	181/187 (97%)	178 (98%)	3 (2%)	66	90
4	F	208/209 (100%)	204 (98%)	4 (2%)	62	89

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	H	207/209 (99%)	207 (100%)	0	100	100
All	All	1404/1480 (95%)	1370 (98%)	34 (2%)	54	85

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	69	LEU
1	A	89	GLN
1	A	116	LEU
1	A	140	MET
1	A	162	LEU
1	A	222	ARG
2	B	58	LYS
2	B	73	THR
2	B	85	VAL
2	B	98	ASP
1	C	9	SER
1	C	26	ASN
1	C	69	LEU
1	C	104	LEU
1	C	136	VAL
1	C	162	LEU
1	C	264	LYS
1	C	266	SER
2	D	13	HIS
2	D	45	ARG
2	D	58	LYS
2	D	70	PHE
3	E	6	ASP
3	E	81	SER
3	E	131	ASP
3	E	168	ASP
4	F	7	THR
4	F	71	LYS
4	F	185	ASN
4	F	201	PHE
3	G	77	ASP
3	G	81	SER
3	G	129	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	127	GLN
1	C	119	GLN
2	D	42	ASN
4	F	226	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

5 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	$\# Z  > 2$
5	NAG	A	501	1	14,14,15	0.81	1 (7%)	15,19,21	0.53	0
5	NAG	A	502	1	14,14,15	0.97	1 (7%)	15,19,21	1.29	2 (13%)
6	42H	A	503	-	34,34,34	0.37	0	38,41,41	0.37	0
5	NAG	C	301	1	14,14,15	0.66	0	15,19,21	0.64	0
6	42H	C	302	-	34,34,34	0.40	0	38,41,41	0.39	0

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
5	NAG	A	501	1	-	0/6/23/26	0/1/1/1
5	NAG	A	502	1	-	0/6/23/26	0/1/1/1
6	42H	A	503	-	-	0/36/36/36	0/0/0/0
5	NAG	C	301	1	-	0/6/23/26	0/1/1/1
6	42H	C	302	-	-	0/36/36/36	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	502	NAG	O5-C1	-2.81	1.39	1.43
5	A	501	NAG	C1-C2	2.69	1.56	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	502	NAG	C1-O5-C5	-3.30	107.61	112.17
5	A	502	NAG	C4-C3-C2	2.32	114.42	111.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	503	42H	6	0
5	C	301	NAG	1	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

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	263/281 (93%)	0.02	11 (4%) 37 26	39, 63, 126, 153	0
1	C	249/281 (88%)	0.07	12 (4%) 31 21	42, 78, 134, 180	0
2	B	98/105 (93%)	0.11	1 (1%) 82 77	52, 95, 147, 171	0
2	D	90/105 (85%)	0.36	3 (3%) 47 36	60, 100, 154, 187	0
3	E	198/207 (95%)	-0.18	0 100 100	28, 49, 99, 150	0
3	G	201/207 (97%)	-0.19	0 100 100	32, 59, 130, 180	0
4	F	243/245 (99%)	-0.09	2 (0%) 86 81	26, 52, 102, 142	0
4	H	243/245 (99%)	-0.23	0 100 100	30, 53, 92, 119	0
All	All	1585/1676 (94%)	-0.06	29 (1%) 69 60	26, 63, 129, 187	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	271	GLN	4.5
4	F	162	VAL	3.9
2	D	42	ASN	3.2
1	A	14	TRP	2.8
1	C	263	VAL	2.8
1	C	276	TYR	2.8
1	A	93	PRO	2.8
2	B	39	LEU	2.7
1	A	85	ALA	2.7
1	A	89	GLN	2.7
1	C	218	VAL	2.6
1	C	204	LEU	2.6
2	D	23	LEU	2.6
1	A	95	GLU	2.5
1	C	232	ARG	2.5
1	A	144	PHE	2.4

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	C	274	VAL	2.3
1	A	78	PHE	2.3
4	F	165	LYS	2.3
1	C	14	TRP	2.2
1	A	119	GLN	2.2
1	A	70	PHE	2.1
1	A	86	HIS	2.1
1	A	27	LEU	2.1
2	D	39	LEU	2.1
1	C	23	TRP	2.0
1	C	275	LEU	2.0
1	C	273	ILE	2.0
1	C	78	PHE	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
6	42H	A	503	35/35	0.86	0.60	5.96	46,97,205,206	0
6	42H	C	302	35/35	0.80	0.54	5.77	55,92,296,297	0
5	NAG	A	501	14/15	0.90	0.28	3.35	80,101,112,113	0
5	NAG	A	502	14/15	0.86	0.25	-	86,101,123,125	0
5	NAG	C	301	14/15	0.82	0.23	-	92,109,120,120	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.