



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 16, 2021 – 07:45 PM EDT

PDB ID : 1RDF  
Title : G50P mutant of phosphonoacetaldehyde hydrolase in complex with substrate analogue vinyl sulfonate  
Authors : Lahiri, S.D.; Zhang, G.; Dunaway-Mariano, D.; Allen, K.N.  
Deposited on : 2003-11-05  
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
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)	:	2.23.2

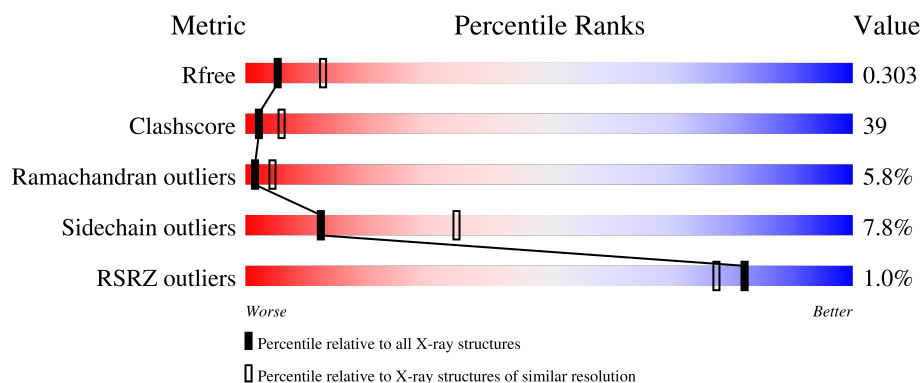
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

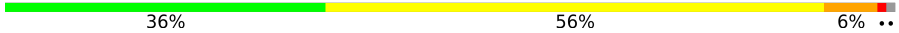
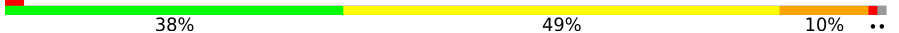
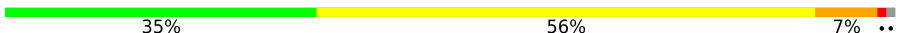


The reported resolution of this entry is 2.80 Å.

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))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	267	
1	B	267	
1	C	267	
1	D	267	
1	E	267	

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Mol	Chain	Length	Quality of chain
1	F	267	

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
3	ESA	D	500	-	X	-	-
3	ESA	E	500	-	X	-	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12927 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 phosphonoacetaldehyde hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2101	1337	348	399	17			
1	B	263	Total	C	N	O	S	0	0	0
			2098	1335	348	398	17			
1	C	263	Total	C	N	O	S	0	0	0
			2100	1337	348	398	17			
1	D	263	Total	C	N	O	S	0	0	0
			2100	1337	348	398	17			
1	E	263	Total	C	N	O	S	0	0	0
			2100	1337	348	398	17			
1	F	263	Total	C	N	O	S	0	0	0
			2100	1337	348	398	17			

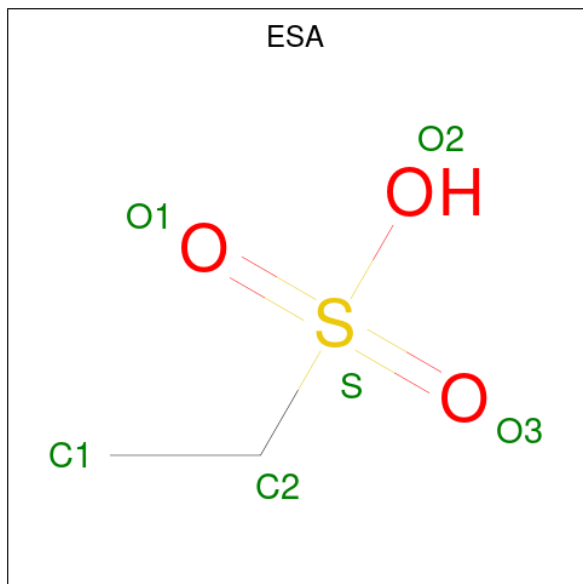
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	PRO	GLY	engineered mutation	UNP O31156
A	107	ALA	GLY	conflict	UNP O31156
B	50	PRO	GLY	engineered mutation	UNP O31156
B	107	ALA	GLY	conflict	UNP O31156
C	50	PRO	GLY	engineered mutation	UNP O31156
C	107	ALA	GLY	conflict	UNP O31156
D	50	PRO	GLY	engineered mutation	UNP O31156
D	107	ALA	GLY	conflict	UNP O31156
E	50	PRO	GLY	engineered mutation	UNP O31156
E	107	ALA	GLY	conflict	UNP O31156
F	50	PRO	GLY	engineered mutation	UNP O31156
F	107	ALA	GLY	conflict	UNP O31156

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	E	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0

- Molecule 3 is ETHANESULFONIC ACID (three-letter code: ESA) (formula:  $C_2H_6O_3S$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O S 6 2 3 1	0	0
3	B	1	Total C O S 6 2 3 1	0	0
3	C	1	Total C O S 6 2 3 1	0	0
3	D	1	Total C O S 6 2 3 1	0	0
3	E	1	Total C O S 6 2 3 1	0	0
3	F	1	Total C O S 6 2 3 1	0	0

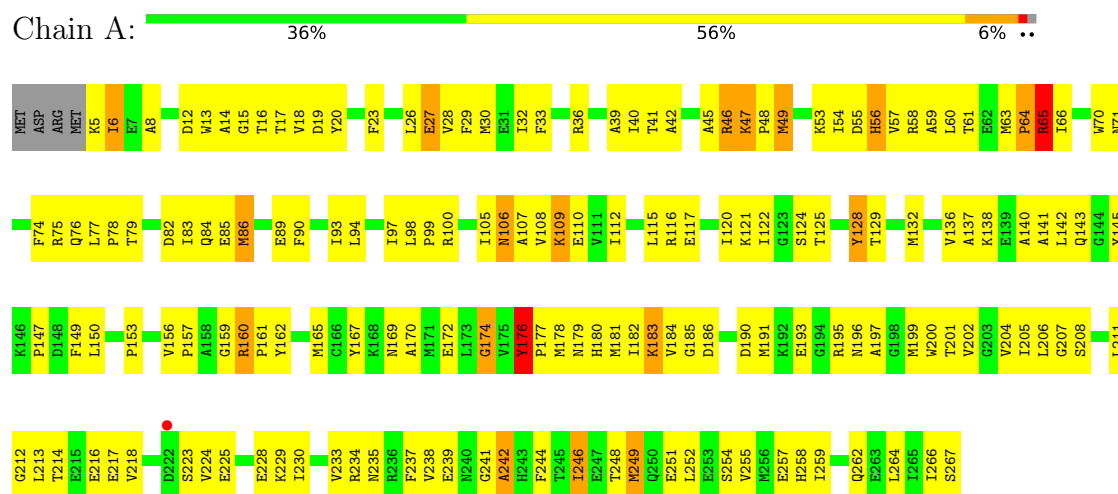
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	52	Total 52	O 52	0	0
4	B	57	Total 57	O 57	0	0
4	C	43	Total 43	O 43	0	0
4	D	50	Total 50	O 50	0	0
4	E	49	Total 49	O 49	0	0
4	F	35	Total 35	O 35	0	0

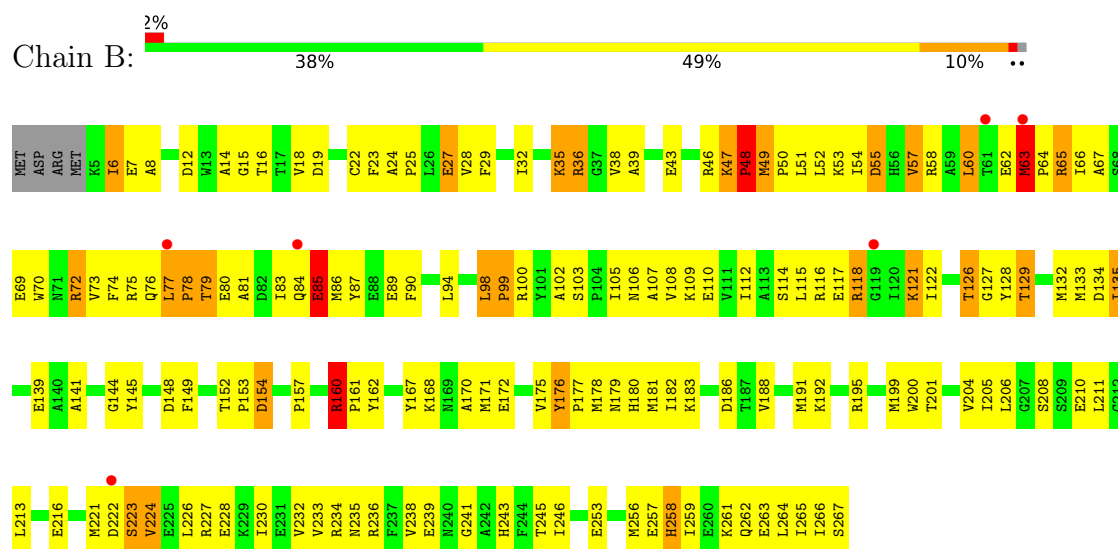
### 3 Residue-property plots

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

#### • Molecule 1: phosphonoacetaldehyde hydrolase

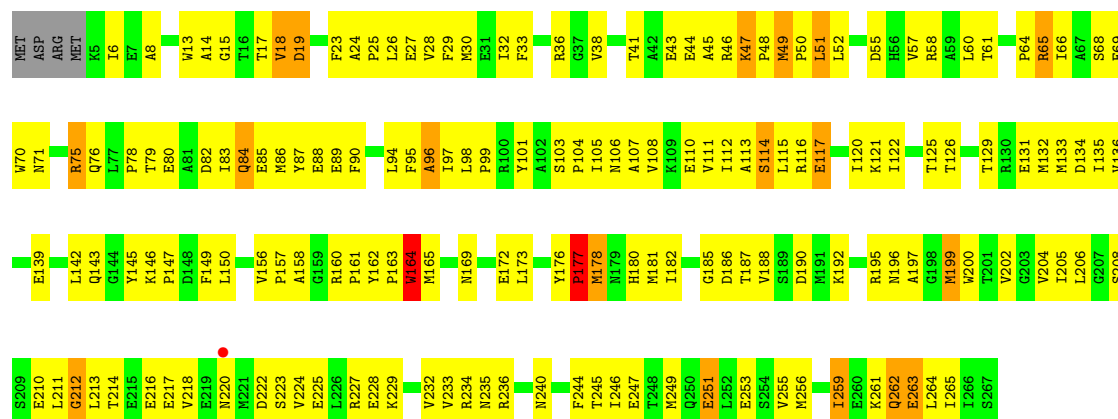


#### • Molecule 1: phosphonoacetaldehyde hydrolase

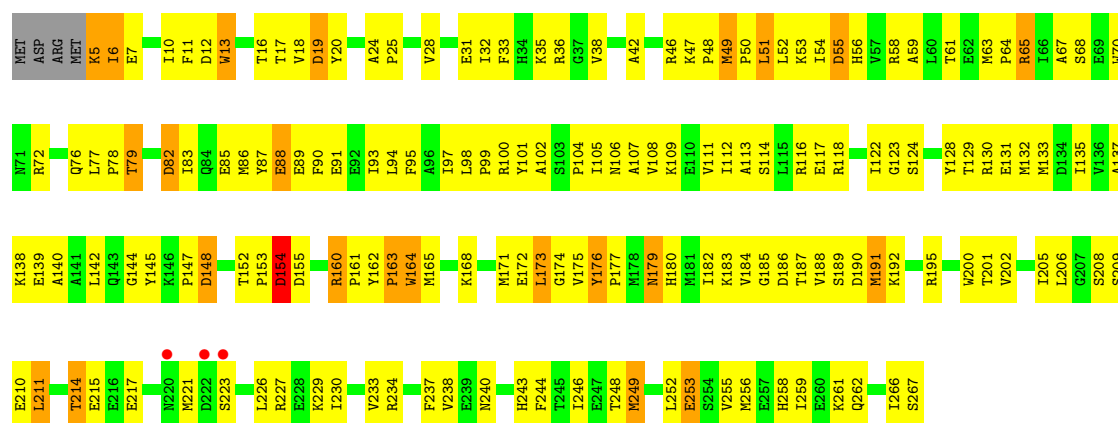


#### • Molecule 1: phosphonoacetaldehyde hydrolase

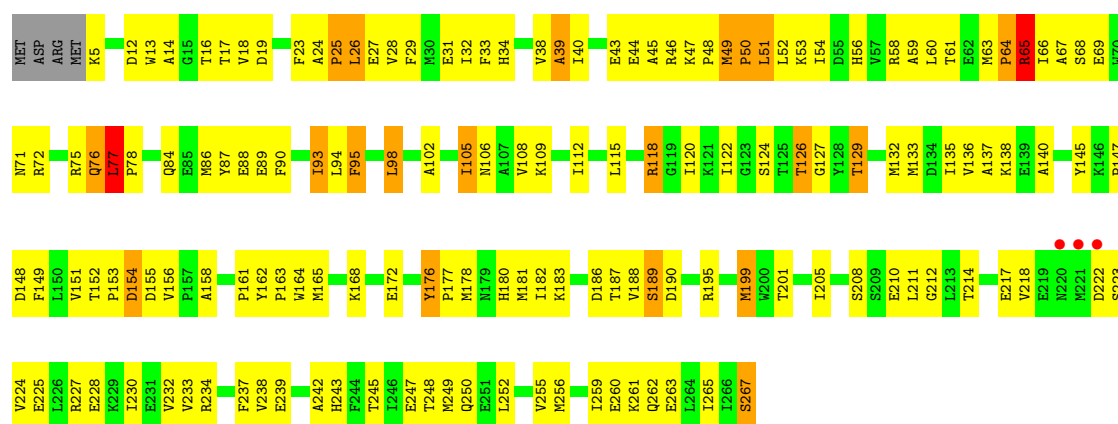




• Molecule 1: phosphonoacetaldehyde hydrolase



• Molecule 1: phosphonoacetaldehyde hydrolase



• Molecule 1: phosphonoacetaldehyde hydrolase





E219	N220	N221	D222	S223	V224	E225	L226	R227	E228	E231	V232	V233	R234	E239	A242	H243	F244	E247	T248	M249	Q250	E251	L252	E253	S254	V255	M256	E257	H258	I259	E260	K261	Q262	E263	L264	I265	I266	S267																
●		●																																																				
D148	F149	L150	V151	T152	P153	V156	P157	A158	G159	R160	P161	Y162	P163	W164	Y167	M171	E172	L173	G174	V175	Y176	P177	H180	M181	I182	K183	V184	G185	D186	T187	V188	S189	D190	M191	K192	E193	G194	M199	W200	T201	V202	G203	V204	T205	L206	G207	E210	T214	E215	E216	E217	V218		
W70	N71	F74	R75	Q76	L77	P78	I83	E89	E92	I93	L94	F95	A96	I97	L98	P99	I105	N106	A107	V108	K109	E110	V111	I112	L115	R116	E117	R118	K121	I122	T125	T126	G127	Y128	E131	M132	M133	D134	I135	V136	E139	A140	A141	I142	Q143	G144	Y145	K146	P147					
MET	ASP	ARG	MET	K5	I6	I10	F11	D12	W13	A14	G15	T16	I17	V18	D19	Y20	G21	C22	F23	A24	P25	L26	E27	V28	E31	I32	F33	R36	G37	V38	E43	E44	A45	R46	K47	P48	M49	P50	L51	I54	D55	H56	V57	R58	A59	L60	T61	E62	M63	P64	R65	I66	A67	

## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	181.71Å 147.35Å 131.31Å 90.00° 125.21° 90.00°	Depositor
Resolution (Å)	34.09 – 2.80 34.09 – 2.70	Depositor EDS
% Data completeness (in resolution range)	77.2 (34.09-2.80) 73.9 (34.09-2.70)	Depositor EDS
$R_{merge}$	0.29	Depositor
$R_{sym}$	0.28	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.91 (at 2.68Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.230 , 0.260 0.245 , 0.303	Depositor DCC
$R_{free}$ test set	5437 reflections (9.49%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.2	Xtriage
Anisotropy	0.208	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 46.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.86	EDS
Total number of atoms	12927	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.25 % 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 7.6614e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: MG, ESA

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	0.41	0/2144	0.69	2/2898 (0.1%)
1	B	0.42	0/2140	0.73	2/2893 (0.1%)
1	C	0.42	0/2143	0.80	3/2898 (0.1%)
1	D	0.42	0/2143	0.68	0/2898
1	E	0.43	0/2143	0.63	0/2898
1	F	0.42	0/2143	0.65	0/2898
All	All	0.42	0/12856	0.70	7/17383 (0.0%)

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	176	TYR	C-N-CD	-22.46	71.18	120.60
1	B	48	PRO	CA-N-CD	-11.61	95.25	111.50
1	C	176	TYR	C-N-CA	10.42	165.77	122.00
1	B	47	LYS	C-N-CD	-5.47	108.56	120.60
1	C	177	PRO	CA-N-CD	-5.39	103.95	111.50

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	2101	0	2084	165	0
1	B	2098	0	2078	174	0
1	C	2100	0	2084	182	0
1	D	2100	0	2084	162	0
1	E	2100	0	2084	161	0
1	F	2100	0	2084	161	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	6	0	3	0	0
3	B	6	0	3	1	0
3	C	6	0	3	3	0
3	D	6	0	3	0	0
3	E	6	0	3	1	0
3	F	6	0	3	1	0
4	A	52	0	0	20	0
4	B	57	0	0	23	0
4	C	43	0	0	17	0
4	D	50	0	0	20	0
4	E	49	0	0	24	0
4	F	35	0	0	6	0
All	All	12927	0	12516	992	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 39.

The worst 5 of 992 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:B:176:TYR:HB3	1:B:177:PRO:HD3	1.24	1.13
1:B:205:ILE:HG21	1:B:230:ILE:HG23	1.41	1.03
1:F:65:ARG:HG2	1:F:66:ILE:H	1.23	1.01
1:E:49:MET:HB3	1:E:50:PRO:HD3	1.44	1.00
1:B:176:TYR:HB3	1:B:177:PRO:CD	1.92	0.98

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	261/267 (98%)	196 (75%)	46 (18%)	19 (7%)	1	2
1	B	261/267 (98%)	201 (77%)	40 (15%)	20 (8%)	1	2
1	C	261/267 (98%)	198 (76%)	48 (18%)	15 (6%)	1	5
1	D	261/267 (98%)	211 (81%)	39 (15%)	11 (4%)	3	9
1	E	261/267 (98%)	208 (80%)	41 (16%)	12 (5%)	2	7
1	F	261/267 (98%)	205 (78%)	42 (16%)	14 (5%)	2	5
All	All	1566/1602 (98%)	1219 (78%)	256 (16%)	91 (6%)	1	4

5 of 91 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	ARG
1	A	47	LYS
1	A	56	HIS
1	A	64	PRO
1	A	128	TYR

### 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	227/231 (98%)	213 (94%)	14 (6%)	18	47
1	B	226/231 (98%)	203 (90%)	23 (10%)	7	22
1	C	227/231 (98%)	213 (94%)	14 (6%)	18	47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	227/231 (98%)	210 (92%)	17 (8%)	13	37
1	E	227/231 (98%)	209 (92%)	18 (8%)	12	34
1	F	227/231 (98%)	207 (91%)	20 (9%)	10	29
All	All	1361/1386 (98%)	1255 (92%)	106 (8%)	12	35

5 of 106 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	88	GLU
1	E	93	ILE
1	F	164	TRP
1	D	114	SER
1	D	191	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 24 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	243	HIS
1	F	84	GLN
1	F	71	ASN
1	F	143	GLN
1	B	240	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

Of 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 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
3	ESA	B	500	2	5,5,5	3.27	1 (20%)	7,7,7	2.34	3 (42%)
3	ESA	A	500	2	5,5,5	3.24	1 (20%)	7,7,7	2.34	2 (28%)
3	ESA	E	500	2	5,5,5	3.17	1 (20%)	7,7,7	2.32	3 (42%)
3	ESA	C	500	2	5,5,5	3.08	1 (20%)	7,7,7	2.30	3 (42%)
3	ESA	F	500	2	5,5,5	3.21	1 (20%)	7,7,7	2.47	2 (28%)
3	ESA	D	500	2	5,5,5	3.01	1 (20%)	7,7,7	2.33	3 (42%)

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
3	ESA	B	500	2	-	0/3/3/3	-
3	ESA	A	500	2	-	0/3/3/3	-
3	ESA	E	500	2	-	3/3/3/3	-
3	ESA	C	500	2	-	0/3/3/3	-
3	ESA	F	500	2	-	0/3/3/3	-
3	ESA	D	500	2	-	3/3/3/3	-

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	500	ESA	C1-C2	-7.14	1.26	1.51
3	A	500	ESA	C1-C2	-7.10	1.26	1.51
3	F	500	ESA	C1-C2	-7.08	1.26	1.51
3	E	500	ESA	C1-C2	-6.90	1.27	1.51
3	C	500	ESA	C1-C2	-6.73	1.28	1.51

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	500	ESA	C1-C2-S	4.84	124.84	113.00
3	D	500	ESA	C1-C2-S	4.69	124.46	113.00
3	A	500	ESA	C1-C2-S	4.52	124.06	113.00
3	C	500	ESA	C1-C2-S	4.46	123.91	113.00
3	E	500	ESA	C1-C2-S	4.34	123.61	113.00

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	500	ESA	C1-C2-S-O1
3	D	500	ESA	C1-C2-S-O2
3	D	500	ESA	C1-C2-S-O3
3	E	500	ESA	C1-C2-S-O1
3	E	500	ESA	C1-C2-S-O3

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	500	ESA	1	0
3	E	500	ESA	1	0
3	C	500	ESA	3	0
3	F	500	ESA	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 [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, 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	263/267 (98%)	-0.44	1 (0%) 92 91	12, 37, 60, 67	0
1	B	263/267 (98%)	-0.34	6 (2%) 60 51	6, 36, 65, 77	0
1	C	263/267 (98%)	-0.37	1 (0%) 92 91	11, 37, 60, 69	0
1	D	263/267 (98%)	-0.47	3 (1%) 80 75	5, 36, 58, 68	0
1	E	263/267 (98%)	-0.46	3 (1%) 80 75	4, 35, 58, 65	0
1	F	263/267 (98%)	-0.48	2 (0%) 86 81	9, 33, 54, 60	0
All	All	1578/1602 (98%)	-0.43	16 (1%) 82 77	4, 36, 60, 77	0

The worst 5 of 16 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	222	ASP	3.4
1	B	77	LEU	2.7
1	D	222	ASP	2.6
1	B	63	MET	2.5
1	E	220	ASN	2.4

### 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 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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	A	501	1/1	0.83	0.16	4,4,4,4	0
2	MG	F	501	1/1	0.95	0.25	18,18,18,18	0
2	MG	C	501	1/1	0.96	0.15	1,1,1,1	0
3	ESA	C	500	6/6	0.96	0.21	39,43,44,44	0
2	MG	E	501	1/1	0.97	0.23	3,3,3,3	0
3	ESA	A	500	6/6	0.98	0.17	30,32,33,34	0
3	ESA	B	500	6/6	0.98	0.19	33,35,35,36	0
2	MG	B	501	1/1	0.98	0.23	8,8,8,8	0
3	ESA	D	500	6/6	0.98	0.14	27,29,29,29	0
3	ESA	E	500	6/6	0.98	0.16	24,25,26,27	0
3	ESA	F	500	6/6	0.98	0.19	35,37,38,39	0
2	MG	D	501	1/1	0.99	0.25	12,12,12,12	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.