



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 7, 2024 – 08:23 am GMT

PDB ID : 5OKF
Title : CH1 chimera of human 14-3-3 sigma with the HSPB6 phosphopeptide in a conformation with self-bound phosphopeptides
Authors : Sluchanko, N.N.; Tugaeva, K.V.; Greive, S.J.; Antson, A.A.
Deposited on : 2017-07-25
Resolution : 3.20 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 : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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

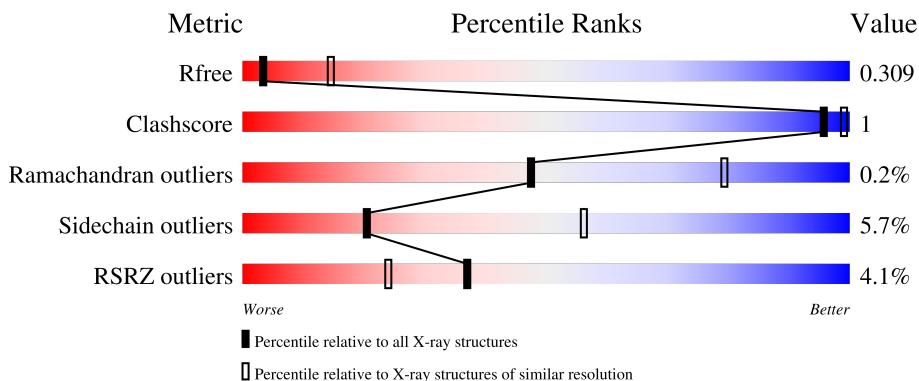
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	246	 5% 87% 5% • 6%
1	B	246	 3% 87% 6% 7%
1	C	246	 4% 83% 11% 6%
1	D	246	 2% 87% 7% 7%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 14370 atoms, of which 7043 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,Heat shock protein beta-6.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
			Total	C	H	N	O	P	S			
1	A	231	3568	1137	1744	314	362	1	10	0	0	0
1	B	228	3569	1129	1759	313	357	1	10	0	0	0
1	C	231	3608	1143	1777	316	361	1	10	0	0	0
1	D	230	3586	1139	1763	315	358	1	10	0	0	0

There are 40 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
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
C	-2	GLY	-	expression tag	UNP P31947

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	5	Total Cd 5 5	0	0
2	B	6	Total Cd 6 6	0	0
2	C	2	Total Cd 2 2	0	0
2	D	4	Total Cd 4 4	0	0

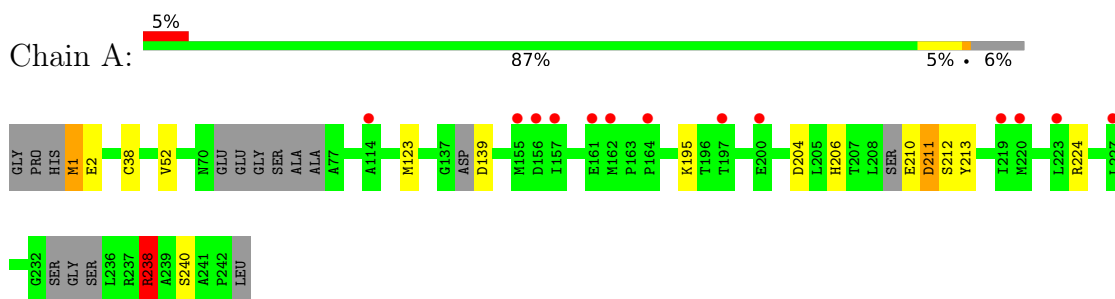
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	5	Total O 5 5	0	0
3	B	9	Total O 9 9	0	0
3	C	6	Total O 6 6	0	0
3	D	2	Total O 2 2	0	0

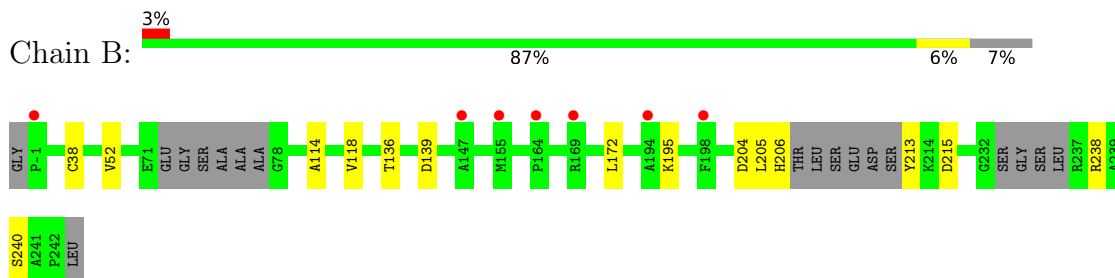
3 Residue-property plots [i](#)

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.

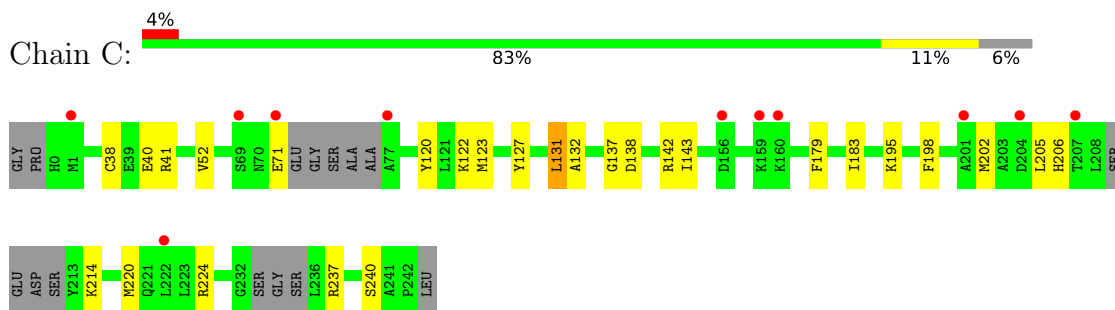
- Molecule 1: 14-3-3 protein sigma,Heat shock protein beta-6



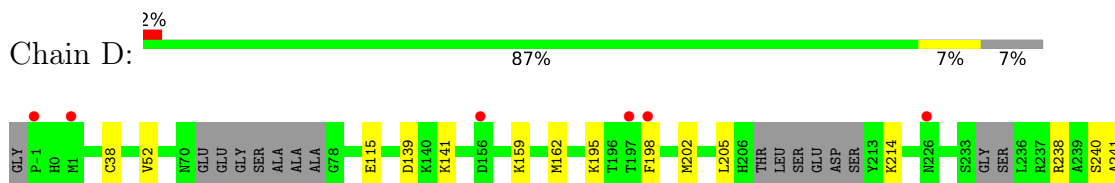
- Molecule 1: 14-3-3 protein sigma,Heat shock protein beta-6



- Molecule 1: 14-3-3 protein sigma,Heat shock protein beta-6



- Molecule 1: 14-3-3 protein sigma,Heat shock protein beta-6



P242
L243

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.40Å 97.76Å 158.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.00 – 3.20 48.22 – 3.13	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.00-3.20) 99.0 (48.22-3.13)	Depositor EDS
R_{merge}	0.46	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 3.12Å)	Xtrriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.247 , 0.279 0.278 , 0.309	Depositor DCC
R_{free} test set	1075 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	51.7	Xtrriage
Anisotropy	0.477	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 68.9	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	0.88	EDS
Total number of atoms	14370	wwPDB-VP
Average B, all atoms (Å ²)	67.0	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 55.44 % 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 3.1636e-05. 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: CD, SEP, CSO

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	0/1828	0.58	0/2451
1	B	0.51	0/1817	0.62	0/2437
1	C	0.50	0/1837	0.63	0/2465
1	D	0.52	0/1830	0.61	0/2455
All	All	0.50	0/7312	0.61	0/9808

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	1824	1744	1804	6	0
1	B	1810	1759	1788	1	0
1	C	1831	1777	1816	6	1
1	D	1823	1763	1809	3	1
2	A	5	0	0	0	0
2	B	6	0	0	0	0
2	C	2	0	0	0	0
2	D	4	0	0	0	0
3	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	9	0	0	1	0
3	C	6	0	0	0	0
3	D	2	0	0	0	0
All	All	7327	7043	7217	16	1

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 1.

All (16) 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:1:MET:HG3	1:A:2:GLU:H	1.03	1.04
1:A:1:MET:HG3	1:A:2:GLU:N	1.79	0.97
1:A:1:MET:CG	1:A:2:GLU:N	2.57	0.62
1:A:1:MET:CG	1:A:2:GLU:H	1.93	0.54
1:D:115:GLU:HA	1:D:162:MET:CE	2.38	0.54
1:B:114:ALA:O	1:B:118:VAL:HG23	2.12	0.50
3:B:404:HOH:O	1:C:142:ARG:HB3	2.15	0.47
1:C:132:ALA:HB2	1:C:143:ILE:HG22	1.97	0.46
1:D:241:ALA:HB1	1:D:243:LEU:HB2	1.98	0.46
1:D:198:PHE:CE1	1:D:202:MET:CE	3.00	0.45
1:C:198:PHE:CE1	1:C:202:MET:CE	3.00	0.44
1:C:127:TYR:O	1:C:131:LEU:HD12	2.18	0.43
1:A:210:GLU:HG3	1:A:211:ASP:H	1.84	0.42
1:C:179:PHE:CE2	1:C:183:ILE:HG21	2.55	0.41
1:C:41:ARG:HD2	1:C:120:TYR:CE1	2.55	0.41
1:A:238:ARG:C	1:A:238:ARG:HD2	2.42	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:40:GLU:OE2	1:D:141:LYS:NZ[3_444]	1.95	0.25

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	219/246 (89%)	211 (96%)	7 (3%)	1 (0%)	29	67
1	B	218/246 (89%)	211 (97%)	7 (3%)	0	100	100
1	C	221/246 (90%)	210 (95%)	10 (4%)	1 (0%)	29	67
1	D	220/246 (89%)	210 (96%)	10 (4%)	0	100	100
All	All	878/984 (89%)	842 (96%)	34 (4%)	2 (0%)	47	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	137	GLY
1	A	238	ARG

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	190/201 (94%)	178 (94%)	12 (6%)	18	52
1	B	189/201 (94%)	178 (94%)	11 (6%)	20	55
1	C	191/201 (95%)	178 (93%)	13 (7%)	16	49
1	D	191/201 (95%)	184 (96%)	7 (4%)	34	68
All	All	761/804 (95%)	718 (94%)	43 (6%)	20	56

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	52	VAL
1	A	123	MET
1	A	139	ASP
1	A	195	LYS
1	A	204	ASP
1	A	206	HIS
1	A	211	ASP
1	A	212	SER
1	A	213	TYR
1	A	224	ARG
1	A	238	ARG
1	B	52	VAL
1	B	136	THR
1	B	139	ASP
1	B	172	LEU
1	B	195	LYS
1	B	204	ASP
1	B	205	LEU
1	B	206	HIS
1	B	213	TYR
1	B	215	ASP
1	B	238	ARG
1	C	52	VAL
1	C	71	GLU
1	C	122	LYS
1	C	123	MET
1	C	131	LEU
1	C	138	ASP
1	C	195	LYS
1	C	205	LEU
1	C	206	HIS
1	C	214	LYS
1	C	220	MET
1	C	224	ARG
1	C	237	ARG
1	D	52	VAL
1	D	139	ASP
1	D	159	LYS
1	D	195	LYS
1	D	205	LEU
1	D	214	LYS
1	D	238	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	206	HIS
1	D	206	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

8 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
1	CSO	C	38	1	3,6,7	1.82	1 (33%)	0,6,8	-	-
1	CSO	B	38	1	3,6,7	4.60	1 (33%)	0,6,8	-	-
1	CSO	A	38	1	3,6,7	2.05	1 (33%)	0,6,8	-	-
1	SEP	B	240	1	8,9,10	1.22	1 (12%)	8,12,14	1.78	2 (25%)
1	SEP	A	240	1	8,9,10	0.76	0	8,12,14	2.10	2 (25%)
1	CSO	D	38	2,1	3,6,7	3.67	1 (33%)	0,6,8	-	-
1	SEP	D	240	1	8,9,10	1.09	1 (12%)	8,12,14	1.83	2 (25%)
1	SEP	C	240	1	8,9,10	1.28	2 (25%)	8,12,14	1.18	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
1	CSO	C	38	1	-	1/1/5/7	-
1	CSO	B	38	1	-	0/1/5/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSO	A	38	1	-	1/1/5/7	-
1	SEP	B	240	1	-	1/5/8/10	-
1	SEP	A	240	1	-	1/5/8/10	-
1	CSO	D	38	2,1	-	1/1/5/7	-
1	SEP	D	240	1	-	0/5/8/10	-
1	SEP	C	240	1	-	0/5/8/10	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	38	CSO	CB-CA	7.86	1.72	1.53
1	D	38	CSO	CB-CA	6.25	1.68	1.53
1	A	38	CSO	CB-CA	3.36	1.61	1.53
1	C	38	CSO	CB-CA	2.98	1.60	1.53
1	B	240	SEP	P-OG	-2.51	1.52	1.60
1	C	240	SEP	P-O1P	2.22	1.57	1.50
1	D	240	SEP	P-O3P	-2.19	1.46	1.54
1	C	240	SEP	P-OG	-2.10	1.53	1.60

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	240	SEP	O2P-P-OG	3.69	116.55	106.73
1	A	240	SEP	O3P-P-OG	3.63	116.40	106.73
1	B	240	SEP	O3P-P-OG	3.08	114.93	106.73
1	B	240	SEP	O3P-P-O2P	3.06	119.34	107.64
1	D	240	SEP	O3P-P-O1P	-3.00	98.95	110.68
1	A	240	SEP	O3P-P-O2P	2.89	118.67	107.64
1	C	240	SEP	O3P-P-O2P	2.53	117.29	107.64

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	38	CSO	N-CA-CB-SG
1	C	38	CSO	N-CA-CB-SG
1	D	38	CSO	N-CA-CB-SG
1	A	240	SEP	CB-OG-P-O1P
1	B	240	SEP	CB-OG-P-O1P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 17 are monoatomic - leaving 0 for Mogul analysis.

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	229/246 (93%)	0.43	13 (5%) 23 13	40, 73, 103, 114	0
1	B	226/246 (91%)	0.31	7 (3%) 49 32	40, 66, 103, 124	0
1	C	229/246 (93%)	0.29	11 (4%) 30 18	27, 64, 98, 131	0
1	D	228/246 (92%)	0.22	6 (2%) 56 40	25, 58, 99, 112	0
All	All	912/984 (92%)	0.31	37 (4%) 37 24	25, 66, 102, 131	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	71	GLU	4.1
1	A	164	PRO	3.8
1	C	207	THR	3.6
1	B	194	ALA	3.5
1	A	156	ASP	3.2
1	B	164	PRO	3.1
1	A	157	ILE	3.1
1	B	155	MET	3.1
1	C	77	ALA	2.9
1	D	-1	PRO	2.9
1	C	160	LYS	2.9
1	A	155	MET	2.8
1	C	159	LYS	2.8
1	D	197	THR	2.7
1	A	227	LEU	2.6
1	A	200	GLU	2.5
1	C	204	ASP	2.5
1	C	1	MET	2.4
1	B	169	ARG	2.4
1	C	69	SER	2.4
1	A	219	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	156	ASP	2.4
1	A	220	MET	2.3
1	D	156	ASP	2.3
1	A	162	MET	2.2
1	A	114	ALA	2.2
1	A	223	LEU	2.2
1	D	226	ASN	2.2
1	D	198	PHE	2.2
1	D	1	MET	2.1
1	A	161	GLU	2.1
1	B	147	ALA	2.1
1	C	201	ALA	2.1
1	B	198	PHE	2.1
1	B	-1	PRO	2.1
1	A	197	THR	2.0
1	C	222	LEU	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
1	CSO	A	38	7/8	0.76	0.26	81,83,84,84	0
1	CSO	B	38	7/8	0.81	0.27	46,49,58,61	0
1	CSO	C	38	7/8	0.84	0.27	60,61,68,70	0
1	CSO	D	38	7/8	0.92	0.20	55,57,60,61	0
1	SEP	C	240	10/11	0.93	0.17	36,39,47,47	0
1	SEP	B	240	10/11	0.95	0.16	45,49,52,52	0
1	SEP	A	240	10/11	0.96	0.16	49,52,55,55	0
1	SEP	D	240	10/11	0.97	0.15	50,51,54,54	0

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
2	CD	B	305	1/1	0.30	0.24	182,182,182,182	0
2	CD	C	301	1/1	0.53	0.16	167,167,167,167	0
2	CD	B	303	1/1	0.75	0.22	157,157,157,157	0
2	CD	A	304	1/1	0.77	0.22	161,161,161,161	0
2	CD	D	301	1/1	0.77	0.27	152,152,152,152	0
2	CD	A	303	1/1	0.81	0.11	178,178,178,178	0
2	CD	A	305	1/1	0.84	0.15	115,115,115,115	0
2	CD	D	304	1/1	0.84	0.12	189,189,189,189	0
2	CD	C	302	1/1	0.89	0.28	186,186,186,186	0
2	CD	B	301	1/1	0.92	0.10	149,149,149,149	0
2	CD	D	302	1/1	0.94	0.12	109,109,109,109	0
2	CD	D	303	1/1	0.95	0.18	170,170,170,170	0
2	CD	B	302	1/1	0.95	0.27	110,110,110,110	0
2	CD	A	302	1/1	0.96	0.16	97,97,97,97	0
2	CD	A	301	1/1	0.97	0.15	134,134,134,134	0
2	CD	B	306	1/1	0.97	0.12	87,87,87,87	0
2	CD	B	304	1/1	0.98	0.24	107,107,107,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.