



Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2020 – 12:06 pm BST

PDB ID : 1UZ6
Title : anti-Lewis X Fab fragment uncomplexed
Authors : Van Roon, A.M.M.; Pannu, N.S.; De Vrind, J.P.M.; Hokke, C.H.; Deelder, A.M.; Van Der marel, G.A.; Van Boom, J.H.; Abrahams, J.P.
Deposited on : 2004-03-05
Resolution : 2.05 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 : 2.11
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) : 2.11

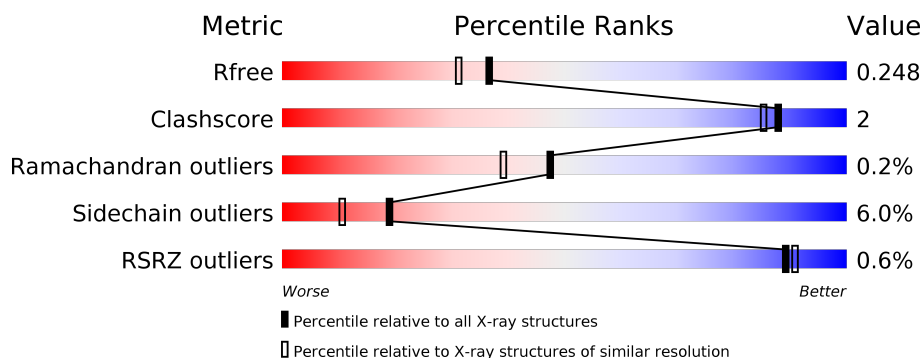
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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	E	218	<div> <div>90%</div> <div>6%</div> <div>.</div> </div>
1	L	218	<div> <div>90%</div> <div>6%</div> <div>.</div> </div>
1	M	218	<div> <div>88%</div> <div>7%</div> <div>..</div> </div>
1	V	218	<div> <div>%</div> <div>91%</div> <div>6%</div> <div>.</div> </div>
2	F	217	<div> <div>81%</div> <div>15%</div> <div>.</div> </div>
2	H	217	<div> <div>%</div> <div>85%</div> <div>11%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
2	P	217	<div><div>%</div><div><div></div><div>84%</div><div>13%</div><div></div></div><div></div></div>
2	W	217	<div><div>%</div><div><div></div><div>80%</div><div>17%</div><div></div></div><div></div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13333 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 IGG FAB (IGG3, KAPPA) LIGHT CHAIN 291-2G3-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	211	Total	C	N	O	S	0	0	1
			1630	1024	271	328	7			
1	L	211	Total	C	N	O	S	0	0	1
			1630	1024	271	328	7			
1	M	212	Total	C	N	O	S	0	0	1
			1638	1028	273	330	7			
1	V	212	Total	C	N	O	S	0	0	0
			1649	1034	276	332	7			

- Molecule 2 is a protein called IGG FAB (IGG3, KAPPA) HEAVY CHAIN 291-2G3-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	209	Total	C	N	O	S	0	1	1
			1592	1015	262	308	7			
2	H	211	Total	C	N	O	S	0	1	1
			1602	1020	264	311	7			
2	P	212	Total	C	N	O	S	0	1	1
			1608	1023	265	313	7			
2	W	212	Total	C	N	O	S	0	1	0
			1619	1029	268	315	7			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	O	S	0	0
			5	4	1		
3	L	1	Total	O	S	0	0
			5	4	1		
3	M	1	Total	O	S	0	0
			5	4	1		
3	V	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

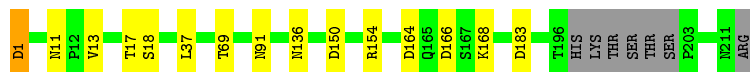
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	30	Total	O	0	0
			30	30		
4	F	63	Total	O	0	0
			63	63		
4	H	72	Total	O	0	0
			72	72		
4	L	35	Total	O	0	0
			35	35		
4	M	36	Total	O	0	0
			36	36		
4	P	46	Total	O	0	0
			46	46		
4	V	29	Total	O	0	0
			29	29		
4	W	34	Total	O	0	0
			34	34		

3 Residue-property plots


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

- Molecule 1: IGG FAB (IGG3, KAPPA) LIGHT CHAIN 291-2G3-A

Chain E: 



- Molecule 1: IGG FAB (IGG3, KAPPA) LIGHT CHAIN 291-2G3-A

Chain L: 



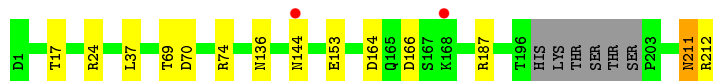
- Molecule 1: IGG FAB (IGG3, KAPPA) LIGHT CHAIN 291-2G3-A

Chain M: 




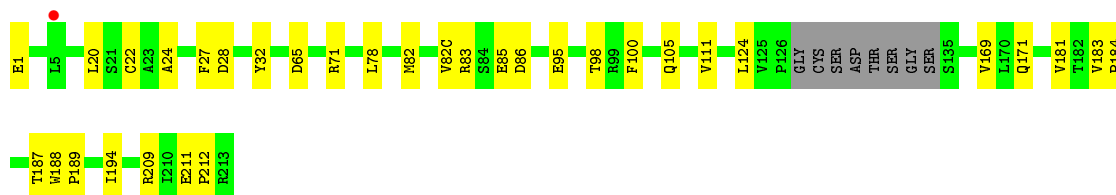
- Molecule 1: IGG FAB (IGG3, KAPPA) LIGHT CHAIN 291-2G3-A

Chain V: 

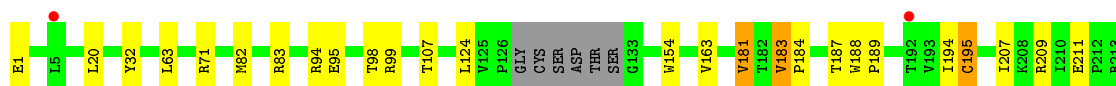
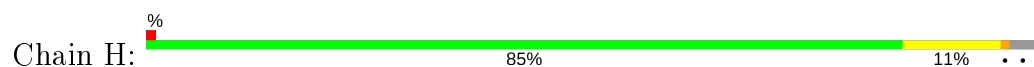


- Molecule 2: IGG FAB (IGG3, KAPPA) HEAVY CHAIN 291-2G3-A

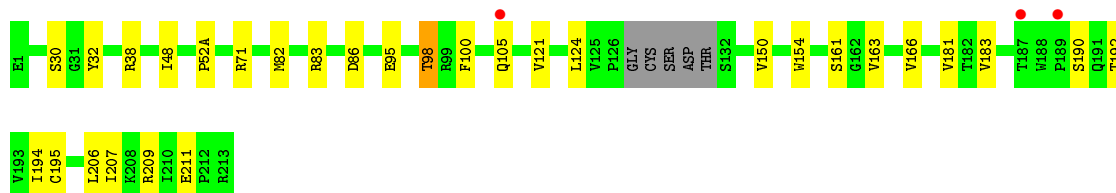
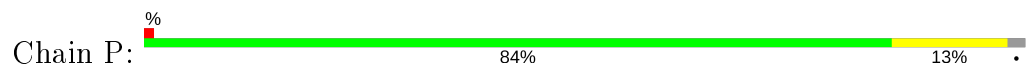
Chain F: 



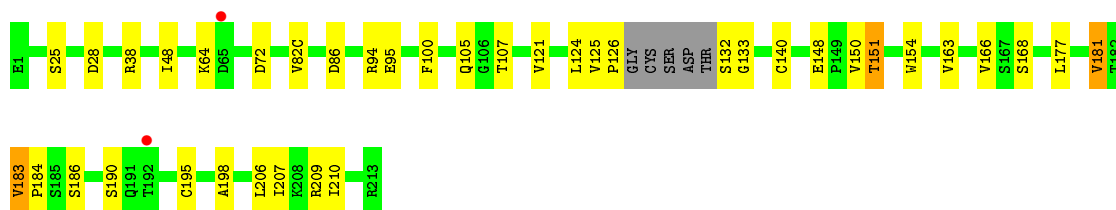
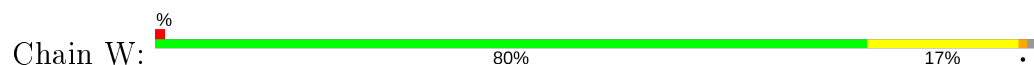
- Molecule 2: IGG FAB (IGG3, KAPPA) HEAVY CHAIN 291-2G3-A



- Molecule 2: IGG FAB (IGG3, KAPPA) HEAVY CHAIN 291-2G3-A



- Molecule 2: IGG FAB (IGG3, KAPPA) HEAVY CHAIN 291-2G3-A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	67.41Å 71.56Å 104.79Å 86.48° 71.32° 83.26°	Depositor
Resolution (Å)	45.18 – 2.05 45.09 – 2.05	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.18-2.05) 82.6 (45.09-2.05)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.2.0000	Depositor
R, R_{free}	0.198 , 0.244 0.203 , 0.248	Depositor DCC
R_{free} test set	5088 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.199 for -h,-k,-h+l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13333	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	E	0.63	0/1666	0.73	5/2263 (0.2%)
1	L	0.62	0/1666	0.74	3/2263 (0.1%)
1	M	0.62	0/1674	0.69	4/2274 (0.2%)
1	V	0.62	0/1685	0.72	5/2286 (0.2%)
2	F	0.71	0/1636	0.72	2/2231 (0.1%)
2	H	0.74	0/1646	0.72	0/2244
2	P	0.66	0/1652	0.71	1/2252 (0.0%)
2	W	0.67	1/1663 (0.1%)	0.71	3/2264 (0.1%)
All	All	0.66	1/13288 (0.0%)	0.72	23/18077 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	W	82(C)	VAL	CB-CG1	-5.35	1.41	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	86	ASP	CB-CG-OD2	7.08	124.67	118.30
2	W	86	ASP	CB-CG-OD2	6.97	124.58	118.30
2	P	86	ASP	CB-CG-OD2	6.15	123.83	118.30
1	E	150	ASP	CB-CG-OD2	6.04	123.74	118.30
1	L	164	ASP	CB-CG-OD2	6.00	123.69	118.30
1	E	183	ASP	CB-CG-OD2	5.91	123.62	118.30
2	F	28	ASP	CB-CG-OD2	5.80	123.52	118.30
1	E	166	ASP	CB-CG-OD2	5.59	123.33	118.30
1	V	166	ASP	CB-CG-OD2	5.55	123.30	118.30
1	E	164	ASP	CB-CG-OD2	5.44	123.19	118.30
1	L	150	ASP	CB-CG-OD2	5.40	123.16	118.30
1	V	187	ARG	NE-CZ-NH1	5.34	122.97	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	166	ASP	CB-CG-OD2	5.27	123.04	118.30
1	M	1	ASP	CB-CG-OD2	5.26	123.03	118.30
1	V	70	ASP	CB-CG-OD2	5.22	123.00	118.30
1	M	166	ASP	CB-CG-OD2	5.22	123.00	118.30
1	M	60	ASP	CB-CG-OD2	5.17	122.95	118.30
2	W	28	ASP	CB-CG-OD2	5.15	122.93	118.30
1	M	164	ASP	CB-CG-OD2	5.14	122.93	118.30
2	W	72	ASP	CB-CG-OD2	5.14	122.92	118.30
1	E	1	ASP	CB-CG-OD2	5.11	122.90	118.30
1	V	164	ASP	CB-CG-OD2	5.08	122.88	118.30
1	V	187	ARG	NE-CZ-NH2	-5.04	117.78	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	E	1630	0	1572	1	0
1	L	1630	0	1572	2	0
1	M	1638	0	1578	7	0
1	V	1649	0	1591	1	0
2	F	1592	0	1571	10	0
2	H	1602	0	1579	8	0
2	P	1608	0	1584	9	0
2	W	1619	0	1597	13	0
3	E	5	0	0	0	0
3	L	5	0	0	1	0
3	M	5	0	0	0	0
3	V	5	0	0	0	0
4	E	30	0	0	0	0
4	F	63	0	0	1	0
4	H	72	0	0	0	0
4	L	35	0	0	0	0
4	M	36	0	0	0	0
4	P	46	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	V	29	0	0	0	0
4	W	34	0	0	0	0
All	All	13333	0	12644	51	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 2.

All (51) 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)
2:F:98:THR:HG23	2:F:100:PHE:H	1.62	0.64
2:P:163:VAL:HG22	2:P:181:VAL:HG12	1.80	0.63
1:M:149:ILE:HD11	1:M:178:LEU:HD21	1.84	0.58
1:M:60:ASP:N	1:M:60:ASP:OD1	2.36	0.57
2:H:194:ILE:HG23	2:H:207:ILE:HG23	1.87	0.55
2:P:98:THR:HG23	2:P:100:PHE:H	1.71	0.55
2:W:121:VAL:HG21	2:W:206:LEU:HD11	1.87	0.55
1:L:192:THR:OG1	1:L:207:SER:OG	2.21	0.55
1:V:211:ASN:HD22	1:V:212:ARG:N	2.06	0.54
2:H:163:VAL:HG22	2:H:181:VAL:HG13	1.91	0.53
1:M:194:GLU:HB3	1:M:205:VAL:HG22	1.90	0.53
2:W:38:ARG:HB3	2:W:48:ILE:HD11	1.91	0.52
2:P:121:VAL:HG21	2:P:206:LEU:HD11	1.91	0.52
2:H:32:TYR:HD1	2:H:95:GLU:HG3	1.75	0.51
2:W:163:VAL:HG22	2:W:181:VAL:CG1	2.40	0.51
2:F:82(C):VAL:HG12	2:F:111:VAL:HG21	1.93	0.50
2:P:194:ILE:HG23	2:P:207:ILE:HG23	1.95	0.49
1:L:188:HIS:NE2	3:L:1211:SO4:O2	2.44	0.47
2:P:38:ARG:HB3	2:P:48:ILE:HD11	1.96	0.47
2:W:148:GLU:OE2	2:W:168:SER:OG	2.28	0.47
1:E:11:ASN:O	1:E:13:VAL:HG23	2.15	0.47
2:P:163:VAL:HG22	2:P:181:VAL:CG1	2.45	0.47
2:W:151:THR:CG2	2:W:198:ALA:HB3	2.45	0.46
2:W:140[B]:CYS:SG	2:W:210:ILE:HD11	2.55	0.46
2:F:32:TYR:HD1	2:F:95:GLU:HG3	1.80	0.46
1:M:144:ASN:HB3	1:M:196:THR:HG22	1.97	0.46
2:P:30:SER:O	2:P:52(A):PRO:HB3	2.17	0.45
1:M:37:LEU:HG	1:M:86:TYR:CE2	2.51	0.45
2:F:22:CYS:HB3	2:F:78:LEU:HB3	1.97	0.45
2:W:163:VAL:HG22	2:W:181:VAL:HG12	1.99	0.44
2:W:154:TRP:CZ3	2:W:195:CYS:HB3	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:188:TRP:CG	2:H:189:PRO:HA	2.53	0.43
2:P:32:TYR:HD2	2:P:95:GLU:HG3	1.84	0.43
2:F:24:ALA:HB1	2:F:27:PHE:CZ	2.54	0.43
2:F:20:LEU:HG	2:F:82:MET:HE2	2.01	0.43
2:H:20:LEU:HD11	2:H:82:MET:CE	2.49	0.43
2:F:95:GLU:O	2:F:98:THR:HG22	2.19	0.42
2:W:94:ARG:HD2	2:W:100:PHE:O	2.19	0.42
2:W:177:LEU:HD12	2:W:177:LEU:C	2.40	0.42
2:F:169:VAL:HG23	4:F:2050:HOH:O	2.19	0.42
2:F:184:PRO:O	2:F:187:THR:OG1	2.35	0.42
1:M:136:ASN:ND2	1:M:173:SER:OG	2.53	0.42
2:H:154:TRP:CZ3	2:H:195:CYS:HB3	2.54	0.41
1:M:132:VAL:HG22	1:M:177:THR:HG23	2.02	0.41
2:F:188:TRP:CG	2:F:189:PRO:HA	2.55	0.41
2:W:132:SER:OG	2:W:133:GLY:N	2.54	0.41
2:W:125:VAL:HG13	2:W:126:PRO:HD2	2.01	0.41
2:H:183:VAL:HG13	2:H:187:THR:OG1	2.21	0.41
2:W:183:VAL:HG22	2:W:184:PRO:HD2	2.03	0.40
2:H:184:PRO:O	2:H:187:THR:OG1	2.34	0.40
2:P:154:TRP:CZ3	2:P:195:CYS:HB3	2.57	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	E	207/218 (95%)	202 (98%)	5 (2%)	0	100	100
1	L	207/218 (95%)	205 (99%)	2 (1%)	0	100	100
1	M	208/218 (95%)	201 (97%)	7 (3%)	0	100	100
1	V	208/218 (95%)	204 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	206/217 (95%)	203 (98%)	2 (1%)	1 (0%)	29	18
2	H	208/217 (96%)	205 (99%)	3 (1%)	0	100	100
2	P	209/217 (96%)	205 (98%)	3 (1%)	1 (0%)	29	18
2	W	209/217 (96%)	205 (98%)	3 (1%)	1 (0%)	29	18
All	All	1662/1740 (96%)	1630 (98%)	29 (2%)	3 (0%)	47	39

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	W	95	GLU
2	P	190	SER
2	F	212	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	E	185/193 (96%)	176 (95%)	9 (5%)	25	17
1	L	185/193 (96%)	177 (96%)	8 (4%)	29	22
1	M	186/193 (96%)	178 (96%)	8 (4%)	29	22
1	V	187/193 (97%)	178 (95%)	9 (5%)	25	18
2	F	180/186 (97%)	167 (93%)	13 (7%)	14	7
2	H	181/186 (97%)	167 (92%)	14 (8%)	13	5
2	P	182/186 (98%)	169 (93%)	13 (7%)	14	7
2	W	183/186 (98%)	169 (92%)	14 (8%)	13	5
All	All	1469/1516 (97%)	1381 (94%)	88 (6%)	19	11

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	1	ASP

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Mol	Chain	Res	Type
1	E	17	THR
1	E	18	SER
1	E	37	LEU
1	E	69	THR
1	E	91	ASN
1	E	136	ASN
1	E	154	ARG
1	E	168	LYS
2	F	1	GLU
2	F	65	ASP
2	F	71	ARG
2	F	83	ARG
2	F	85	GLU
2	F	105	GLN
2	F	124	LEU
2	F	171	GLN
2	F	181	VAL
2	F	183	VAL
2	F	194	ILE
2	F	209	ARG
2	F	211	GLU
2	H	1	GLU
2	H	63	LEU
2	H	71	ARG
2	H	83	ARG
2	H	94	ARG
2	H	98	THR
2	H	99	ARG
2	H	107	THR
2	H	124	LEU
2	H	181	VAL
2	H	183	VAL
2	H	195	CYS
2	H	209	ARG
2	H	211	GLU
1	L	17	THR
1	L	24	ARG
1	L	37	LEU
1	L	136	ASN
1	L	152	SER
1	L	154	ARG
1	L	189	ASN

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Mol	Chain	Res	Type
1	L	196	THR
1	M	37	LEU
1	M	54	LEU
1	M	60	ASP
1	M	67	SER
1	M	69	THR
1	M	194	GLU
1	M	196	THR
1	M	207	SER
2	P	71	ARG
2	P	82	MET
2	P	83	ARG
2	P	98	THR
2	P	105	GLN
2	P	124	LEU
2	P	150	VAL
2	P	161	SER
2	P	166	VAL
2	P	183	VAL
2	P	192	THR
2	P	209	ARG
2	P	211	GLU
1	V	17	THR
1	V	24	ARG
1	V	37	LEU
1	V	69	THR
1	V	74	ARG
1	V	136	ASN
1	V	144	ASN
1	V	153	GLU
1	V	211	ASN
2	W	25	SER
2	W	64	LYS
2	W	105	GLN
2	W	107	THR
2	W	124	LEU
2	W	150	VAL
2	W	151	THR
2	W	166	VAL
2	W	181	VAL
2	W	183	VAL
2	W	186	SER

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Mol	Chain	Res	Type
2	W	190	SER
2	W	207	ILE
2	W	209	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	50	GLN
1	E	91	ASN
1	E	136	ASN
1	E	209	ASN
1	L	50	GLN
1	L	136	ASN
1	M	123	GLN
1	M	136	ASN
1	V	50	GLN
1	V	136	ASN
1	V	209	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	L	1211	-	4,4,4	0.12	0	6,6,6	0.36	0
3	SO4	M	1212	-	4,4,4	0.33	0	6,6,6	0.32	0
3	SO4	E	1211	-	4,4,4	0.33	0	6,6,6	0.27	0
3	SO4	V	1213	-	4,4,4	0.35	0	6,6,6	0.26	0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	1211	SO4	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, 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	E	211/218 (96%)	-0.07	0 100 100	37, 53, 64, 67	0
1	L	211/218 (96%)	-0.02	0 100 100	39, 53, 63, 68	0
1	M	212/218 (97%)	-0.12	0 100 100	35, 53, 63, 68	0
1	V	212/218 (97%)	-0.01	2 (0%) 84 86	40, 54, 63, 75	0
2	F	209/217 (96%)	-0.05	1 (0%) 91 92	30, 50, 61, 69	0
2	H	211/217 (97%)	-0.05	2 (0%) 84 86	30, 49, 60, 69	0
2	P	212/217 (97%)	0.05	3 (1%) 75 78	30, 51, 61, 67	0
2	W	212/217 (97%)	-0.02	2 (0%) 84 86	31, 51, 62, 70	0
All	All	1690/1740 (97%)	-0.04	10 (0%) 89 91	30, 52, 63, 75	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	P	189	PRO	3.0
2	P	105	GLN	2.6
2	W	192	THR	2.5
1	V	168	LYS	2.5
2	F	5	LEU	2.4
2	P	187	THR	2.4
2	H	5	LEU	2.3
2	W	65	ASP	2.3
1	V	144	ASN	2.3
2	H	192	THR	2.2

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(\AA^2)	Q<0.9
3	SO4	L	1211	5/5	0.96	0.12	56,56,58,61	0
3	SO4	M	1212	5/5	0.98	0.07	57,58,60,60	0
3	SO4	V	1213	5/5	0.98	0.07	54,56,58,58	0
3	SO4	E	1211	5/5	0.99	0.08	54,56,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.