



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 9, 2020 – 05:04 PM GMT

PDB ID : 6T80
Title : Human 14-3-3 sigma fused to the AANAT peptide including phosphoserine-205
Authors : Sluchanko, N.N.; Tugaeva, K.V.; Titterington, J.; Antson, A.A.
Deposited on : 2019-10-23
Resolution : 2.99 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 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 : 2.14.6
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) : 2.14.6

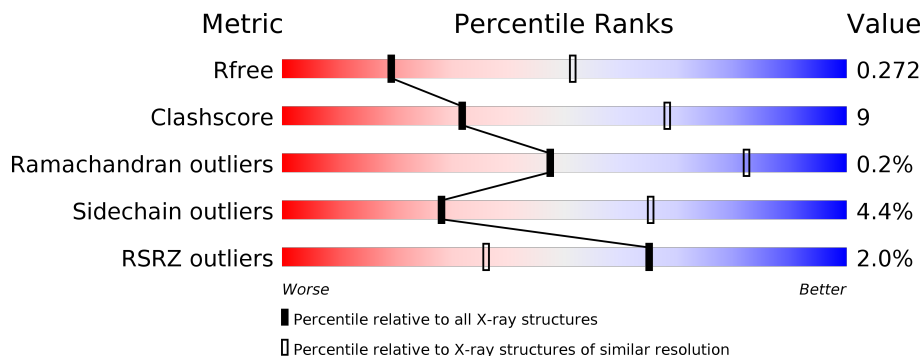
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.99 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.



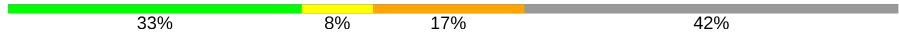
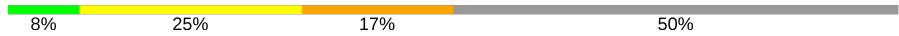
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 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	239	
1	B	239	
1	C	239	
1	D	239	
2	E	12	
2	F	12	

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Mol	Chain	Length	Quality of chain
2	G	12	
2	H	12	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7240 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 14-3-3 protein sigma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	233	1811	1131	304	366	10	0	0	0
1	B	228	1768	1103	298	358	9	0	0	0
1	C	220	1701	1066	284	341	10	0	0	0
1	D	226	1738	1085	292	352	9	0	0	0

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P31947
A	-1	PRO	-	expression tag	UNP P31947
A	0	HIS	-	expression tag	UNP P31947
A	75	ALA	GLU	engineered mutation	UNP P31947
A	76	ALA	GLU	engineered mutation	UNP P31947
A	77	ALA	LYS	engineered mutation	UNP P31947
A	232	GLY	-	linker	UNP P31947
A	233	SER	-	linker	UNP P31947
A	234	GLY	-	linker	UNP P31947
A	235	SER	-	linker	UNP P31947
A	236	LEU	-	linker	UNP P31947
B	-2	GLY	-	expression tag	UNP P31947
B	-1	PRO	-	expression tag	UNP P31947
B	0	HIS	-	expression tag	UNP P31947
B	75	ALA	GLU	engineered mutation	UNP P31947
B	76	ALA	GLU	engineered mutation	UNP P31947
B	77	ALA	LYS	engineered mutation	UNP P31947
B	232	GLY	-	linker	UNP P31947
B	233	SER	-	linker	UNP P31947
B	234	GLY	-	linker	UNP P31947
B	235	SER	-	linker	UNP P31947

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Chain	Residue	Modelled	Actual	Comment	Reference
B	236	LEU	-	linker	UNP P31947
C	-2	GLY	-	expression tag	UNP P31947
C	-1	PRO	-	expression tag	UNP P31947
C	0	HIS	-	expression tag	UNP P31947
C	75	ALA	GLU	engineered mutation	UNP P31947
C	76	ALA	GLU	engineered mutation	UNP P31947
C	77	ALA	LYS	engineered mutation	UNP P31947
C	232	GLY	-	linker	UNP P31947
C	233	SER	-	linker	UNP P31947
C	234	GLY	-	linker	UNP P31947
C	235	SER	-	linker	UNP P31947
C	236	LEU	-	linker	UNP P31947
D	-2	GLY	-	expression tag	UNP P31947
D	-1	PRO	-	expression tag	UNP P31947
D	0	HIS	-	expression tag	UNP P31947
D	75	ALA	GLU	engineered mutation	UNP P31947
D	76	ALA	GLU	engineered mutation	UNP P31947
D	77	ALA	LYS	engineered mutation	UNP P31947
D	232	GLY	-	linker	UNP P31947
D	233	SER	-	linker	UNP P31947
D	234	GLY	-	linker	UNP P31947
D	235	SER	-	linker	UNP P31947
D	236	LEU	-	linker	UNP P31947

- Molecule 2 is a protein called AANAT peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
2	E	6	Total	C	N	O	P	S	0	0	0
			50	24	13	11	1	1			
2	F	7	Total	C	N	O	P	S	0	0	0
			58	30	14	12	1	1			
2	G	7	Total	C	N	O	P	S	0	0	0
			58	30	14	12	1	1			
2	H	6	Total	C	N	O	P		0	0	0
			54	28	13	12	1				

- Molecule 3 is water.

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

3 Residue-property plots [i](#)


These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

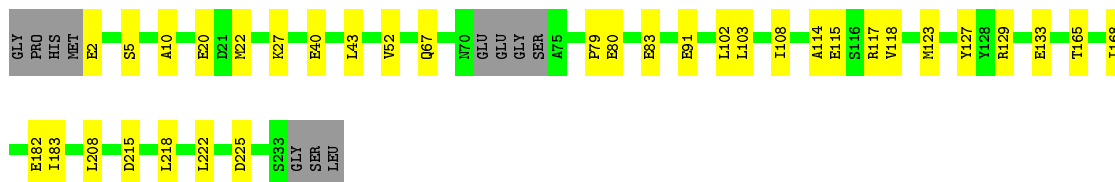
- Molecule 1: 14-3-3 protein sigma

Chain A: 




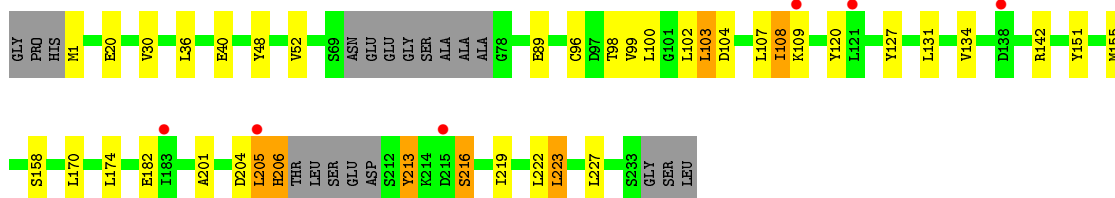
- Molecule 1: 14-3-3 protein sigma

Chain B: 




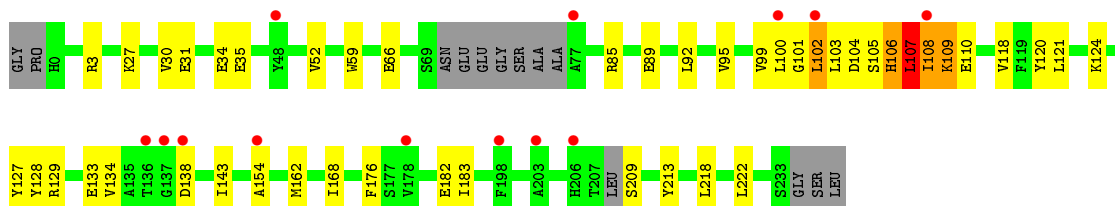
- Molecule 1: 14-3-3 protein sigma

Chain C: 




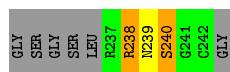
- Molecule 1: 14-3-3 protein sigma

Chain D: 




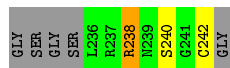
- Molecule 2: AANAT peptide

Chain E: 



- Molecule 2: AANAT peptide

Chain F: 




- Molecule 2: AANAT peptide

Chain G: 



- Molecule 2: AANAT peptide

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.26Å 74.10Å 102.09Å 90.00° 96.03° 90.00°	Depositor
Resolution (Å)	63.51 – 2.99 74.84 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.6 (63.51-2.99) 99.6 (74.84-2.99)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 3.01Å)	Xtrriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.200 , 0.251 0.224 , 0.272	Depositor DCC
R_{free} test set	1080 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	63.3	Xtrriage
Anisotropy	0.641	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 58.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7240	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.53	0/1838	0.69	0/2478
1	B	0.52	0/1793	0.68	0/2421
1	C	0.54	0/1724	0.78	3/2323 (0.1%)
1	D	0.49	0/1762	0.71	2/2378 (0.1%)
2	E	0.38	0/38	0.78	0/46
2	F	0.53	0/46	0.78	0/57
2	G	0.50	0/46	0.79	0/57
2	H	0.77	0/43	1.02	0/55
All	All	0.52	0/7290	0.72	5/9815 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	205	LEU	CB-CA-C	-7.77	95.43	110.20
1	D	107	LEU	N-CA-C	-6.86	92.49	111.00
1	C	205	LEU	N-CA-C	-6.46	93.56	111.00
1	C	158	SER	CB-CA-C	5.76	121.04	110.10
1	D	102	LEU	CB-CA-C	5.10	119.89	110.20

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	1811	0	1752	17	0
1	B	1768	0	1704	26	0
1	C	1701	0	1630	32	0
1	D	1738	0	1666	56	0
2	E	50	0	42	5	0
2	F	58	0	53	1	0
2	G	58	0	53	3	0
2	H	54	0	50	2	0
3	A	1	0	0	0	0
3	E	1	0	0	0	0
All	All	7240	0	6950	126	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 9.

All (126) 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:D:27:LYS:HG3	1:D:102:LEU:CD1	1.66	1.24
1:D:27:LYS:HG3	1:D:102:LEU:HD11	1.25	1.12
1:D:30:VAL:HG11	1:D:107:LEU:HD11	1.28	1.06
1:D:30:VAL:HG11	1:D:107:LEU:CD1	1.89	1.02
1:D:95:VAL:O	1:D:99:VAL:HG23	1.63	0.99
1:D:107:LEU:HD23	1:D:120:TYR:CZ	1.98	0.98
1:D:27:LYS:CG	1:D:102:LEU:HD11	1.94	0.95
1:D:108:ILE:HG12	1:D:120:TYR:HB3	1.50	0.94
1:D:31:GLU:HA	1:D:106:HIS:HE1	1.36	0.90
1:D:30:VAL:CG1	1:D:107:LEU:HD11	2.01	0.89
1:D:27:LYS:HG3	1:D:102:LEU:HD13	1.54	0.87
1:D:31:GLU:HA	1:D:106:HIS:CE1	2.10	0.87
1:C:99:VAL:O	1:C:103:LEU:HB2	1.77	0.85
1:D:118:VAL:CG1	1:D:168:ILE:HG22	2.08	0.83
1:D:118:VAL:HG11	1:D:168:ILE:HG22	1.64	0.79
1:D:27:LYS:HE2	1:D:102:LEU:HD11	1.64	0.79
2:E:238:ARG:HH12	2:E:240:SEP:HB3	1.48	0.79
1:C:170:LEU:HB3	1:C:219:ILE:HG21	1.67	0.77
1:D:30:VAL:CG1	1:D:107:LEU:CD1	2.60	0.77
1:C:103:LEU:HA	1:C:107:LEU:HD12	1.69	0.74
1:D:27:LYS:CD	1:D:102:LEU:HD11	2.18	0.73
1:D:27:LYS:CG	1:D:102:LEU:CD1	2.55	0.73
1:C:227:LEU:CA	1:C:227:LEU:CG	2.68	0.72
1:B:129:ARG:HG3	1:B:183:ILE:HG13	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:92:LEU:HD23	1:D:92:LEU:C	2.13	0.69
1:B:108:ILE:HG23	1:B:117:ARG:HH22	1.55	0.69
1:C:30:VAL:HG22	1:C:36:LEU:HD11	1.74	0.68
1:D:209:SER:O	1:D:213:TYR:CB	2.40	0.68
1:D:59:TRP:CD1	1:D:89:GLU:HB2	2.30	0.67
1:D:107:LEU:HD23	1:D:120:TYR:OH	1.93	0.67
1:D:27:LYS:CE	1:D:102:LEU:HD11	2.23	0.67
1:C:170:LEU:HB3	1:C:219:ILE:CG2	2.26	0.65
1:D:108:ILE:HA	1:D:120:TYR:HD2	1.62	0.65
2:E:238:ARG:NH1	2:E:240:SEP:HB3	2.12	0.64
1:D:129:ARG:HG3	1:D:183:ILE:HG13	1.79	0.63
1:C:174:LEU:HD13	1:C:223:LEU:HG	1.81	0.62
1:A:0:HIS:CE1	1:B:79:PRO:HG3	2.36	0.61
1:B:27:LYS:HG3	1:B:102:LEU:HD21	1.83	0.60
1:C:205:LEU:O	1:C:206:HIS:HB2	2.00	0.60
1:A:0:HIS:ND1	1:B:79:PRO:HG3	2.17	0.60
1:B:225:ASP:HB2	2:H:239:ASN:HD21	1.67	0.60
1:D:30:VAL:HG11	1:D:107:LEU:HD12	1.83	0.59
1:C:213:TYR:N	1:C:213:TYR:HD1	2.00	0.59
1:D:31:GLU:OE2	1:D:102:LEU:HD21	2.03	0.58
1:D:35:GLU:HB3	1:D:110:GLU:OE2	2.03	0.58
1:C:213:TYR:N	1:C:213:TYR:CD1	2.72	0.57
1:D:107:LEU:C	1:D:120:TYR:HE2	2.08	0.57
1:D:108:ILE:HD11	1:D:121:LEU:HG	1.87	0.55
1:C:52:VAL:HG21	1:C:127:TYR:HE2	1.72	0.55
1:A:129:ARG:HG3	1:A:183:ILE:HG13	1.89	0.55
1:C:213:TYR:HA	1:C:216:SER:HB2	1.89	0.55
1:C:96:CYS:O	1:C:100:LEU:N	2.39	0.55
1:B:52:VAL:HG21	1:B:127:TYR:HE1	1.71	0.54
1:C:213:TYR:H	1:C:213:TYR:HD1	1.54	0.54
1:C:100:LEU:HD12	1:C:131:LEU:HD11	1.90	0.54
1:A:9:LYS:HE2	1:B:80:GLU:HA	1.89	0.54
1:A:182:GLU:CD	2:G:238:ARG:HE	2.11	0.53
1:D:121:LEU:O	1:D:124:LYS:HB3	2.08	0.53
1:B:108:ILE:CG2	1:B:117:ARG:HH22	2.21	0.53
1:D:118:VAL:CG1	1:D:168:ILE:CG2	2.84	0.53
1:D:35:GLU:OE2	1:D:110:GLU:OE1	2.27	0.53
1:D:118:VAL:HG12	1:D:168:ILE:CG2	2.39	0.53
1:D:118:VAL:HG12	1:D:168:ILE:HG22	1.89	0.52
1:D:118:VAL:HG21	1:D:162:MET:SD	2.50	0.52
1:C:48:TYR:CD1	1:C:99:VAL:CG2	2.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:ILE:HG23	1:B:117:ARG:NH2	2.25	0.51
1:D:182:GLU:CD	2:F:238:ARG:HE	2.14	0.51
1:B:40:GLU:HA	1:B:43:LEU:HD12	1.93	0.51
1:D:129:ARG:O	1:D:133:GLU:HG3	2.11	0.51
1:D:59:TRP:CE2	1:D:134:VAL:HG12	2.46	0.51
1:A:52:VAL:HG21	1:A:127:TYR:HE2	1.76	0.51
1:B:182:GLU:CD	2:H:238:ARG:HE	2.14	0.51
1:D:106:HIS:CD2	1:D:106:HIS:N	2.79	0.50
1:D:30:VAL:CG1	1:D:107:LEU:HD12	2.40	0.50
1:D:101:GLY:O	1:D:104:ASP:N	2.45	0.50
1:D:52:VAL:HG21	1:D:127:TYR:HE1	1.75	0.50
1:D:3:ARG:CZ	1:D:34:GLU:HB2	2.42	0.50
1:C:170:LEU:CB	1:C:219:ILE:HG21	2.37	0.49
1:D:107:LEU:CA	1:D:120:TYR:HE2	2.25	0.49
1:C:182:GLU:CD	2:E:238:ARG:HE	2.15	0.49
1:A:-1:PRO:HD2	1:A:0:HIS:HD2	1.78	0.48
1:A:103:LEU:HA	1:A:107:LEU:HB2	1.96	0.48
1:D:66:GLU:OE1	1:D:85:ARG:NE	2.45	0.48
1:D:107:LEU:C	1:D:120:TYR:CE2	2.85	0.48
1:C:98:THR:O	1:C:102:LEU:HB2	2.14	0.47
1:A:18:ARG:HH22	1:B:91:GLU:CD	2.18	0.47
1:A:40:GLU:HA	1:A:43:LEU:HD12	1.98	0.46
1:C:89:GLU:HG3	1:C:134:VAL:HB	1.98	0.46
1:B:103:LEU:HD11	1:B:123:MET:HG2	1.97	0.45
1:A:0:HIS:ND1	1:B:79:PRO:CG	2.79	0.45
1:C:96:CYS:O	1:C:100:LEU:HB2	2.16	0.45
1:B:215:ASP:HA	1:B:218:LEU:HD12	1.99	0.45
1:C:98:THR:O	1:C:102:LEU:CB	2.64	0.45
1:C:104:ASP:HA	1:C:108:ILE:CG1	2.46	0.45
1:B:2:GLU:O	1:B:5:SER:N	2.50	0.45
1:D:99:VAL:O	1:D:103:LEU:HG	2.17	0.45
1:B:165:THR:HB	1:B:208:LEU:HD11	2.00	0.44
1:A:0:HIS:HB2	1:B:83:GLU:OE1	2.17	0.44
1:D:100:LEU:HD21	1:D:127:TYR:HB2	2.00	0.44
1:A:222:LEU:HD13	2:G:239:ASN:HB3	2.00	0.44
1:B:129:ARG:O	1:B:133:GLU:HG3	2.18	0.44
1:A:165:THR:HB	1:A:208:LEU:HD11	2.00	0.43
1:B:27:LYS:HA	1:B:102:LEU:HD11	2.01	0.43
1:A:9:LYS:HG3	1:B:80:GLU:HB3	2.00	0.43
1:D:128:TYR:HD1	1:D:143:ILE:HG23	1.83	0.43
1:A:30:VAL:HG11	1:A:102:LEU:HD22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:222:LEU:CD2	2:E:239:ASN:HB3	2.49	0.43
1:C:222:LEU:HD22	2:E:239:ASN:HB3	1.99	0.43
1:B:114:ALA:O	1:B:118:VAL:HG23	2.19	0.42
1:C:151:TYR:O	1:C:155:MET:HG2	2.19	0.42
1:B:117:ARG:HA	1:B:117:ARG:HD2	1.81	0.42
1:C:104:ASP:O	1:C:104:ASP:OD1	2.38	0.42
1:C:201:ALA:O	1:C:204:ASP:CG	2.58	0.42
1:D:95:VAL:O	1:D:99:VAL:CG2	2.52	0.42
1:D:107:LEU:CA	1:D:120:TYR:CE2	3.03	0.41
1:B:115:GLU:CD	1:B:168:ILE:HD13	2.41	0.41
1:C:103:LEU:CD2	1:C:120:TYR:HB3	2.51	0.41
1:D:108:ILE:CA	1:D:120:TYR:HD2	2.32	0.41
1:B:10:ALA:O	1:B:22:MET:HG3	2.21	0.41
1:C:48:TYR:CD1	1:C:99:VAL:HG21	2.56	0.41
1:D:100:LEU:HA	1:D:103:LEU:HD12	2.02	0.41
1:D:121:LEU:HB3	1:D:154:ALA:HB2	2.02	0.41
1:C:108:ILE:HG12	1:C:108:ILE:H	1.61	0.40
1:D:109:LYS:NZ	1:D:110:GLU:HB2	2.36	0.40
1:A:222:LEU:HD22	2:G:239:ASN:HD22	1.86	0.40
1:C:219:ILE:HA	1:C:222:LEU:HD12	2.02	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	229/239 (96%)	223 (97%)	6 (3%)	0	100	100
1	B	224/239 (94%)	212 (95%)	12 (5%)	0	100	100
1	C	214/239 (90%)	200 (94%)	13 (6%)	1 (0%)	29	68
1	D	220/239 (92%)	204 (93%)	15 (7%)	1 (0%)	29	68

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	3/12 (25%)	3 (100%)	0	0	100	100
2	F	4/12 (33%)	4 (100%)	0	0	100	100
2	G	4/12 (33%)	3 (75%)	1 (25%)	0	100	100
2	H	4/12 (33%)	4 (100%)	0	0	100	100
All	All	902/1004 (90%)	853 (95%)	47 (5%)	2 (0%)	47	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	107	LEU
1	C	109	LYS

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	189/198 (96%)	187 (99%)	2 (1%)	73	90
1	B	183/198 (92%)	180 (98%)	3 (2%)	62	86
1	C	174/198 (88%)	164 (94%)	10 (6%)	20	56
1	D	179/198 (90%)	170 (95%)	9 (5%)	24	60
2	E	4/7 (57%)	3 (75%)	1 (25%)	0	3
2	F	5/7 (71%)	3 (60%)	2 (40%)	0	0
2	G	5/7 (71%)	3 (60%)	2 (40%)	0	0
2	H	5/7 (71%)	1 (20%)	4 (80%)	0	0
All	All	744/820 (91%)	711 (96%)	33 (4%)	28	65

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	192	SER
1	A	233	SER

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Mol	Chain	Res	Type
1	B	20	GLU
1	B	67	GLN
1	B	222	LEU
1	C	1	MET
1	C	20	GLU
1	C	40	GLU
1	C	103	LEU
1	C	108	ILE
1	C	142	ARG
1	C	206	HIS
1	C	213	TYR
1	C	216	SER
1	C	223	LEU
1	D	105	SER
1	D	106	HIS
1	D	107	LEU
1	D	108	ILE
1	D	109	LYS
1	D	138	ASP
1	D	176	PHE
1	D	218	LEU
1	D	222	LEU
2	E	238	ARG
2	F	238	ARG
2	F	242	CYS
2	G	238	ARG
2	G	239	ASN
2	H	236	LEU
2	H	237	ARG
2	H	238	ARG
2	H	239	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	106	HIS
1	B	55	GLN
1	C	55	GLN
1	D	55	GLN
1	D	106	HIS
2	G	239	ASN

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Mol	Chain	Res	Type
2	H	239	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SEP	E	240	2	8,9,10	0.93	0	8,12,14	4.21	2 (25%)
2	SEP	F	240	2	8,9,10	1.23	1 (12%)	8,12,14	2.71	5 (62%)
2	SEP	G	240	2	8,9,10	1.16	1 (12%)	8,12,14	2.00	4 (50%)
2	SEP	H	240	2	8,9,10	1.48	1 (12%)	8,12,14	1.41	1 (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	SEP	E	240	2	-	4/5/8/10	-
2	SEP	F	240	2	-	2/5/8/10	-
2	SEP	G	240	2	-	2/5/8/10	-
2	SEP	H	240	2	-	2/5/8/10	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	240	SEP	P-O1P	3.21	1.60	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	240	SEP	P-OG	-2.43	1.52	1.60
2	G	240	SEP	OG-CB	-2.38	1.35	1.44

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	240	SEP	OG-CB-CA	11.03	118.88	108.14
2	F	240	SEP	OG-CB-CA	5.47	113.47	108.14
2	E	240	SEP	O2P-P-OG	4.19	117.88	106.73
2	F	240	SEP	O2P-P-OG	2.79	114.16	106.73
2	H	240	SEP	OG-CB-CA	2.76	110.83	108.14
2	G	240	SEP	OG-P-O1P	2.75	114.19	106.47
2	G	240	SEP	O3P-P-O2P	2.71	117.98	107.64
2	F	240	SEP	P-OG-CB	-2.70	110.86	118.30
2	G	240	SEP	O3P-P-OG	2.44	113.24	106.73
2	G	240	SEP	O2P-P-O1P	-2.41	101.23	110.68
2	F	240	SEP	O2P-P-O1P	2.24	119.46	110.68
2	F	240	SEP	O3P-P-O2P	-2.03	99.87	107.64

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	240	SEP	CB-OG-P-O1P
2	E	240	SEP	CB-OG-P-O2P
2	E	240	SEP	CB-OG-P-O3P
2	G	240	SEP	CB-OG-P-O1P
2	F	240	SEP	CB-OG-P-O2P
2	E	240	SEP	N-CA-CB-OG
2	F	240	SEP	N-CA-CB-OG
2	G	240	SEP	CB-OG-P-O3P
2	H	240	SEP	CB-OG-P-O2P
2	H	240	SEP	CA-CB-OG-P

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	240	SEP	2	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	233/239 (97%)	-0.09	0 100 100	45, 67, 109, 130	0
1	B	228/239 (95%)	-0.04	0 100 100	43, 66, 113, 147	0
1	C	220/239 (92%)	0.04	6 (2%) 54 26	70, 96, 140, 165	0
1	D	226/239 (94%)	0.26	13 (5%) 23 7	79, 112, 163, 192	0
2	E	5/12 (41%)	0.50	0 100 100	114, 124, 124, 144	0
2	F	6/12 (50%)	-0.33	0 100 100	97, 105, 111, 117	0
2	G	6/12 (50%)	-0.10	0 100 100	74, 82, 94, 105	0
2	H	5/12 (41%)	-0.05	0 100 100	65, 75, 100, 101	0
All	All	929/1004 (92%)	0.04	19 (2%) 65 36	43, 85, 144, 192	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	108	ILE	5.3
1	C	215	ASP	5.2
1	C	138	ASP	3.7
1	D	137	GLY	3.5
1	D	100	LEU	3.0
1	D	136	THR	3.0
1	D	154	ALA	2.8
1	C	205	LEU	2.7
1	C	183	ILE	2.7
1	D	48	TYR	2.4
1	D	206	HIS	2.4
1	D	203	ALA	2.4
1	D	102	LEU	2.3
1	D	138	ASP	2.2
1	C	109	LYS	2.2
1	D	198	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	178	VAL	2.1
1	C	121	LEU	2.0
1	D	77	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	SEP	E	240	10/11	0.91	0.15	104,109,121,122	0
2	SEP	F	240	10/11	0.93	0.11	91,95,102,103	0
2	SEP	G	240	10/11	0.97	0.16	59,62,71,73	0
2	SEP	H	240	10/11	0.97	0.18	37,51,58,77	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.