



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:23 PM GMT

PDB ID : 4L3O  
Title : Crystal Structure of SIRT2 in complex with the macrocyclic peptide S2iL5  
Authors : Yamagata, K.; Nishimasu, H.; Ishitani, R.; Nureki, O.  
Deposited on : 2013-06-06  
Resolution : 2.52 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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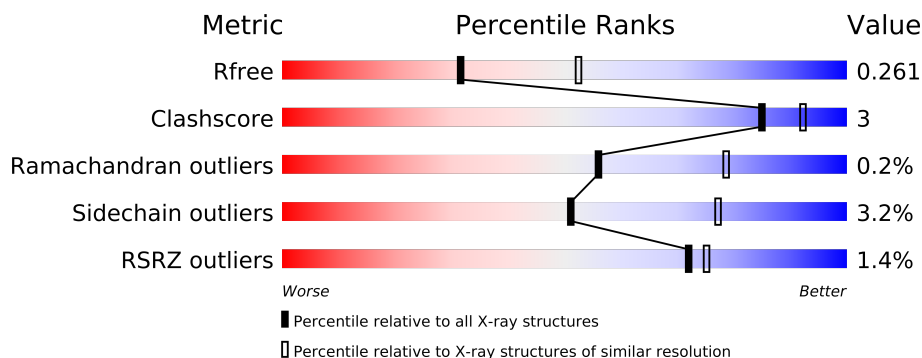
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.52 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))
$R_{free}$	66092	3240 (2.54-2.50)
Clashscore	79885	4080 (2.54-2.50)
Ramachandran outliers	78287	3990 (2.54-2.50)
Sidechain outliers	78261	3992 (2.54-2.50)
RSRZ outliers	66119	3241 (2.54-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	302	
1	B	302	
1	C	302	
1	D	302	
2	E	16	
2	F	16	
2	G	16	
2	H	16	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	MES	A	502	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
4	MES	B	502	-	X
4	MES	D	502	-	X
5	EDO	C	504	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10482 atoms, of which 0 are hydrogen and 0 are deuterium.

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 NAD-dependent protein deacetylase sirtuin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	1	0
			2368	1520	393	436	19			
1	B	295	Total	C	N	O	S	0	0	0
			2329	1497	388	425	19			
1	C	294	Total	C	N	O	S	0	0	0
			2320	1491	387	423	19			
1	D	293	Total	C	N	O	S	0	0	0
			2314	1485	383	427	19			

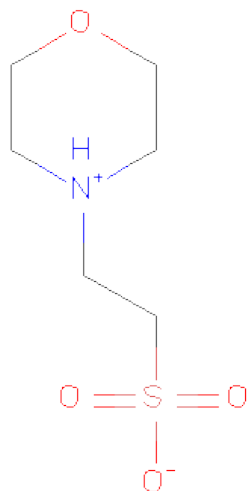
- Molecule 2 is a protein called cyclic peptide S2iL5.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	E	16	Total 144	C 90	F 3	N 27	O 23	S 1	0	0	1
2	F	16	Total 144	C 90	F 3	N 27	O 23	S 1	0	0	1
2	G	16	Total 144	C 90	F 3	N 27	O 23	S 1	0	0	1
2	H	16	Total 144	C 90	F 3	N 27	O 23	S 1	0	0	1

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

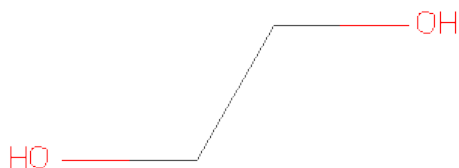
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONICACID (three-letter code: MES) (formula:  $C_6H_{13}NO_4S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	G	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is water.

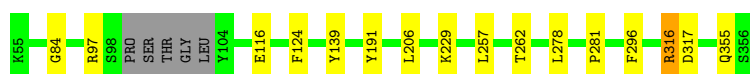
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	132	Total	O	0	0
			132	132		
6	B	103	Total	O	0	0
			103	103		
6	C	122	Total	O	0	0
			122	122		
6	D	96	Total	O	0	0
			96	96		
6	E	17	Total	O	0	0
			17	17		
6	F	15	Total	O	0	0
			15	15		
6	G	12	Total	O	0	0
			12	12		
6	H	10	Total	O	0	0
			10	10		

### 3 Residue-property plots

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

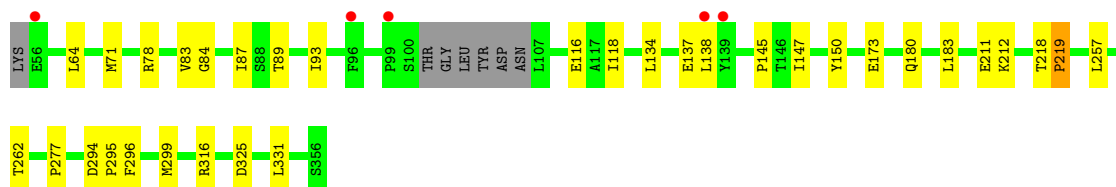
- Molecule 1: NAD-dependent protein deacetylase sirtuin-2

Chain A: 



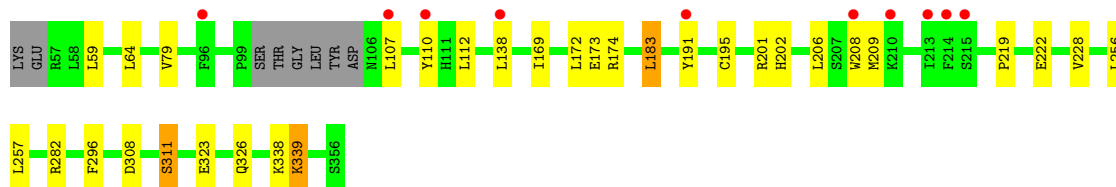
- Molecule 1: NAD-dependent protein deacetylase sirtuin-2

Chain B: 



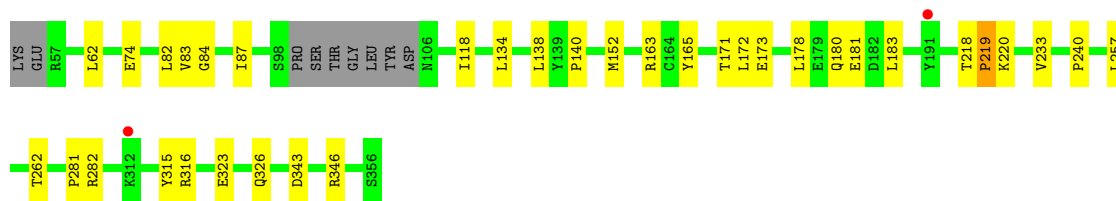
- Molecule 1: NAD-dependent protein deacetylase sirtuin-2

Chain C: 



- Molecule 1: NAD-dependent protein deacetylase sirtuin-2

Chain D: 



- Molecule 2: cyclic peptide S2iL5

Chain E: 



- Molecule 2: cyclic peptide S2iL5

Chain F: 



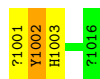
- Molecule 2: cyclic peptide S2iL5

Chain G: 



- Molecule 2: cyclic peptide S2iL5

Chain H: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.18Å 135.61Å 148.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.59 – 2.52 46.59 – 2.52	Depositor EDS
% Data completeness (in resolution range)	97.1 (46.59-2.52) 97.2 (46.59-2.52)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.80 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, $R_{free}$	0.211 , 0.263 0.209 , 0.261	Depositor DCC
$R_{free}$ test set	6134 reflections (10.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.9	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 14.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 61355 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	10482	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, EDO, FAK, MES, NH2, ACY

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	0.22	0/2425	0.37	0/3269
1	B	0.22	0/2383	0.37	0/3215
1	C	0.22	0/2374	0.38	0/3204
1	D	0.22	0/2367	0.37	0/3195
2	E	0.22	0/129	0.42	0/173
2	F	0.22	0/129	0.38	0/173
2	G	0.23	0/129	0.42	0/173
2	H	0.25	0/129	0.42	0/173
All	All	0.22	0/10065	0.37	0/13575

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2368	0	2353	6	0
1	B	2329	0	2314	14	0
1	C	2320	0	2303	14	0
1	D	2314	0	2286	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	144	0	111	0	0
2	F	144	0	111	2	0
2	G	144	0	111	2	0
2	H	144	0	110	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	12	0	12	0	0
4	B	12	0	12	1	0
4	C	12	0	12	0	0
4	D	12	0	12	1	0
5	B	4	0	6	1	0
5	C	8	0	12	0	0
5	G	4	0	6	1	0
6	A	132	0	0	0	0
6	B	103	0	0	0	0
6	C	122	0	0	1	0
6	D	96	0	0	0	0
6	E	17	0	0	0	0
6	F	15	0	0	0	0
6	G	12	0	0	0	0
6	H	10	0	0	0	0
All	All	10482	0	9771	56	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (56) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:343:ASP:OD1	1:D:346:ARG:NH2	2.27	0.68
2:G:1010:ARG:H	5:G:1101:EDO:H22	1.59	0.67
1:C:138:LEU:HD21	1:C:169:ILE:HG13	1.81	0.62
1:D:257:LEU:HB2	1:D:282:ARG:HG2	1.83	0.60
1:B:71:MET:HE3	1:B:331:LEU:HD11	1.84	0.59
1:D:323:GLU:HB2	1:D:326:GLN:HG2	1.85	0.57
1:C:219:PRO:HB2	1:C:228:VAL:HB	1.88	0.56
1:C:308:ASP:OD1	1:C:311:SER:HB3	2.08	0.54
1:C:323:GLU:HB2	1:C:326:GLN:HG2	1.90	0.53
1:C:257:LEU:HB2	1:C:282:ARG:HG2	1.91	0.53
2:H:1001:ACY:C	2:H:1003:HIS:H	2.22	0.53
1:B:277:PRO:O	1:B:316:ARG:NH1	2.43	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:84:GLY:HA3	1:A:262:THR:HB	1.91	0.52
1:B:118:ILE:HD13	1:B:134:LEU:HD13	1.91	0.52
1:C:209:MET:HE1	1:C:219:PRO:HG3	1.91	0.51
1:D:233:VAL:HG21	1:D:240:PRO:HD3	1.93	0.51
1:C:79:VAL:HG22	1:C:256:LEU:HB3	1.93	0.51
2:H:1001:ACY:O	2:H:1003:HIS:N	2.35	0.50
1:A:191:TYR:HA	1:A:206:LEU:HD13	1.93	0.50
1:C:338:LYS:HB3	1:C:339:LYS:HE3	1.92	0.50
1:B:116:GLU:HG3	2:F:1010:ARG:HH21	1.76	0.49
1:C:174:ARG:NH2	6:C:683:HOH:O	2.46	0.48
1:B:147:ILE:HD12	1:B:325:ASP:HB3	1.96	0.48
1:D:62:LEU:HD21	1:D:315:TYR:O	2.14	0.47
2:H:1001:ACY:O	2:H:1003:HIS:ND1	2.47	0.47
1:D:83:VAL:HB	1:D:87:ILE:HD13	1.97	0.47
1:C:195:CYS:HB3	1:C:201:ARG:HA	1.96	0.47
1:D:281:PRO:HA	1:D:316:ARG:HB2	1.98	0.46
1:B:83:VAL:HB	1:B:87:ILE:HD13	1.96	0.46
1:C:191:TYR:HA	1:C:206:LEU:HD12	1.98	0.46
1:D:82:LEU:HD23	1:D:165:TYR:HB2	1.99	0.45
1:D:163:ARG:NH1	1:D:181:GLU:O	2.49	0.44
1:C:110:TYR:HB2	1:C:112:LEU:HG	1.99	0.44
1:B:299:MET:HE1	2:F:1004:THR:H	1.82	0.44
1:B:294:ASP:HA	1:B:295:PRO:HD3	1.89	0.43
1:D:118:ILE:HD13	1:D:134:LEU:HD13	2.00	0.43
2:G:1001:ACY:C	2:G:1003:HIS:H	2.29	0.43
1:B:212:LYS:HD3	1:B:219:PRO:HA	2.01	0.42
1:B:78:ARG:HG2	5:B:503:EDO:H21	2.00	0.42
1:A:97:ARG:NH2	1:A:116:GLU:OE1	2.53	0.42
1:A:281:PRO:HA	1:A:316:ARG:HB2	2.01	0.42
1:B:84:GLY:HA3	1:B:262:THR:HB	2.02	0.42
1:C:173:GLU:HB2	1:C:183:LEU:HD21	2.02	0.42
1:D:173:GLU:O	1:D:178:LEU:HB2	2.20	0.42
1:D:218:THR:HA	1:D:219:PRO:HD3	1.87	0.41
1:B:218:THR:HA	1:B:219:PRO:HD3	1.94	0.41
4:D:502:MES:H51	4:D:502:MES:H81	1.87	0.41
1:A:316:ARG:HD2	1:A:317:ASP:OD1	2.20	0.41
4:B:502:MES:H51	4:B:502:MES:H81	1.74	0.41
1:C:208:TRP:HH2	1:C:222:GLU:HG2	1.86	0.40
1:D:84:GLY:HA3	1:D:262:THR:HB	2.02	0.40
1:D:62:LEU:HD21	1:D:315:TYR:C	2.42	0.40
1:A:124:PHE:CE2	1:A:229:LYS:HB3	2.57	0.40
1:B:89:THR:HA	1:B:93:ILE:O	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:145:PRO:HB2	1:B:150:TYR:CE2	2.56	0.40
1:D:140:PRO:HB3	1:D:171:THR:OG1	2.22	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	294/302 (97%)	288 (98%)	6 (2%)	0	100	100
1	B	291/302 (96%)	286 (98%)	4 (1%)	1 (0%)	50	72
1	C	290/302 (96%)	280 (97%)	10 (3%)	0	100	100
1	D	289/302 (96%)	279 (96%)	9 (3%)	1 (0%)	50	72
2	E	13/16 (81%)	10 (77%)	3 (23%)	0	100	100
2	F	13/16 (81%)	11 (85%)	2 (15%)	0	100	100
2	G	13/16 (81%)	11 (85%)	2 (15%)	0	100	100
2	H	13/16 (81%)	11 (85%)	1 (8%)	1 (8%)	1	1
All	All	1216/1272 (96%)	1176 (97%)	37 (3%)	3 (0%)	56	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	1002	TYR
1	D	219	PRO
1	B	219	PRO

### 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	259/263 (98%)	253 (98%)	6 (2%)	63	87
1	B	253/263 (96%)	244 (96%)	9 (4%)	47	72
1	C	252/263 (96%)	243 (96%)	9 (4%)	47	72
1	D	252/263 (96%)	245 (97%)	7 (3%)	56	82
2	E	13/13 (100%)	12 (92%)	1 (8%)	18	32
2	F	13/13 (100%)	13 (100%)	0	100	100
2	G	13/13 (100%)	12 (92%)	1 (8%)	18	32
2	H	13/13 (100%)	12 (92%)	1 (8%)	18	32
All	All	1068/1104 (97%)	1034 (97%)	34 (3%)	51	77

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

Mol	Chain	Res	Type
1	A	139	TYR
1	A	257	LEU
1	A	278	LEU
1	A	296	PHE
1	A	316	ARG
1	A	355	GLN
1	B	64	LEU
1	B	137	GLU
1	B	138	LEU
1	B	173	GLU
1	B	180	GLN
1	B	183	LEU
1	B	211	GLU
1	B	257	LEU
1	B	296	PHE
1	C	59	LEU
1	C	64	LEU
1	C	107	LEU
1	C	172	LEU
1	C	183	LEU
1	C	202	HIS
1	C	296	PHE
1	C	311	SER
1	C	339	LYS
1	D	74	GLU
1	D	138	LEU
1	D	152	MET

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Mol	Chain	Res	Type
1	D	172	LEU
1	D	180	GLN
1	D	183	LEU
1	D	220	LYS
2	E	1002	TYR
2	G	1002	TYR
2	H	1002	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	167	GLN
1	A	225	GLN
1	A	326	GLN
1	B	194	HIS
1	B	202	HIS
1	C	142	GLN
1	C	355	GLN
1	D	167	GLN
1	D	180	GLN
1	D	202	HIS
1	D	326	GLN
2	F	1006	HIS
2	G	1006	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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	FAK	E	1008	2	14,14,15	4.81	3 (21%)	16,18,20	2.06	2 (12%)
2	FAK	F	1008	2	14,14,15	4.79	3 (21%)	16,18,20	1.61	2 (12%)
2	FAK	G	1008	2	14,14,15	4.85	3 (21%)	16,18,20	1.91	2 (12%)
2	FAK	H	1008	2	14,14,15	4.85	3 (21%)	16,18,20	1.75	2 (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	FAK	E	1008	2	-	0/13/16/18	0/0/0/0
2	FAK	F	1008	2	-	0/13/16/18	0/0/0/0
2	FAK	G	1008	2	-	0/13/16/18	0/0/0/0
2	FAK	H	1008	2	-	0/13/16/18	0/0/0/0

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	1008	FAK	O-C	16.74	1.22	1.11
2	G	1008	FAK	O-C	16.73	1.22	1.11
2	E	1008	FAK	O-C	16.73	1.22	1.11
2	F	1008	FAK	O-C	16.58	1.22	1.11
2	H	1008	FAK	CB-CA	-5.13	1.49	1.53
2	G	1008	FAK	CB-CA	-5.06	1.49	1.53
2	F	1008	FAK	CB-CA	-4.86	1.49	1.53
2	E	1008	FAK	CB-CA	-4.63	1.49	1.53
2	E	1008	FAK	CH-NZ	4.20	1.43	1.34
2	F	1008	FAK	CH-NZ	4.12	1.43	1.34
2	G	1008	FAK	CH-NZ	4.10	1.43	1.34
2	H	1008	FAK	CH-NZ	4.07	1.43	1.34

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1008	FAK	CT-CH-NZ	6.56	120.63	115.32
2	G	1008	FAK	CT-CH-NZ	5.98	120.15	115.32
2	F	1008	FAK	CT-CH-NZ	5.00	119.36	115.32
2	H	1008	FAK	CT-CH-NZ	4.92	119.30	115.32
2	H	1008	FAK	C-CA-N	-3.85	109.98	113.83
2	E	1008	FAK	C-CA-N	-3.37	110.46	113.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	1008	FAK	C-CA-N	-3.09	110.74	113.83
2	F	1008	FAK	C-CA-N	-2.52	111.31	113.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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$
4	MES	A	502	-	12,12,12	2.18	1 (8%)	16,16,16	2.24	6 (37%)
4	MES	B	502	-	12,12,12	2.11	1 (8%)	16,16,16	1.96	3 (18%)
5	EDO	B	503	-	3,3,3	0.54	0	2,2,2	0.32	0
4	MES	C	502	-	12,12,12	2.17	1 (8%)	16,16,16	2.40	6 (37%)
5	EDO	C	503	-	3,3,3	0.55	0	2,2,2	0.31	0
5	EDO	C	504	-	3,3,3	0.56	0	2,2,2	0.33	0
4	MES	D	502	-	12,12,12	2.14	1 (8%)	16,16,16	2.07	4 (25%)
5	EDO	G	1101	-	3,3,3	0.53	0	2,2,2	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
4	MES	A	502	-	-	0/6/14/14	0/1/1/1
4	MES	B	502	-	-	0/6/14/14	0/1/1/1
5	EDO	B	503	-	-	0/1/1/1	0/0/0/0
4	MES	C	502	-	-	0/6/14/14	0/1/1/1
5	EDO	C	503	-	-	0/1/1/1	0/0/0/0
5	EDO	C	504	-	-	0/1/1/1	0/0/0/0
4	MES	D	502	-	-	0/6/14/14	0/1/1/1
5	EDO	G	1101	-	-	0/1/1/1	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	502	MES	C8-S	-6.59	1.66	1.78
4	C	502	MES	C8-S	-6.54	1.66	1.78
4	D	502	MES	C8-S	-6.48	1.66	1.78
4	B	502	MES	C8-S	-6.41	1.66	1.78

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	502	MES	C8-C7-N4	-5.15	103.63	112.44
4	A	502	MES	C8-C7-N4	-4.95	103.96	112.44
4	A	502	MES	C5-N4-C3	4.69	119.05	109.75
4	C	502	MES	C5-N4-C3	4.33	118.34	109.75
4	D	502	MES	C5-N4-C3	4.29	118.26	109.75
4	C	502	MES	C6-C5-N4	-4.27	104.49	109.96
4	B	502	MES	C5-N4-C3	4.08	117.84	109.75
4	D	502	MES	C8-C7-N4	-4.01	105.57	112.44
4	B	502	MES	C8-C7-N4	-3.73	106.06	112.44
4	D	502	MES	C6-C5-N4	-3.43	105.56	109.96
4	A	502	MES	C6-C5-N4	-3.28	105.75	109.96
4	B	502	MES	C6-C5-N4	-3.25	105.79	109.96
4	C	502	MES	C2-C3-N4	-2.56	106.67	109.96
4	C	502	MES	C7-N4-C3	2.49	118.17	111.66
4	C	502	MES	C7-N4-C5	2.39	117.89	111.66
4	A	502	MES	C7-N4-C3	2.12	117.19	111.66
4	A	502	MES	C7-N4-C5	2.10	117.14	111.66
4	A	502	MES	C2-C3-N4	-2.06	107.31	109.96
4	D	502	MES	C7-N4-C5	2.04	116.99	111.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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, 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	297/302 (98%)	-0.35	0 100 100	13, 25, 58, 71	0
1	B	295/302 (97%)	-0.08	5 (1%) 67 70	15, 38, 70, 101	0
1	C	294/302 (97%)	-0.15	10 (3%) 43 45	12, 29, 73, 91	0
1	D	293/302 (97%)	-0.10	2 (0%) 84 87	17, 42, 65, 82	0
2	E	15/16 (93%)	-0.36	0 100 100	15, 21, 26, 27	0
2	F	15/16 (93%)	-0.24	0 100 100	16, 20, 26, 26	0
2	G	15/16 (93%)	-0.22	0 100 100	15, 23, 29, 29	0
2	H	15/16 (93%)	-0.11	0 100 100	17, 26, 34, 37	0
All	All	1239/1272 (97%)	-0.17	17 (1%) 72 75	12, 32, 67, 101	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	99	PRO	4.3
1	C	110	TYR	4.0
1	C	214	PHE	3.4
1	C	107	LEU	3.2
1	C	138	LEU	3.0
1	C	215	SER	2.9
1	C	210	LYS	2.6
1	C	96	PHE	2.6
1	C	191	TYR	2.5
1	B	138	LEU	2.5
1	B	96	PHE	2.5
1	B	56	GLU	2.4
1	C	208	TRP	2.3
1	D	312	LYS	2.2
1	C	213	ILE	2.2
1	B	139	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	191	TYR	2.1

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

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	FAK	G	1008	15/16	0.14	0.17	11,20,27,29	0
2	FAK	F	1008	15/16	0.13	-0.03	15,18,32,39	0
2	FAK	E	1008	15/16	0.12	-0.29	10,15,23,30	0
2	FAK	H	1008	15/16	0.11	-1.00	16,23,36,37	0

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	EDO	C	504	4/4	0.38	12.47	30,31,40,42	0
4	MES	D	502	12/12	0.21	3.51	47,55,67,74	0
4	MES	A	502	12/12	0.19	2.85	35,59,64,67	0
4	MES	B	502	12/12	0.19	2.42	38,58,68,72	0
5	EDO	C	503	4/4	0.16	0.90	24,28,29,38	0
4	MES	C	502	12/12	0.14	0.73	40,52,63,68	0
5	EDO	G	1101	4/4	0.14	0.38	27,33,34,37	0
5	EDO	B	503	4/4	0.15	0.29	35,36,42,42	0
3	ZN	C	501	1/1	0.08	-2.08	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZN	D	501	1/1	0.05	-2.33	58,58,58,58	0
3	ZN	B	501	1/1	0.08	-2.57	58,58,58,58	0
3	ZN	A	501	1/1	0.04	-2.71	66,66,66,66	0

## 6.5 Other polymers

There are no such residues in this entry.